

**Edwards Aquifer Habitat Conservation Plan
Expanded Water Quality Monitoring Report**

December 2015

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APPENDICES

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1.0 INTRODUCTION

The Edwards Aquifer Authority (EAA) and its predecessor agency, the Edwards Underground Water District (EUWD), in cooperation with the U.S. Geological Survey (USGS) and the Texas Water Development Board (TWDB) have maintained a water quality sampling program since 1968. Analyses of these data have been used by the EAA to assess aquifer water quality. This routine or historical sampling program involves the analyses of a broad spectrum of parameters in wells, springs, and streams across the region. The EAA's existing sampling program was expanded with the adoption of the Edwards Aquifer Habitat Conservation Program (EAHCP) to include collection of additional samples and sample types in the immediate vicinity of Comal and San Marcos Springs. The expanded water quality sampling program was developed in accordance with the directives of the EAHCP to identify and assess potential impairments to water quality within the Comal River and headwaters of the San Marcos River systems. The expanded EAHCP sampling requirements are described in the *2015 Water Quality Monitoring Program Strategy for Comal Springs and San Marcos Springs in Support of the Edwards Aquifer Habitat Conservation Plan* (EAHCP Workplan) (EAA, 2014) which is included in Appendix A of this document.

Based on the requirements of the EAHCP Workplan, the expanded sampling program requires the collection of the following sample types:

1. Surface water (base flow) samples
2. Sediment samples
3. Real time instrument (RTI) water quality monitoring
4. Stormwater sampling
5. Passive Diffusion Samplers (PDS)

SWCA Environmental Consultants, Inc. (SWCA) was contracted by the EAA to execute the expanded sampling program in 2014 and 2015, with the exception of RTI water quality monitoring, which is still conducted by the EAA. Surface water quality monitoring was enhanced for 2014 and 2015 by the addition of passive diffusion samplers (PDS) for trace organic compounds analyses, which was conducted by SWCA. A groundwater sampling element was also included in the sampling program, which was to be conducted during periods of extremely low spring flow from Comal and San Marcos springs. The groundwater sampling element of the sampling program was only to be conducted if spring flow rates dropped below 30 cubic feet per second (cfs) at Comal Springs or below 50 cfs at San Marcos Springs. Spring flow rates remained above 30 cfs at Comal Springs and above 50 cfs at San Marcos Springs during 2015 and; therefore, the groundwater sampling element was not conducted. Hydrographs of spring flow are presented in Appendix B.

Prior to the implementation of the EAHCP, the historical sampling program had not specifically addressed surface water quality, sediment quality, real time changes for basic water quality parameters, or stormwater impacts along the Comal River or headwaters of the San Marcos River. Therefore, this expanded sampling program was designed to gather data specific to all of the new parameters. This report presents the surface water, sediment, stormwater and PDS data collected by SWCA in 2015. The data set represents the third year of the program and is not sufficient to establish any long-term trends or patterns.

For purposes of this report, the Comal River may also be referred to as Comal Springs or Comal Springs complex, and the San Marcos River headwaters may also be referred to as San Marcos Springs or San Marcos Springs complex. An overview of surface water, sediment, and stormwater sample locations for Comal and San Marcos springs are shown in Figures 1 and 2. Figures 3–6 provide detailed location data for sample points at the Comal Springs complex. Figures 7–10 provide detailed locations for the sample points at the San Marcos Springs complex.

1.1 Surface Water (Base Flow) Samples

Surface water (base flow) samples are collected twice annually at each spring complex. The Comal Springs complex has five sample locations along the Comal system from the upstream end of Landa Lake (where Blieders Creek empties into the headwaters of Landa Lake) to the south end of the Comal River, upstream of the confluence with the Guadalupe River. In the San Marcos system, surface water samples are collected at seven locations. Sample sites begin at Sink Creek upstream of the headwaters of Spring Lake on the north end of the system and end downstream of Capes Dam on the south end of the system.

Surface water sample locations are designed to provide water quality data for the majority of the surface waters of each spring system and river reach of concern. Samples are collected both above and below where each system's surface waters are influenced by springflow as well as other potential surface water inputs (such as Dry Comal Creek or Purgatory Creek). Surface water samples are analyzed for a broad spectrum of parameters as outlined in Table 1. Surface water samples were collected in March and September 2015.

Regulatory standards for surface water quality vary depending upon type of use. For this report, surface water results are compared to drinking water quality standards (30 TAC, Chapter 290, Subchapter F) for detected constituents of concern. These guidelines were selected for use since in general they provide the most stringent quality standards. For detections of interest that do not have an established maximum contaminant level (MCL) under 30 TAC 290, the Texas Risk Reduction Program (TRRP) from 30 TAC 350 was substituted. The TRRP standards used are the Tier I, residential standards and are referred to as protective concentration levels (PCLs). Pharmaceutical and Personal Care Products (PPCP) including caffeine can be chemicals of concern because they can indicate the presence of contamination from anthropogenic sources including wastewater discharge (U.S. Environmental Protection Agency [EPA], 2012). Currently, there are no regulatory standards to compare caffeine detections against, but results are listed in this report to provide an indication of potential anthropogenic impacts. Additionally, bacteriological results were compared with Texas Surface Water Quality Standards for primary recreational waters (30 TAC, Chapter 307, Rule 307.7). The 30 TAC 307.7 standards are typically applied to waters affected by anthropogenic sources and are used here solely to provide a reference level for bacterial counts. Other guidelines may be more useful or appropriate for particular research; however, for the scope of this report these standards provide an appropriate and applicable guideline with regard to water quality.

Table 1. Listing of Analytical Parameters by Sample Type

| Analytical Parameter | Surface Water (Base Flow) Samples | Stormwater Samples | Sediment Samples | PDS |
|--|---|-----------------------|------------------------|------|
| Volatile Organic Compounds (VOCs) | Yes | Yes | Yes | Yes* |
| Semi-volatile Organic Compounds (SVOCs) | Yes | Yes | Yes | Yes* |
| Organochlorine Pesticides | Yes | Yes | Yes | Yes* |
| Polychlorinated Biphenyls (PCBs) | Yes | Yes | Yes | No |
| Organophosphorous Pesticides | Yes | Yes | Yes | No |
| Herbicides | Yes | Yes | Yes | No |
| Metals (Al, Sb, As, Ba, Be, Cd, Cr [total], Cu, Fe, Pb, Mn, Hg, Ni, Se, Ag, Tl, and Zn) | Yes | Yes | Yes | No |
| General water quality parameters (GWQP), total alkalinity (as CaCO ₃), bicarbonate alkalinity (as CaCO ₃), carbonate alkalinity (as CaCO ₃); Cl, Br, NO ₃ , SO ₄ , F, pH, total dissolved solids (TDS), total suspended solids (TSS), Ca, Mg, Na, K, Si, Sr, CO ₃ , | Yes | Yes | No TDS, TSS, or TKN | No |
| Phosphorus (total) | Yes | Yes | Yes | No |
| Total Organic Carbon (TOC), | Yes | Yes | Yes | No |
| Dissolved Organic Carbon (DOC) | Yes | Yes | Yes | No |
| Total Kjeldahl Nitrogen (TKN) | Yes | Yes | Yes | No |
| Bacteria (<i>E. coli</i>) | Yes | Yes | No | No |
| Field Parameters (DO, pH, Conductivity, Turbidity, Temperature) | Yes | Yes | No | No |
| Caffeine | Yes | Yes | No | No |

* Passive diffusion samplers (PDS) samplers are analyzed for a modified set of VOCs, SVOCs, and organochlorine pesticides

Figure 1. EAHCP expanded water quality monitoring program, Comal Springs and River.

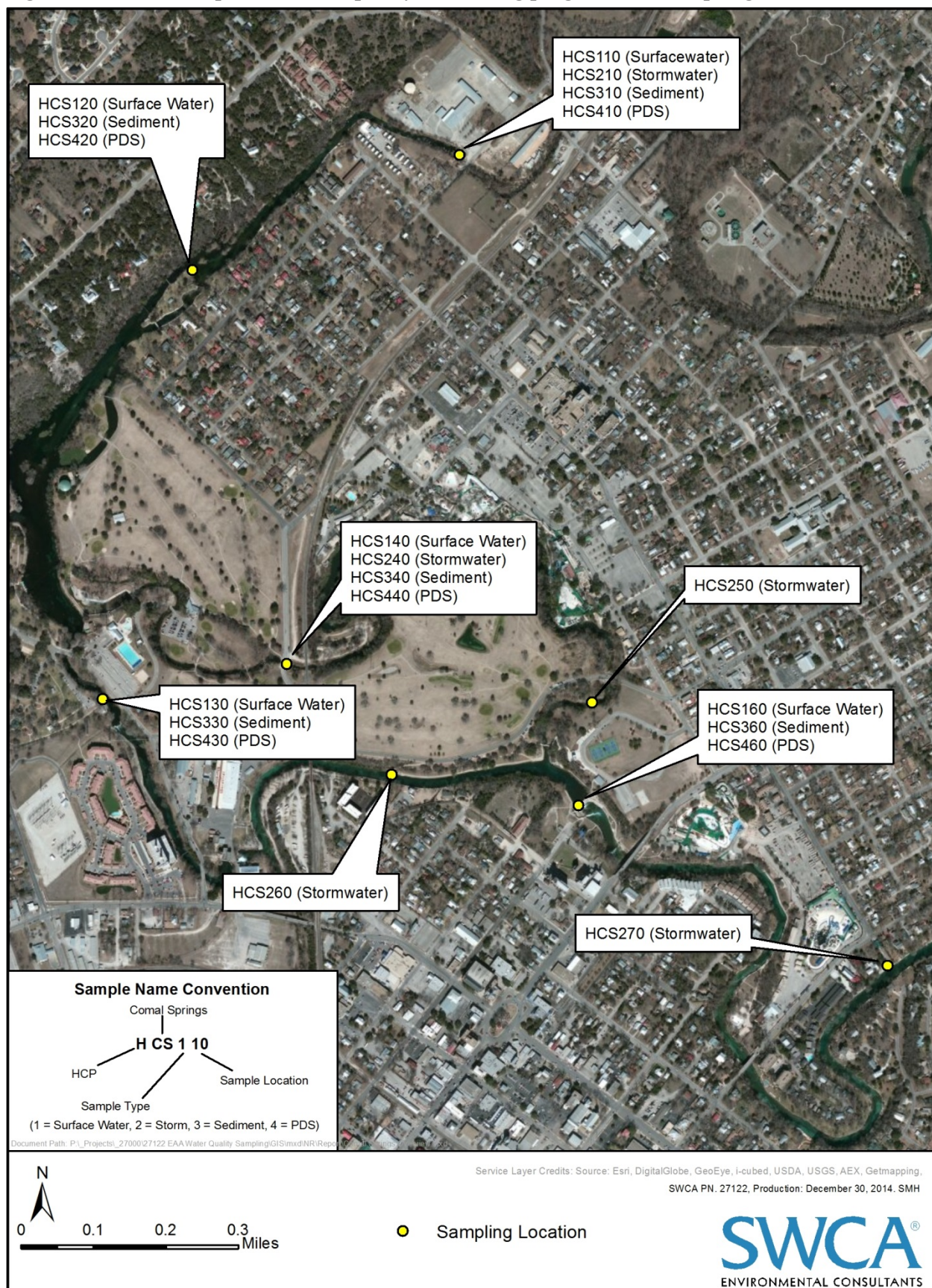
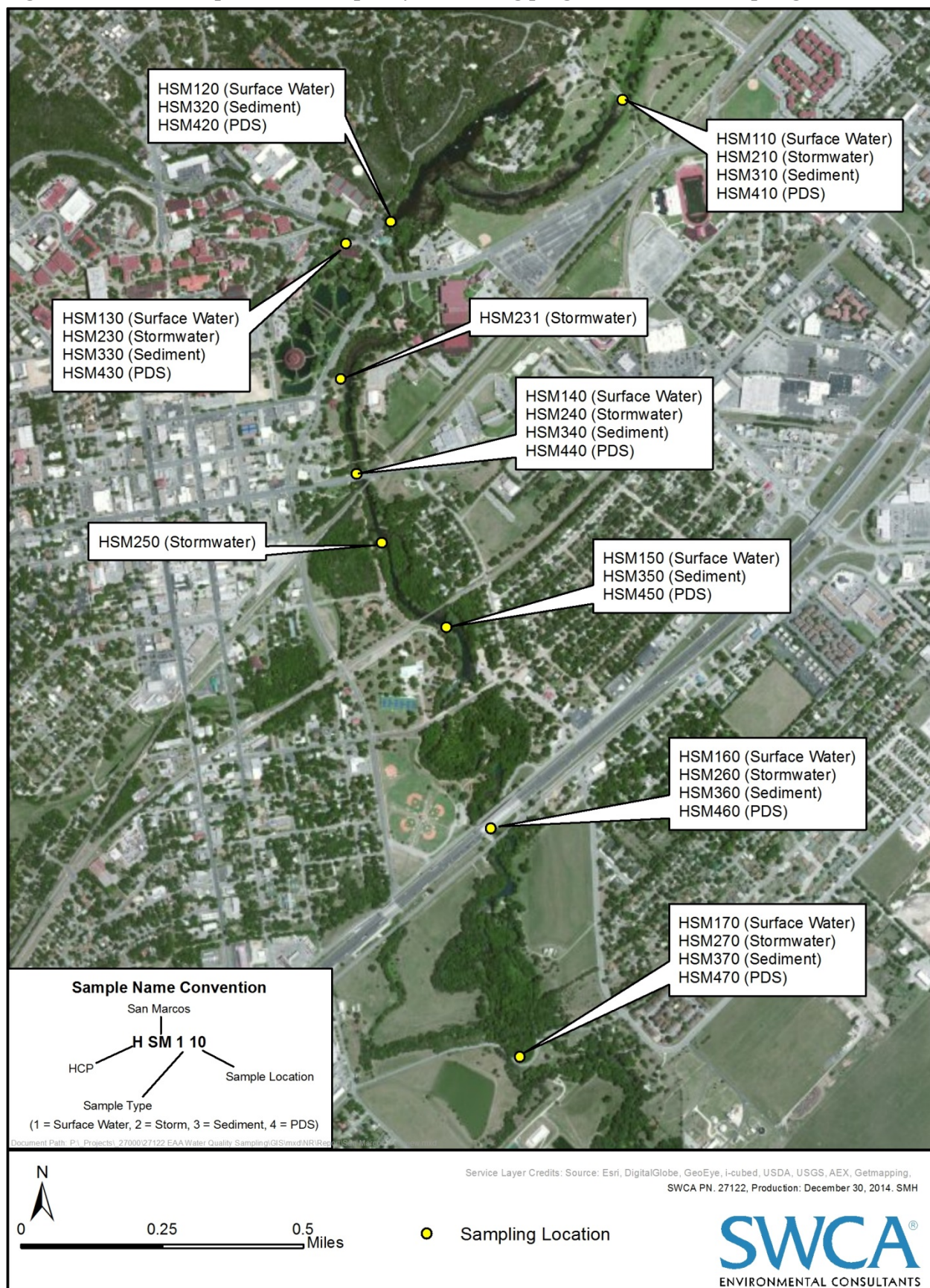


Figure 2. EAHCP expanded water quality monitoring program, San Marcos Springs and River.



1.2 Sediment Samples

Collection of sediment samples within each spring system was included in the program to help ascertain potential effects on listed species via direct or indirect exposure to sediments. Designated sediment sample locations were coincident with surface water (base flow) sample locations at each spring complex. Specifically, five sediment samples were collected from the Comal Springs area, and seven locations were sampled within the San Marcos area. In the first two years of the program sediment samples were collected from the sediment surface to approximately 18 inches below the surface. For 2015, the EAHCP workplan was changed to include sampling only the first three inches below the surface of the sediment. Samples were analyzed for the parameters listed in Table 1.

Sediment samples were collected as close to each associated surface water sample locations as possible. However, for some of the samples, collection points were moved slightly to find adequate sediment or to avoid rocky substrates that prevented collection of adequate sample volume. Appendix C of this report discusses sample locations where any significant deviations from this approach occurred.

Analytical results for sediment samples are compared to the sediment quality guidelines published in *Development and Evaluation of Consensus-Based Sediment Quality Guidelines for Freshwater Ecosystems* (MacDonald, Ingersoll, and Berger, 2000). These guidelines are based on determination of probable sediment toxicity in freshwater ecosystems and provide a numerical sediment quality guideline for 28 chemicals of concern. The guidance provides two basic standards for comparison: 1) threshold effect concentration (TEC), and 2) probable effect concentration (PEC). PEC values for additional chemicals of concern that were not included in MacDonald et al. (2000) were taken from *Conducting Ecological Risk Assessments at Remediation Sites in Texas* developed by the Texas Commission on Environmental Quality (TCEQ, 2014a). Analytical results with a concentration below the TEC are predicted to be non-toxic (on sediment-dwelling organisms), while results with a concentration above the PEC are indicated as having a probable toxic effect on sediment-dwelling organisms. Detected compounds with concentrations between the TEC and PEC are considered equally likely to be toxic or non-toxic. While numerous other guidelines for sediment quality exist, these guidelines provide a good reference for the scope of the current investigation. Future researchers may find other guidelines more specific to particular concerns or interest.

1.3 Stormwater Samples

Stormwater sampling was performed at five Comal Springs locations and at seven San Marcos Springs locations. Stormwater sample collection was adopted as part of the expanded water quality monitoring effort to assess potential contaminants that may be present in surface water runoff generated by storm events. The stormwater sampling effort was designed to assess what changes in water quality occur within each surface water system during a storm event. Storm samples were collected in association with various surface water inputs along each spring complex within the study area. Appendix C of this report discusses details of each stormwater sample location and any deviations from the EAHCP workplan. Stormwater samples were analyzed for the same parameters as surface water (base flow) samples as outlined in Table 1.

Stormwater samples were collected at three points across the storm hydrograph for each stormwater sampling site. Sample collection was targeted for the rising limb, peak, and receding limb of the storm

hydrograph. Timing for sample collection was generally determined using the RTI system's conductivity and turbidity parameters rather than the flow measurements from the USGS streamflow gauges. The USGS gauges are only updated on an hourly basis whereas data from the RTIs was available on 15-minute intervals and provided more timely information. Automated sample collection equipment was not utilized for stormwater sample collection due to sample volume, preservation, and analysis limitations. Therefore, sampling was conducted manually. Each spring group was sampled twice for stormwater events during calendar year 2015 per the EAHCP Workplan.

As previously mentioned, standards for surface water quality vary dependent upon type of use. For this report, stormwater results are compared to drinking water quality standards (30 TAC, Chapter 290, Subchapter F) for detected chemicals of concern. These guidelines were selected for use since, in general, they provide the most stringent quality standards. For detections of interest that do not have an established MCL under 30 TAC 290, the TRRP PCLs from 30 TAC 350 was substituted. The TRRP standards used are the Tier I, residential standards. Currently, there are no regulatory standards to compare caffeine detections against, but results are listed in this report to provide an indication of anthropogenic contamination. Bacterial counts were compared with Texas Surface Water Quality Standards (30 TAC 307). Other guidelines may be more useful or appropriate for particular research; however, for the scope of this report these standards provide an appropriate and applicable guideline with regard to water quality.

1.4 Surface Water Passive Sampling

Amplified Geochemical Imaging (AGI) LLC, PDSs were deployed in both spring complexes to measure trace organic constituents. Samplers consisted of a sorbent solid phase material that concentrates compounds from the environment. Following collection, the analytes of interest were eluted and analyzed by gas chromatography coupled with a mass spectrometry detector (GC-MS). The increased contact time associated with long-term deployment of the collection material allowed the analytes to be greatly concentrated beyond what is typically found in water samples. Therefore, the PDS provides greater sensitivity to trace level constituents. Analyzed parameters can be found in Table 1.

PDS samplers were deployed to each of the 12 sample sites for two-week periods in February, April, June, August, October and December 2015. Sample points coincided with surface water collection points unless prevented by field conditions, and any alterations are discussed in Appendix C.

2.0 SAMPLE LOCATION DETAIL

Details of individual sample locations are provided in the following figures. Figures 3–6 show sample location details for the Comal Springs area. Figures 7–10 provide sample location details for the San Marcos Springs area.

Figure 3. EAHCP Comal Springs detailed map indicating sample locations 110, 210, 310, 120, and 320.



Figure 4. EAHCP Comal Springs detailed map indicating sample locations 130 and 330.

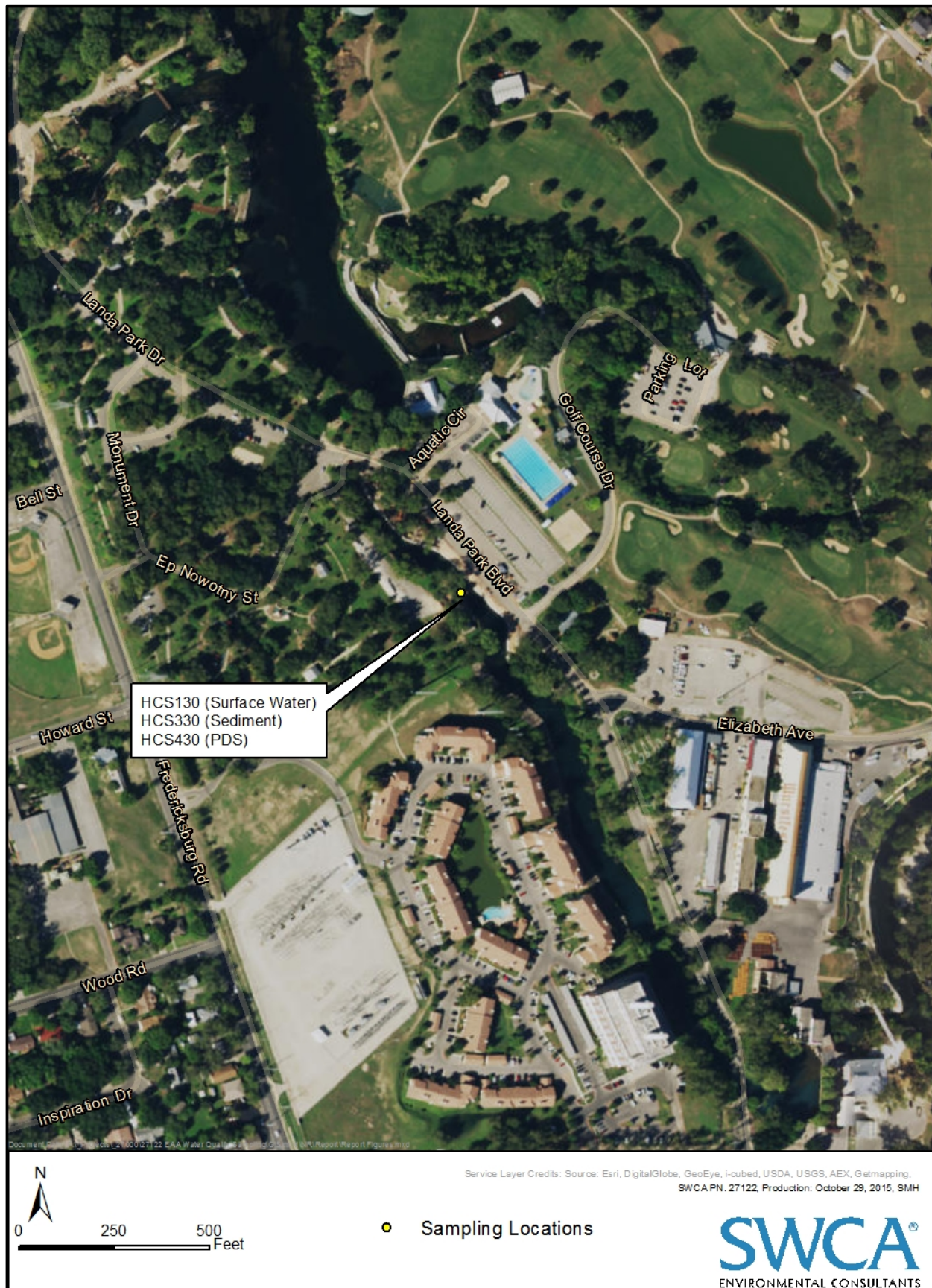


Figure 5. EAHCP Comal Springs detailed map indicating sample locations 140, 240, 340, 250, 160, 260, and 360.

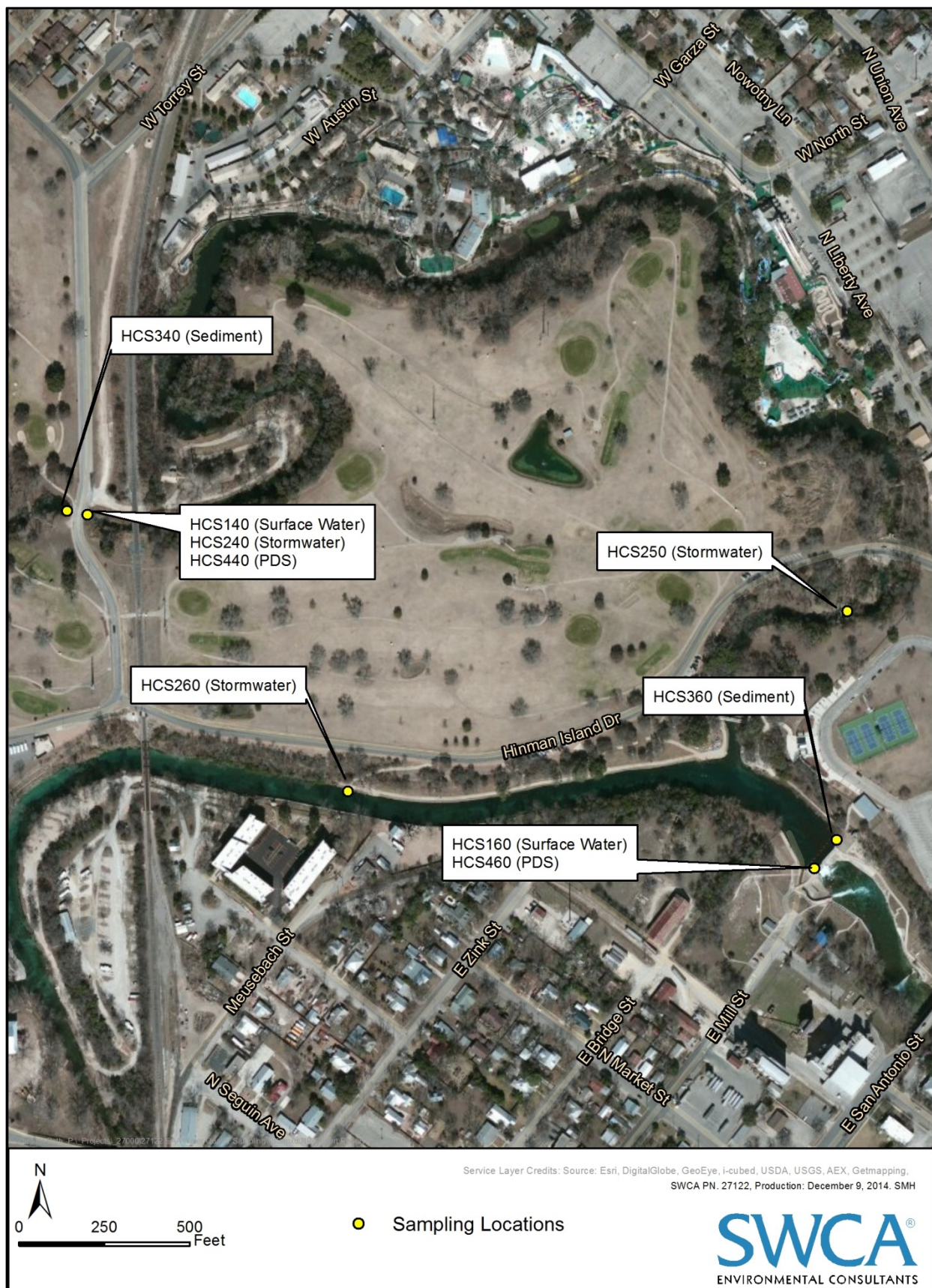


Figure 6. EAHCP Comal Springs detailed map indicating sample location 270.

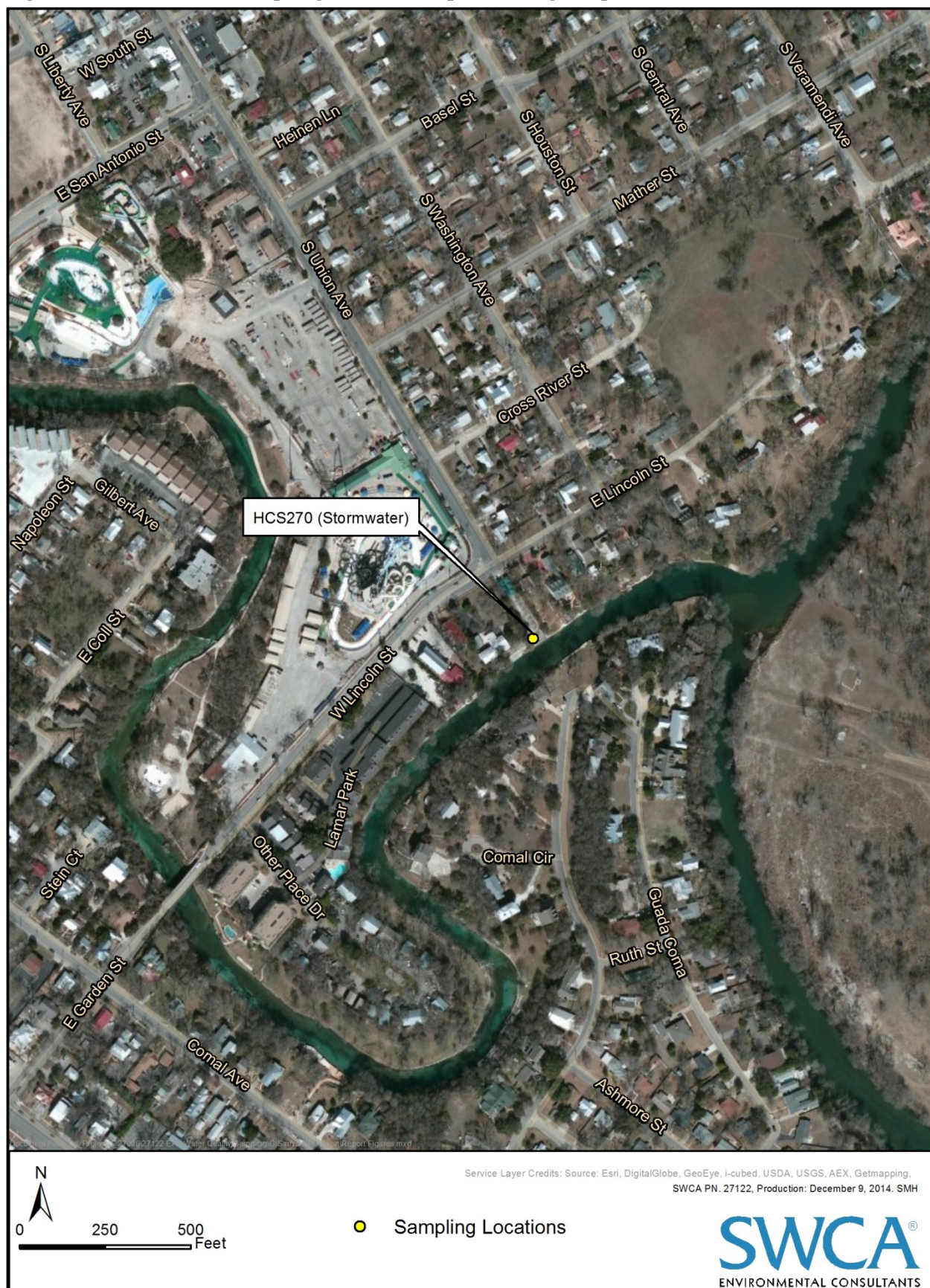


Figure 7. EAHCP San Marcos Springs detailed map indicating sample locations 110, 210, 310, 120, and 320.

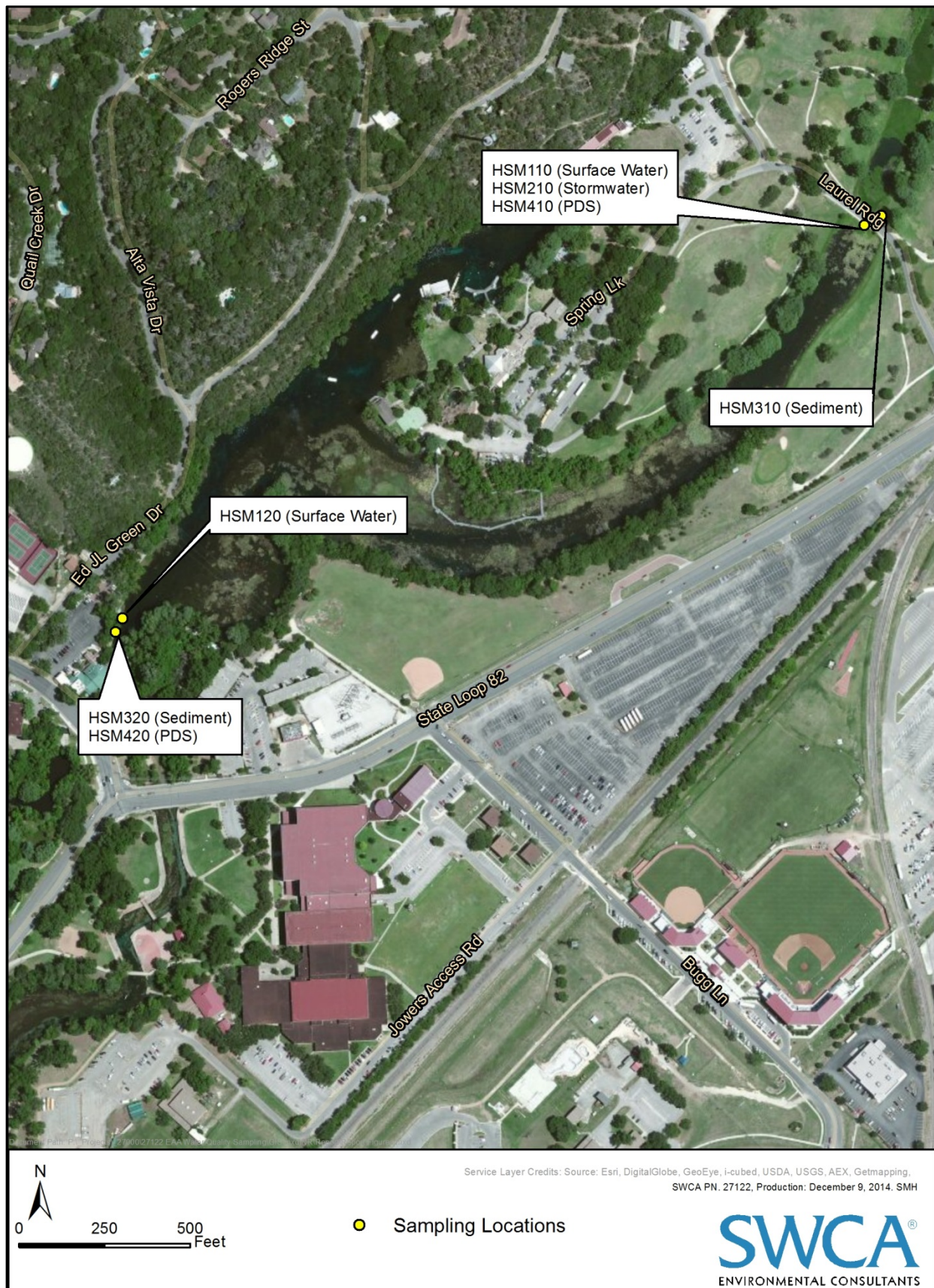


Figure 8. EAHCP San Marcos Springs detailed map indicating sample locations 130, 230, 330, 231, 140, 240, and 340.

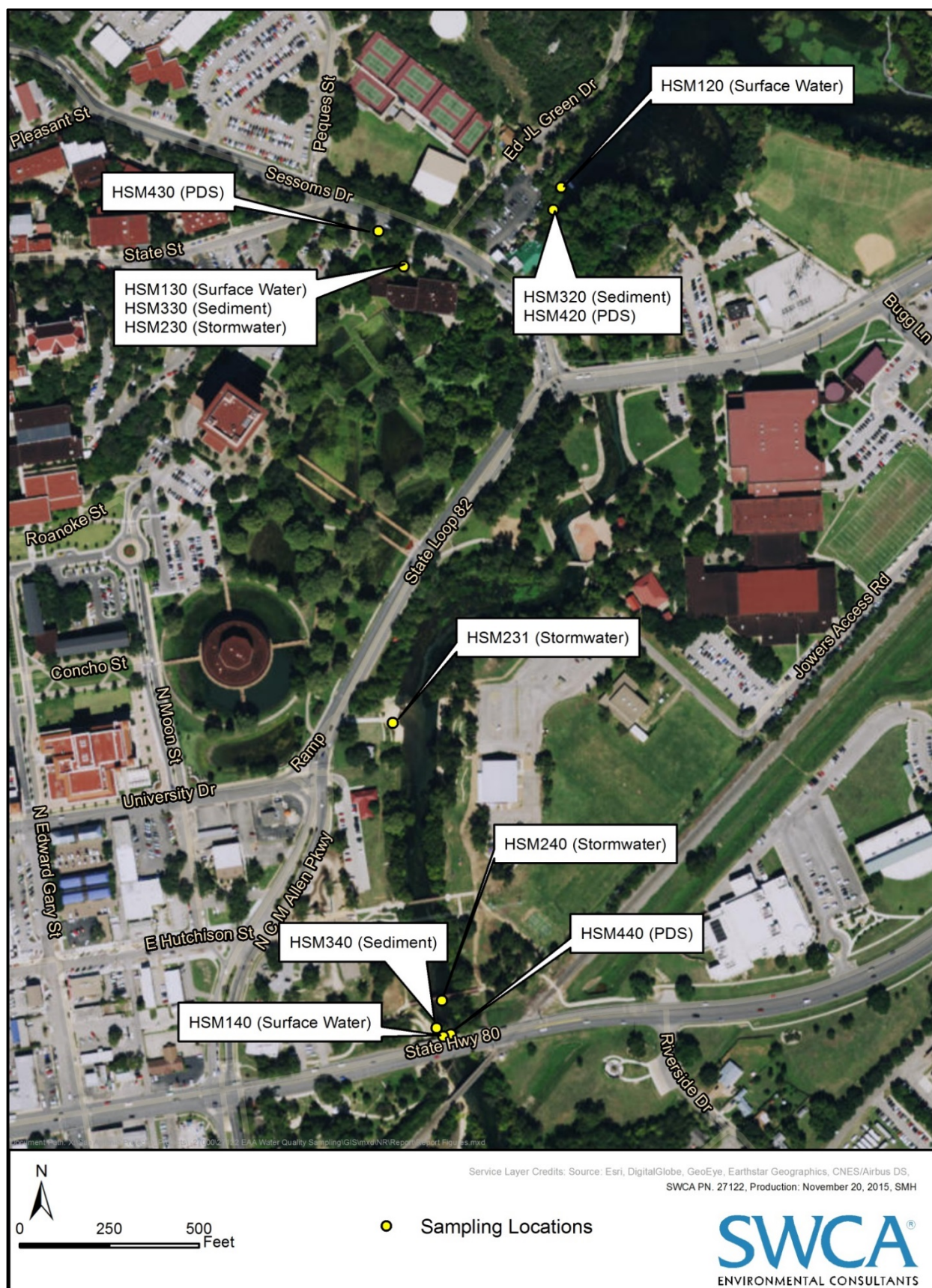


Figure 9. EAHCP San Marcos Springs detailed map indicating sample locations 150, 250, and 350.



Figure 10. EAHCP San Marcos Springs detailed indicating map sample locations 160, 260, 360, 170, 270, and 370.



3.0 MODIFICATIONS OF ACTIVITIES DUE TO DROUGHT

Sampling activities continued to be affected by drought conditions within the project area. Significant rainfall occurred during the first half of 2015, which contrasted severe drought conditions experienced in 2014. However, rainfall was very sparse from July through much of October. Rain events were generally scattered in nature and often too small in magnitude to generate sufficient runoff to sample. To collect the stormwater sample events for the second half of 2015, both spring complexes were sampled on the same day during the October 23, 2015, storm event. Due to the high probability that the October 23, 2015, storm event would be of the size and magnitude sufficient to sample, additional personnel and resources were made available to sample at both spring complexes on the same day.

Also, due to abundant rainfall in the spring of 2015, flow rates did not drop below 30 cfs at Comal Springs, or below 50 cfs at San Marcos Springs. No extreme low-flow sampling was initiated at water wells (Sections 6.4.3.3 and 6.4.4.3 of the Edwards Aquifer Habitat Conservation Plan).

4.0 AFFECT ON COVERED SPECIES

The implementation of the EAHCP water quality and sediment sampling program provided baseline data along the Comal River and upper reaches of the San Marcos River system. Water quality grab samples were collected twice from each river during base flow conditions and during two storm events. Sediment samples were also collected from both systems. PDSs were used to evaluate trace organic compounds six times throughout the year.

The collection and analysis of water quality and sediment samples aids evaluation of the habitat of species by providing base flow, storm flow, and sediment quality data. The data included water quality discharging directly from the springs and water discharging into the Comal and San Marcos Rivers below the springs.

In Section 7, analytical results are compared to various water quality and sediment standards as guidelines to identify any existing problems and create a body of baseline data to ascertain any long-term sediment and water quality trends.

5.0 LOGISTICS

To accommodate the needs of the EAHCP's expanded water quality monitoring program, a significant amount of resources are required. These resources, including sampling equipment, safety gear, trained staff, and sampling schedules, are all key components to the program. In addition, the development of sampling strategies and planning of each sampling event is required to insure that resources are used efficiently, and collection is completed within the scheduled time frame. The strategies must account for the unpredictable nature of storm events. Below is a short synopsis of events and tasks undertaken to accomplish the necessary logistics for the EAHCP sampling program.

5.1 Surface Water Sampling Program

Prior to each sampling event, SWCA staff acquired necessary supplies and equipment including laboratory sample kits, disposable bailers, and 0.45-micron filters.

5.2 Sediment Sampling Program

In May 2015, SWCA staff purchased trowels for sediment sampling and acquired sample containers from the contract laboratory.

5.3 Stormwater Program

Prior to each sampling event, laboratory sample kits were acquired and prepared for use in the field. All other sampling and safety supplies were kept stocked and ready for mobilization in the event a storm occurred. Weather forecasts are monitored on a regular basis to determine if teams would be mobilized for a potential sampling event. Prior to mobilization many logistical concerns have to be addressed including, but not limited to, personnel availability, safety, base camp reservation, vehicle availability, sonde rental, laboratory notifications, and preparing checks for laboratory payment.

5.4 Surface Water Passive Sampling Program

PDS were acquired from the contract laboratory approximately two weeks prior to each sampling event. Sample deployment devices were constructed in 2014 but were decontaminated and placed inside clean plastic bags the day before or day of each deployment.

6.0 SAMPLE COLLECTION METHODOLOGY

6.1 Surface Water / Base Flow Sampling Program

Surface water quality base flow grab samples were collected from five sites throughout the Comal Springs complex and seven sites throughout the San Marcos Springs complex biannually. According to the EAHCP workplan, the sample dates were to be six months apart. The preferred method for obtaining a surface water sample is to reach the sampling location from the shoreline or wade to the sample location, obtain field parameters (pH, specific conductivity, dissolved oxygen, and temperature) and then insert the sample bottle directly into the water or use a sample bottle and pole assembly. SWCA collected samples directly within sample bottles as opposed to using a pole assembly. Samples collected while wading were collected on the upstream side of the sampler. Samples were collected in accordance with the criteria set forth in the *EAA Groundwater Quality Monitoring Plan*.

Filtration for methods 6010B (metals), 6020 (metals), 7470A (mercury) and field alkalinity were performed at the sample location by using a 0.45-micron high capacity cartridge filter attached to a weighted, single-sample disposable bailer. Preservatives were placed in the bottles (as appropriate) by the contracted laboratory. Ice was placed into the cooler prior to or immediately after sampling and later shipped to the contract laboratory. When not in use or after collection, sampling equipment and/or coolers containing samples were secured inside the SWCA vehicles to maintain appropriate sample custody and security.

The *EAA Groundwater Quality Monitoring Plan* required the collection of one field duplicate sample for each spring complex per sampling event. The field duplicate was sampled immediately after the parent water quality sample and in the same manner as the parent water quality sample.

Analyses for field alkalinity were conducted at SWCA's San Antonio office. The method used for field alkalinity is discussed in detail in the *EAA Groundwater Quality Monitoring Plan* (Appendix D). Representative photographs of field activities are included in Appendix E.

6.2 Sediment Sampling Program

Sediment samples were collected once annually from the first three inches of sediment below the streambed surface at each of the 12 sampling locations. Sediment sample collection points generally coincided with the surface water collection points at each of the 12 sample locations in the spring complexes but varied slightly based on field conditions. Based on the amount of available sediment at each site, the location and area sampled varied. Sample collection location variations are discussed in Appendix C. Sediment sample collection methods were consistent with the guidelines established in the *EAA Groundwater Quality Monitoring Plan*. The majority of samples were collected using stainless steel hand trowels. The trowel was inserted into the sediment three inches, and the sample was scooped into sample containers provided by the contract laboratory. One 8-ounce jar and one 2-ounce jar for volatile organic compound (VOC) analysis were collected at each location. Samples were composed of sediment collected at three locations at each sample point and were individually homogenized at the contract laboratory prior to analysis. The water depth at HCS330 made it impossible to collect a sample using hand trowels; instead the sample was collected using a hand core sampler consisting of a two-inch-diameter, 20-inch-long stainless steel barrel with a plastic tube liner. Samples were extruded from the sample tube and into the sample containers.

In compliance with the *EAA Groundwater Quality Monitoring Plan* and consistent with the EAA practices of 2013, two field duplicates and two equipment blanks were collected. One field duplicate sample is required for each spring complex. The field duplicates were collected at the same locations as two of the field samples using the same methods as the field samples. Two equipment blanks were prepared in the laboratory of SWCA's San Antonio office. To collect one of the blanks, American Society for Testing and Materials (ASTM) Type II Reagent Grade water was poured through a new plastic sampling tube into sample collection containers. The second equipment blank was collected by pouring ASTM Type II Reagent Grade water over a decontaminated trowel into sample collection containers. The samples were containerized in the same manner as a surface water sample using the same types of containers and preservatives. Sample portions for metals analyses requiring field filtration were filtered using a 0.45-micron high capacity cartridge filter and disposable bailer. The equipment blanks were not analyzed for the following analytes: field parameters, turbidity, field alkalinity, and bacteria.

All samples were labeled and put on ice immediately upon collection for later shipment to the contract laboratory. Samples were secured inside locked SWCA vehicles during field operations and appropriate custody was maintained at all times. Representative photographs of field activities are included in Appendix E.

6.3 Stormwater Sampling Program

Stormwater samples are designated by the EAHCP Workplan (Appendix A) for collection twice annually from each spring complex. Stormwater samples were collected when rainfall amounts were adequate to initiate at least a five percent rise at the respective USGS gauging locations for each spring complex. Samples were collected across the storm-affected stream hydrograph at the rise, peak, and recession limb

of the associated stream hydrograph. As with the other sample types, five locations at Comal Springs and seven locations at San Marcos Springs were sampled. In general, the turbidity and conductivity data from the RTIs at each site were used as a surrogate for the stream hydrograph due to immediate availability of the data. Stream hydrograph data is only updated hourly on the USGS website. The RTI data is updated every fifteen minutes, which provides greater resolution regarding the effect of the storm event on the streams and facilitates quicker sampling response times. Graphs showing water quality parameters during each storm event are included in Appendix B.

Stormwater sample collection was affected by the ongoing drought conditions across the region during the second half of 2015. Often, storms that materialized were insufficient to create adequate runoff for sample collection. As a result, members of the stormwater sampling team spent many more hours on-call than were expected. In general, when rainfall probabilities exceeded 20 percent for a given time period, the team was placed on-call for sample collection. The team was mobilized when rainfall probabilities exceeded 50 percent. Multiple scenarios involved a mobilization that was later cancelled due to insufficient rainfall or storms that dissipated prior to reaching the sample area. Storm team duty is summarized and documented in Appendix F of this document.

Due to the inherently unsafe conditions associated with stormwater flow, SWCA field staff used disposable single-use bailers when needed in order to safely obtain water samples during stormwater sampling events. Field parameters were collected first by inserting the sonde probe as close to the sample location as possible. In May and October 2015, sample location HSM240 was sampled using disposable bailers. Bailers were lowered from the bridge above the sample location. A rope affixed to the bailer was used for retrieval. New bailers and rope were used for each sample point. After retrieval the water was transferred to the sample containers. Only new disposable equipment was used for stormwater sampling events.

Stormwater sampling efforts conformed to the protocols outlined in the *EAA Groundwater Quality Monitoring Plan* for sample collection, handling, and decontamination. Filtration for methods 6010B (metals), 6020 (metals), and field alkalinity were performed using a 0.45-micron high capacity cartridge filter and peristaltic pump. Preservatives were placed in the bottles (as needed) by the contract laboratory prior to sample collection. All samples were immediately placed into coolers with ice after sampling and later shipped to the contract laboratory. When not in use or after collection, sampling equipment and/or coolers containing samples were secured inside locked SWCA vehicles to maintain appropriate sample custody and security.

According to the *EAA Groundwater Quality Monitoring Plan*, two field duplicates were collected for the Comal Springs complex and three field duplicates for the San Marcos Springs complex per rain event. Field duplicates were sampled after collection of the parent sample and in the same manner as the field sample. No equipment blanks were required for stormwater samples as all equipment used was new and disposable.

Analyses for field alkalinity were performed at the field staging area or at SWCA's San Antonio office. The method used for field alkalinity is discussed in detail in the *EAA Groundwater Quality Monitoring Plan*. Representative photographs of field activities are included in Appendix E.

6.4 Surface Water Passive Samplers

The PDSs were deployed at each of the 12 sample locations during the months of February, April, June, August, October and December 2015. In general, PDS locations corresponded to surface water sampling points unless prevented by field conditions. Lost PDSs, human tampering, and any variations in deployment locations are discussed in Appendix C.

Deployment devices were constructed by SWCA staff at SWCA's San Antonio office in June 2014. Two-inch thick, 18-inch diameter concrete disks were poured and a stainless steel silverware cup was set approximately one inch deep in the center of the disk. Handles were formed by inserting both ends of an 18-inch length of vinyl-coated stainless steel cable into each side of the disk. Site numbers were marked in the wet concrete to dedicate each device to a sample location. The concrete was allowed to cure, and each device was decontaminated following the *EAA Groundwater Quality Monitoring Plan* guidelines and placed in a clean plastic bag prior to the first deployment. The same decontamination procedures were followed for subsequent sampling events. The deployment device is pictured in Figure 11.

Upon arrival at the sample location, the PDS was removed from a dedicated vial and affixed inside of a second stainless steel silverware cup with a plastic cable tie. This cup was inverted and placed on top of the cup set in the concrete sampling device enclosing the PDS inside the two cups. The two cups were secured to one another with additional plastic cable ties. The device was then gently lowered into the water. Installation date and time and PDS identification numbers were noted in the field notebook and on the PDS vial. To retrieve the PDS, the devices were simply removed from the water and the cable ties cut. The PDS was then immediately placed back in the dedicated vial and retrieval date and time were notated. Deployment devices were secured at SWCA offices when PDSs were not deployed.

Field duplicates were collected as directed by the *EAA Groundwater Quality Monitoring Plan*. To collect field duplicates, a second PDS was installed inside selected deployment devices. Field PDSs were always accompanied by test blank samplers to monitor for VOC contamination. Deployment devices were dedicated to each sample location to avoid cross contamination and were decontaminated following the *EAA Groundwater Quality Monitoring Plan* guidelines prior to each use. Representative photographs of field activities are included in Appendix E.

Figure 11. PDS deployment device prior to installation at site HSM440.



7.0 SAMPLE RESULTS

Results from the sampling efforts related to the EAHCP sampling program are discussed in the paragraphs that follow. Results are discussed by sample type for Comal Springs, followed by a separate discussion by sample type for San Marcos Springs. Sample events are listed in the order of surface water (base flow) samples, sediment samples, stormwater samples, and PDS. Laboratory analyses and field parameters are provided in Appendix G of this document. The laboratory data were reviewed by SWCA staff with the results of that review provided as Appendix H (data validation discussion) of this document. Each sample location (latitude/longitude), name, and other location information are also summarized in Appendix I of this document.

7.1 Comal Springs Sample Results

The Comal Springs complex was sampled for water quality during surface water base flow conditions in March and September 2015. In general, few detections were noted. As discussed previously, surface water base flow samples are compared to the drinking water standards for water quality in this report.

Sediments at the Comal Springs complex were sampled in June 2015. Sediment results are compared to the standards developed by McDonald et al. (2000) and TCEQ (2014a). These standards are based on the probability that a detected compound has a toxic effect on sediment-dwelling organisms and are referred to as the TEC and PEC. Detections below the TEC are not considered to be toxic, while detections above the PEC are considered to be toxic to sediment dwelling organisms. Detections above the TEC but less than the PEC are considered equally likely to be toxic or non-toxic.

Stormwater events were sampled at the Comal Springs complex in January and October 2015. Stormwater results did not indicate a significant number of detections of concern.

PDS sampling events were conducted at the Comal Springs complex in February, April, June, August, October, and December 2015. Generally speaking, various VOCs and total petroleum hydrocarbons (TPH) were detected at various sample locations, but only tetrachlorethene and TPH were relatively consistently detected.

7.1.1 Comal Springs Surface Water / Base Flow Sampling

The Comal Springs complex was sampled on March 16 and September 9, 2015, for surface water (base flow) events.

7.1.1.1 Surface Water / Base Flow - Bacteria

Bacteria results for surface water (base flow) associated with the Comal Springs complex ranged from 4 MPN/100 mL (most probable number of colony-forming units per 100 milliliters of water) to 800 MPN/100 mL for *E. coli*. Because of the presence of various fauna in surface water collection sites, positive detections are not uncommon. The 2014 Texas Surface Water Quality Standard for *E. coli* in primary recreation waters is a geometric mean of 126 MPN/100 mL with no individual sample exceeding 399 MPN/100 mL (30 TAC 307.7). The geometric mean for surface water samples collected from the Comal Springs complex during 2015 was 79.2 MPN/100 mL. Two samples exceeded the maximum limit of 399 MPN/100 mL, as described below.

- HCS110, sampled in September 2015, had a concentration of 800 MPN/mL
- HCS130, sampled in September 2015, had a concentration of 460 MPN/mL

Surface water (base flow) bacteria counts are summarized in Table 2.

Table 2. Surface Water Samples – Bacteria Counts - Comal Springs Complex

| Location | Date | Count (MPN/100 mL) |
|----------|-----------|--------------------|
| HCS110 | 3/16/2015 | 4 |
| | 9/9/2015 | 800 |
| HCS120 | 3/16/2015 | 47 |
| | 9/9/2015 | 67 |
| FDHCS120 | 3/16/2015 | 51 |
| | 9/9/2015 | 54 |
| HCS130 | 3/16/2015 | 59 |
| | 9/9/2015 | 460 |
| HCS140 | 3/16/2015 | 80 |
| | 9/9/2015 | 140 |
| HCS160 | 3/16/2015 | 60 |
| | 9/9/2015 | 120 |

MPN/100 mL – Most probable number per 100 milliliters of water.

7.1.1.2 Surface Water / Base Flow - Volatile Organic Compounds (VOCs)

No VOCs were detected at any of the five sampling sites from the Comal Springs complex during the March or September 2015 sampling events. However, there was a detection of VOCs in the VOC Trip Blank during the September sampling event. The Trip Blank tested positive for 0.25 micrograms per liter (µg/L) for toluene, which was below the laboratory reporting limit and not detected in any field samples.

7.1.1.3 Surface Water / Base Flow - Semi-volatile Organic Compounds (SVOCs)

Generally, Semi-volatile Organic Compounds (SVOCs) were analyzed because their detection can indicate the presence of chemicals originating from anthropogenic sources and therefore be used to evaluate potential impacts on water quality. No SVOCs were detected at any of the five sampling sites in the Comal Springs complex during the March or September 2015 sampling events.

7.1.1.4 Surface Water / Base Flow - Pesticides

Surface water samples were analyzed for pesticides because their detection can indicate the presence of chemicals originating from anthropogenic sources and therefore be used to evaluate potential impacts on water quality. No pesticides were detected in any of the samples collected for either the March or September 2015 sampling events at all five sites for the Comal Springs complex.

7.1.1.5 Surface Water / Base Flow - Herbicides

Surface water samples were analyzed for herbicides because their detection can indicate the presence of chemicals originating from anthropogenic sources and therefore be used to evaluate potential impacts on water quality. Herbicide analyses indicated no detections for both the March and September 2015 sampling events at all five sites for the Comal Springs complex.

7.1.1.6 Surface Water / Base Flow - Polychlorinated Biphenyls (PCBs)

Surface water samples were analyzed for the various Aroclor compounds that are collectively referred to as Polychlorinated Biphenyls (PCBs). PCB detection can indicate the presence of chemicals originating from anthropogenic sources and therefore, be used to evaluate potential impacts on water quality. No PCBs were detected during both the March and September 2015 sampling events at all five sites for the Comal Springs complex.

7.1.1.7 Surface Water / Base flow - Metals

Surface water samples were analyzed for metals that may indicate the presence of chemicals originating from anthropogenic sources. Although metals were detected for the March and September 2015 sampling events at all five sites for the Comal Springs complex, no metals were detected at a concentration in excess of the drinking water standards. The metals aluminum, antimony, arsenic, barium, chromium, copper, iron, lead, manganese, nickel, selenium, thallium and zinc were the detected metals of concern; however, none of their concentrations approached the MCL. These detections are listed below in Table 3, and note many detections are “J” flagged, indicating the detected concentration is below the laboratory reporting limit, but above the method detection limit. Also note that some metals are naturally occurring in rock, soil, groundwater, and surface water and may not indicate an anthropogenic source.

Table 3. Surface Water Samples – Metal detections - Comal Springs Complex

| Location | Date Collected | Antimony (mg/L) | Arsenic (mg/L) | Barium (mg/L) | Chromium (mg/L) | Copper (mg/L) | Lead (mg/L) | Nickel (mg/L) | Selenium (mg/L) | Thallium (mg/L) | Zinc (mg/L) | Aluminum (mg/L) | Iron (mg/L) | Manganese (mg/L) |
|----------|----------------|-----------------|----------------|---------------|-----------------|---------------|-------------|---------------|-----------------|-----------------|-------------|-----------------|-------------|------------------|
| HCS110 | 3/16/2015 | 0.000118 J | <0.000386 | 0.0371 | <0.000402 | 0.000547 J | <0.0000898 | 0.0017 | 0.000295 J | <0.000101 | 0.0245 | <0.00331 | 0.052 | 0.00243 |
| | 9/9/2015 | <0.0000995 | 0.000505 J | 0.0582 | 0.000492 J | 0.00457 | <0.0000898 | 0.00199 | 0.000503 J | <0.000101 | 0.0314 | 0.00923 J | 0.0424 J | 0.00481 |
| HCS120 | 3/16/2015 | <0.0000995 | <0.000386 | 0.0555 | <0.000402 | 0.000538 J | <0.0000898 | 0.00201 | 0.000334 J | <0.000101 | 0.0239 | <0.00331 | 0.046 J | 0.00178 |
| | 9/9/2015 | <0.0000995 | 0.000388 J | 0.0549 | 0.000524 J | 0.00504 | <0.0000898 | 0.00182 | 0.000521 J | 0.000121 J | 0.00984 | 0.00699 J | 0.0364 J | 0.00082 J |
| FDHCS120 | 3/16/2015 | 0.000166 J | <0.000386 | 0.0562 | 0.000426 J | 0.000783 J | 0.000117 J | 0.00225 | 0.000414 J | <0.000101 | 0.0391 | 0.00459 J | 0.0555 | 0.00607 |
| | 9/9/2015 | <0.0000995 | <0.000386 | 0.0537 | 0.000482 J | 0.00078 J | <0.0000898 | 0.00187 | 0.000932 J | <0.000101 | 0.00773 | 0.00954 J | 0.0406 J | 0.00064 J |
| HCS130 | 3/16/2015 | <0.0000995 | <0.000386 | 0.0574 | <0.000402 | 0.00057 J | <0.0000898 | 0.00184 | 0.000347 J | <0.000101 | 0.00384 J | <0.00331 | 0.0468 J | 0.00459 |
| | 9/9/2015 | <0.0000995 | <0.000386 | 0.0553 | 0.000555 J | 0.0016 | <0.0000898 | 0.00204 | 0.000744 J | 0.000109 J | 0.0159 | 0.00749 J | 0.0465 J | 0.00114 |
| HCS140 | 3/16/2015 | <0.0000995 | <0.000386 | 0.0584 | <0.000402 | 0.000907 J | 0.0000905 J | 0.00197 | 0.000446 J | <0.000101 | 0.0118 | <0.00331 | 0.0399 J | 0.00632 |
| | 9/9/2015 | 0.000106 J | <0.000386 | 0.0555 | 0.000583 J | 0.00276 | 0.000112 J | 0.00228 | 0.00229 | <0.000101 | 0.035 | 0.00923 J | 0.0451 J | 0.00137 |
| HCS160 | 3/16/2015 | 0.000107 J | <0.000386 | 0.0598 | <0.000402 | 0.000607 J | <0.0000898 | 0.00216 | 0.000334 J | <0.000101 | 0.0265 | <0.00331 | 0.0508 | 0.00991 |
| | 9/9/2015 | <0.0000995 | <0.000386 | 0.055 | 0.000645 J | 0.0066 | <0.0000898 | 0.00235 | 0.00193 | <0.000101 | 0.00714 | 0.00801 J | 0.0399 J | 0.00269 |
| MCL | | 0.006 | 0.01 | 2 | 0.1 | NE | NE | NE | 0.05 | 0.002 | NE | NE | NE | NE |
| PCL | | -- | -- | -- | -- | 1.3 | 0.015 | 0.49 | -- | -- | 7.3 | 24 | NE | 1.1 |

-- – Not applicable
J – Detection is above the method detection limit, but below the reporting limit
MCL – maximum contaminant level
mg/L – milligrams per liter
NE – Not established
MCL – maximum contaminant level
PCL – protective concentration levels

7.1.1.8 Surface Water / Base Flow - Nitrates

Surface water samples were analyzed for nitrate-nitrite as nitrogen. Laboratory analyses indicated a limited range of nitrate-nitrite as nitrogen in surface water samples. Of the 12 surface water samples (ten environmental samples and two field duplicates) collected for the two sample events, concentrations ranged from 0.35 milligrams per liter (mg/L) to 1.8 mg/L. None of the nitrate concentrations detected exceeded the MCL of 10 mg/L for drinking water. The highest nitrate concentration in surface water at the Comal Springs complex, 1.8 mg/L, was detected in five samples. These five samples were HCS130 sampled on March 16, 2015, and HCS120, FDHCS120, HCS130 and HCS160 all collected on September 9, 2015. Nitrate-nitrogen results are summarized in Table 4.

**Table 4. Surface Water Samples – Nitrate Detections
- Comal Springs Complex**

| Location | Date | Concentration (mg/L) |
|----------|-----------|-------------------------|
| HCS110 | 3/16/2015 | 0.35 |
| | 9/9/2015 | 1.2 |
| HCS120 | 3/16/2015 | 1.7 |
| | 9/9/2015 | 1.8 |
| FDHCS120 | 3/16/2015 | 1.7 |
| | 9/9/2015 | 1.8 |
| HCS130 | 3/16/2015 | 1.8 |
| | 9/9/2015 | 1.8 |
| HCS140 | 3/16/2015 | 1.7 |
| | 9/9/2015 | 1.7 |
| HCS160 | 3/16/2015 | 1.7 |
| | 9/9/2015 | 1.8 |

mg/L – milligrams per liter

7.1.1.9 Surface Water / Base Flow – Caffeine

Surface water base flows were analyzed for caffeine, which can indicate an anthropogenic source. Caffeine may enter surface water from leaking sewer or septic systems or it may be present in the aquifer from similar sources in the recharge zone (EPA, 2012). Potential ecological effects are currently unknown but could include reduced reproductive success in aquatic species (EPA, 2012). Caffeine detections in surface water samples from Comal Springs in 2015 ranged from 16 to 41 nanograms per liter (ng/L). Caffeine was only detected in surface water samples from two locations in the Comal Springs system in 2015, HCS110 in March and September and the field duplicate collected at HCS120 in September. There is no regulatory standard or expected value for comparison. Results are shown in Table 5.

**Table 5. Surface Water Samples – Caffeine
Detections - Comal Springs Complex**

| Location | Date Collected | Caffeine (ng/L) |
|-----------------|---------------------------|----------------------------|
| HCS110 | 3/16/2015 | 39 |
| | 9/9/2015 | 41 |
| HCS120 | 3/16/2015 | <3.5 |
| | 9/9/2015 | <4 |
| FDHCS120 | 3/16/2015 | <3.5 |
| | 9/9/2015 | 16 |
| HCS130 | 3/16/2015 | <3.5 |
| | 9/9/2015 | <4 |
| HCS140 | 3/16/2015 | <3.5 |
| | 9/9/2015 | <4 |
| HCS160 | 3/16/2015 | <3.5 |
| | 9/9/2015 | <4 |

ng/L – nanograms per liter

7.1.2 Comal Springs Sediment Sampling

7.1.2.1 Sediment - Volatile Organic Compounds (VOCs)

No VOCs were detected in any of the sediment samples collected at Comal Springs during 2015.

7.1.2.2 Sediment - Semi-volatile Organic Compounds (SVOCs)

Only one SVOC compound was detected in the sediment samples collected in the Comal Springs system in 2015. The compound, bis(2-ethylhexyl) phthalate, was detected at HCS320 at a concentration of 0.94mg/kg. There is no PEC or TEC established for this compound.

7.1.2.3 Sediment - Pesticides

Sediment samples were analyzed for both organochlorine and organophosphorous pesticides. No pesticides were detected in any of the sediment samples collected in the Comal Springs complex.

7.1.2.4 Sediment - Herbicides

Sediments were analyzed for herbicide compounds to further assess sediment quality at the Comal Springs complex. No herbicides were detected in any of the sediment samples collected from the Comal Springs complex.

7.1.2.5 Sediment - Polychlorinated Biphenyls

Sediments were analyzed for PCB compounds to further assess sediment quality at the Comal Springs complex. No PCB compounds were detected in any of the sediment samples collected from the Comal Springs complex.

7.1.2.6 Sediment - Metals

Many metals are naturally occurring within soil, rock, and sediment. Sediment sample results for the Comal Springs complex tested positive for several metals, generally at low concentrations. Metals detected above the method detection limit and subsequently evaluated in this report for potential toxic effects using the TEC and/or PEC standards are: arsenic, chromium, copper, iron, lead, manganese, nickel, and zinc. Other metals detected that do not have a TEC or PEC value available are aluminum and barium. These metals (aluminum and barium) were compared to Texas-specific background concentrations (TSBC) (TCEQ, 2014b) for soil. None of these exceeded the listed background concentration.

Zinc exceeded the TEC at HCS360 and in the field duplicate, FDHCS360. Metal detections are listed in Table 6 below. Zinc detections above the established TEC value are displayed graphically in Figure 12. No metals were detected above a PEC.

Figure 12. Zinc in Comal Springs Complex Sediments Compared to the Threshold Effect Concentration (TEC) and Probable Effect Concentration (PEC) Values

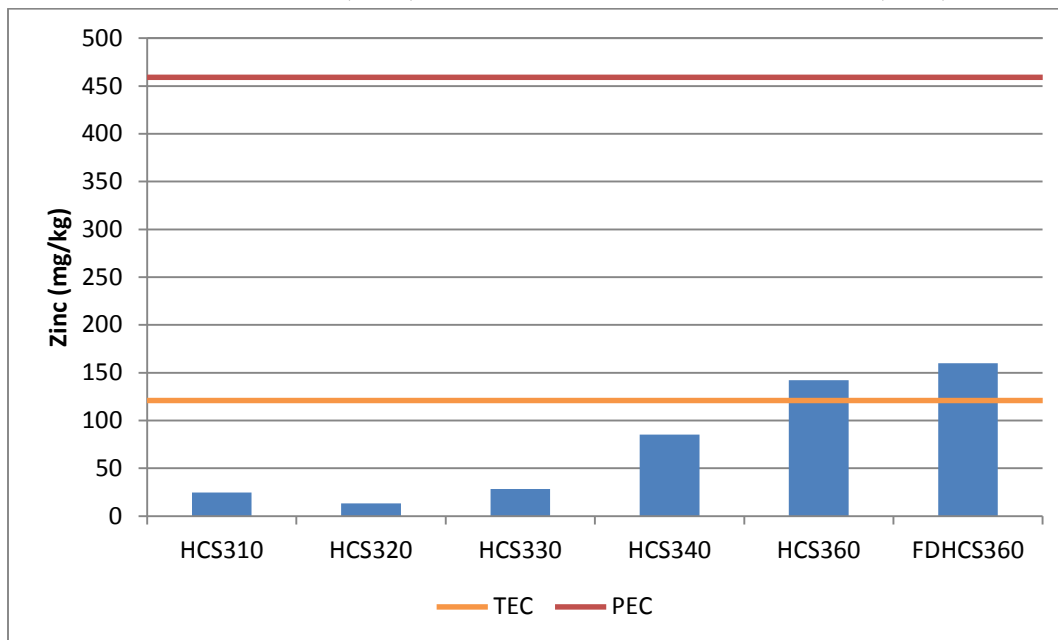


Table 6. Sediment Samples – Metal Detections - Comal Springs Complex

| | Date | Arsenic | Barium | Chromium | Copper | Lead | Nickel | Zinc | Aluminum | Iron | Manganese |
|-----------------------|-----------|---------|---------|----------|---------|---------|---------|---------|----------|---------|-----------|
| Location | Collected | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) |
| HCS310 | 6/4/2015 | <0.482 | 16 | 3.38 | 3.65 | 5.53 | 14.5 | 24.6 | 1390 | 1710 | 31.6 |
| HCS320 | 6/4/2015 | <0.549 | 23.5 | 6.41 | 3.55 | 6.33 | 19.1 | 13.1 | 1580 | 2620 | 47.8 |
| HCS330 | 6/4/2015 | 4.92 | 61 | 15.3 | 9.74 | 10.7 | 20.5 | 28.2 | 10300 | 10300 | 195 |
| HCS340 | 6/4/2015 | 3.14 | 67.9 | 11.7 | 12.9 | 13.5 | 21.1 | 85.1 | 4880 | 9730 | 108 |
| HCS360 | 6/4/2015 | 3.83 | 68.9 | 16.1 | 14.3 | 23.9 | 17 | 142 | 10800 | 9750 | 216 |
| FDHCS360 | 6/4/2015 | 4.15 | 72.9 | 18 | 16 | 26.2 | 18.4 | 160 | 11900 | 10700 | 227 |
| TEC | | 9.79 | NE | 43.4 | 31.6 | 35.8 | 22.7 | 121 | NE | 20000 | 460 |
| PEC | | 33 | NE | 111 | 149 | 128 | 48.6 | 459 | NE | 40000 | 1100 |
| Soil background level | | 5.9 | 300 | 30 | 15 | 15 | 10 | 30 | 30000 | 15000 | 300 |

mg/kg – milligrams per kilograms

PEC – probable effect concentration

TEC – threshold effect concentration

7.1.3 Comal Springs Stormwater Sampling

Stormwater samples were collected during two storm events at the Comal Springs complex. The events were sampled according to the guidelines in the EAHCP workplan. The events occurred on January 22–23, 2015, and October 23, 2015. Total rainfall for the January 2015 event was approximately 2.00 to 2.49 inches (National Oceanic and Atmospheric Administration [NOAA], 2015) causing streamflow measured at USGS Gauge 08169000 to increase from approximately 141 cfs to a peak of 420 cfs (USGS, 2015). Total rainfall for the October 2015 event was approximately 2.00 to 2.49 inches in the 24-hour period beginning at 7:00 AM on October 23, 2015 (NOAA, 2015). Sampling efforts occurred during the first pulse of the storm when precipitation is estimated to have been approximately 1 inch. During sampling efforts, streamflow measurements from the USGS gauge increased from approximately 213 cfs to a peak of 314 cfs (USGS, 2015).

7.1.3.1 Stormwater - Bacteria Detections

Stormwater samples collected and analyzed for bacteria analyses generally tested positive for high levels of bacteria. Bacterial analyses were performed for *E. coli*, using a most probable number method. The 2014 Texas Surface Water Quality Standard for *E. coli* in primary recreation waters is a geometric mean of 126 MPN/100 mL with no individual sample exceeding 399 MPN/100 mL (30 TAC 307.7). The geometric mean for stormwater samples collected from the Comal Springs complex during January 2015 was 1342.5 MPN/100 mL. Bacteria counts from January 2015 ranged from 110 MPN/100 mL to 10,000 MPN/100 mL with several samples exceeding the individual sample limit. The geometric mean for stormwater samples collected from the Comal Springs complex during October 2015 was 17,905.4 MPN/100 mL. Bacteria counts from October 2015 ranged from 2000 MPN/100 mL to 77,000 MPN/100 mL with all samples exceeding the individual sample limit. Individual detections are listed below in Table 7 and shown in relation to stream discharge and specific conductivity in Figures 13 and 14. During October 2015, HCS210 Peak was not analyzed due to insufficient sample volume after failure of the container seal during transit to the laboratory. Due to the timing of storm events and laboratory working hours, it was not possible to deliver all samples to the laboratory within sample holding time of 8 hours, as discussed in Appendix C. These samples were included in the range and geometric mean calculations.

**Table 7. Stormwater Samples – Bacteria Counts -
Comal Springs Complex**

| Location | Date | Count (MPN/100 mL) |
|----------------|------------|-----------------------|
| HCS210 Lead | 1/22/2015 | 2200 H |
| | 10/23/2015 | 77000 |
| HCS210 Peak | 1/22/2015 | 1600 H |
| | 10/23/2015 | NA |
| HCS210 Trail | 1/23/2015 | 880 H |
| | 10/23/2015 | 69000 H |
| HCS240 Lead | 1/22/2015 | 910 H |
| | 10/23/2015 | 20000 |
| HCS240 Peak | 1/22/2015 | 550 H |
| | 10/23/2015 | 2000 H |
| HCS240 Trail | 1/23/2015 | 110 H |
| | 10/23/2015 | 2900 H |
| HCS250 Lead | 1/22/2015 | 630 H |
| | 10/23/2015 | 25000 |
| HCS250 Peak | 1/22/2015 | 1800 H |
| | 10/23/2015 | 17000 H |
| HCS250 Trail | 1/23/2015 | 340 H |
| | 10/23/2015 | 16000 H |
| HCS260 Lead | 1/22/2015 | 370 H |
| | 10/23/2015 | 17000 |
| HCS260 Peak | 1/22/2015 | 10000 H |
| | 10/23/2015 | 20000 H |
| HCS260 Trail | 1/23/2015 | 2300 H |
| | 10/23/2015 | 20000 H |
| FDHCS260 Trail | 1/23/2015 | 3400 H |
| | 10/23/2015 | 24000 H |
| HCS270 Lead | 1/22/2015 | 1300 H |
| | 10/23/2015 | 24000 |
| HCS270 Peak | 1/22/2015 | 5500 H |
| | 10/23/2015 | 20000 H |
| HCS270 Trail | 1/23/2015 | 4400 H |
| | 10/23/2015 | 20000 H |
| FDHCS270 Trail | 1/23/2015 | 2500 H |
| | 10/23/2015 | 17000 H |

H – Analyzed outside hold time, result included for comparison but not considered valid

MPN/100 mL – Most probable number per 100 milliliters of water

NA – Not analyzed

Figure 13. Stormwater Samples – January 2015 Bacteria Counts in Relation to Stream Discharge and Specific Conductivity - Comal Springs Complex

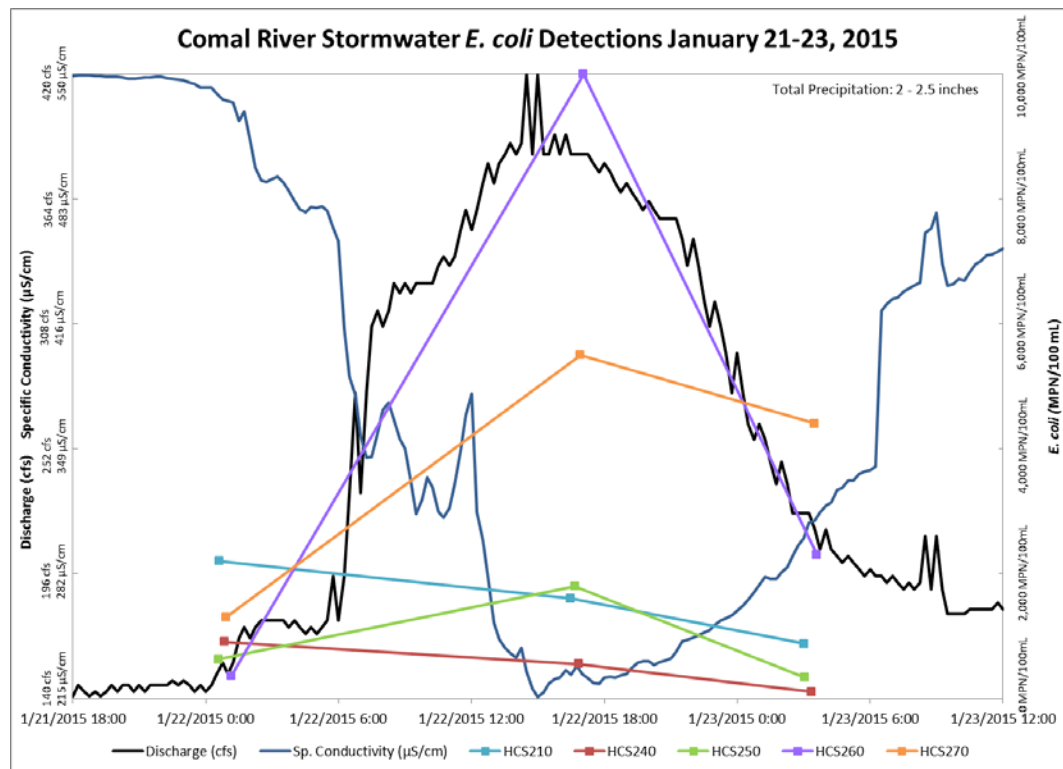
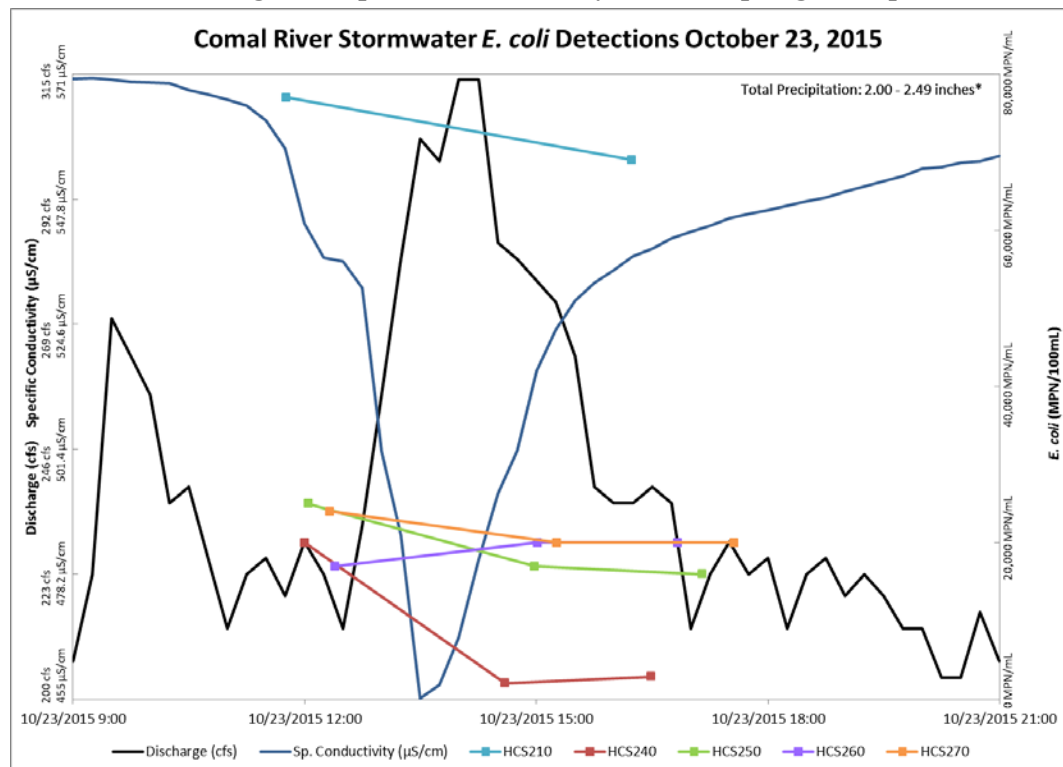


Figure 14. Stormwater Samples – October 2015 Bacteria Counts in Relation to Stream Discharge and Specific Conductivity - Comal Springs Complex



7.1.3.2 Stormwater - Volatile Organic Compounds (VOCs)

VOCs were detected in two samples, HCS210 Lead and HCS240 Lead, in the Comal Springs storm system during the January 2015 event. Acetone and 2-butanone were detected at HCS210 and carbon disulfide was detected at HCS240. All three detections were below the laboratory reporting limit and the established PCLs. Acetone was also detected in sample HCS210 Peak during the October storm event. This detection was also below the laboratory reporting limit and the established PCL. Detections are shown in Table 8.

Table 8. Stormwater Samples – Volatile Organic Compound Detections - Comal Springs Complex

| Location | Date Collected | Acetone (µg/L) | 2-Butanone (µg/L) | Carbon disulfide (µg/L) |
|-------------|----------------|-------------------|----------------------|----------------------------|
| HCS210 Lead | 1/22/2015 | 11 J | 3.7 J | <0.41 |
| HCS210 Peak | 10/23/2015 | 12 J | <2.2 | <0.41 |
| HCS240 Lead | 1/22/2015 | <10 | <2.2 | 0.41 J |
| MCL | | NE | NE | NE |
| PCL | | 22000 | 15000 | 2400 |

J – Detection is above the method detection limit, but below the reporting limit

µg/L – milligrams per Liter

MCL – maximum contaminant level

NE – Not established

PCL – protective concentration levels

7.1.3.3 Stormwater - Semi-volatile Organic Compounds (SVOCs)

Stormwater samples collected during January 2015 were analyzed for SVOCs and were generally non-detect for these compounds, with the exception of one sample, HCS270 Peak, which tested positive for butyl benzyl phthalate at a concentration of 3.4 µg. This detection is below the laboratory reporting limit and the established drinking water PCL. No SVOCs were detected during the October 2015 event.

7.1.3.4 Stormwater - Herbicides and Pesticides

No herbicides or organophosphorous pesticides were detected for the storm event in the Comal Springs complex sampled during January or October 2015.

7.1.3.5 Stormwater - Polychlorinated Biphenyls (PCBs)

Stormwater samples were analyzed for the various Aroclor compounds that are generally referred to collectively as PCBs. None of the stormwater samples from the Comal Springs complex indicated positive detections of PCB compounds during the January or October 2015 sampling events.

7.1.3.6 Stormwater - Metals

Stormwater samples were analyzed for metals in accordance with the EAHCP workplan. Several positive metal detections were noted in the sample set; however, no samples contained a metal at a concentration in excess of the drinking water MCL and most detections were below laboratory reporting limits.

7.1.3.7 Stormwater - Nitrates

Stormwater samples were analyzed for nitrate-nitrite as nitrogen in accordance with the EAHCP workplan. All nitrate results were below the MCL of 10 mg/L. For the January 2015 event, the range of nitrate results was 0.21 mg/L to 1.8 mg/L, with an average of 1.2 mg/L. During the October 2015 event, nitrate concentrations ranged from 0.44 mg/L to 1.8 mg/L, with an average of 1.4 mg/L. For comparison the average nitrate in spring water samples at Comal Springs for calendar year 2014 was 1.91 mg/L (EAA, 2015). Nitrate results are summarized in Table 9.

Table 9. Stormwater Samples – Nitrate Detections - Comal Springs Complex

| Location | Date | Concentration (mg/L) |
|----------------|------------|----------------------|
| HCS210 Lead | 1/22/2015 | 0.55 |
| | 10/23/2015 | 0.44 |
| HCS210 Peak | 1/22/2015 | 0.26 |
| | 10/23/2015 | 0.61 |
| HCS210 Trail | 1/23/2015 | 0.21 |
| | 10/23/2015 | 0.68 H |
| HCS240 Lead | 1/22/2015 | 1.7 |
| | 10/23/2015 | 1.2 |
| HCS240 Peak | 1/22/2015 | 1.7 |
| | 10/23/2015 | 1.7 |
| HCS240 Trail | 1/23/2015 | 1.8 |
| | 10/23/2015 | 1.8 H |
| HCS250 Lead | 1/22/2015 | 1.6 |
| | 10/23/2015 | 1.3 |
| HCS250 Peak | 1/22/2015 | 1.5 |
| | 10/23/2015 | 1.6 |
| HCS250 Trail | 1/23/2015 | 1.7 |
| | 10/23/2015 | 1.6 H |
| HCS260 Lead | 1/22/2015 | 1.7 |
| | 10/23/2015 | 1.7 |
| HCS260 Peak | 1/22/2015 | 0.94 |
| | 10/23/2015 | 1.6 H |
| HCS260 Trail | 1/23/2015 | 1.2 |
| | 10/23/2015 | 1.7 H |
| FDHCS260 Trail | 1/23/2015 | 1.2 |
| | 10/23/2015 | 1.7 H |

Table 9. Stormwater Samples – Nitrate Detections - Comal Springs Complex

| Location | Date | Concentration (mg/L) |
|----------------|------------|----------------------|
| HCS270 Lead | 1/22/2015 | 1.6 |
| | 10/23/2015 | 1.4 |
| HCS270 Peak | 1/22/2015 | 0.98 |
| | 10/23/2015 | 1.6 H |
| HCS270 Trail | 1/23/2015 | 1.2 |
| | 10/23/2015 | 1.6 H |
| FDHCS270 Trail | 1/23/2015 | 1.2 |
| | 10/23/2015 | 1.6 H |

H – Analyzed outside hold time, result included for comparison but not considered valid

mg/L – milligrams per liter

7.1.3.8 Stormwater – Caffeine

Stormwater was analyzed for caffeine, which can indicate an anthropogenic source. Caffeine may enter surface water from leaking sewer or septic systems or it may be present in the aquifer from similar sources in the recharge zone (EPA, 2012). Potential ecological effects are currently unknown but could include reduced reproductive success of aquatic organisms (EPA, 2012). Caffeine detections in stormwater samples from Comal Springs in January 2015 ranged from 11 ng/L to 480 ng/L. In October 2015, caffeine detections ranged from 11 ng/L to 850 ng/L. There is no regulatory standard or expected value for comparison. These results are shown in Table 10.

Table 10. Stormwater Samples – Caffeine Detections - Comal Springs Complex

| Location | Date Collected | Caffeine (ng/L) |
|--------------|----------------|-----------------|
| HCS210 Lead | 1/22/2015 | 480 |
| | 10/23/2015 | 430 |
| HCS210 Peak | 1/22/2015 | 95 |
| | 10/23/2015 | 180 |
| HCS210 Trail | 1/23/2015 | 140 |
| | 10/23/2015 | 150 |
| HCS240 Lead | 1/22/2015 | 47 |
| | 10/23/2015 | 120 |
| HCS240 Peak | 1/22/2015 | <3.6 |
| | 10/23/2015 | 26 |
| HCS240 Trail | 1/23/2015 | 18 |
| | 10/23/2015 | 11 |

Table 10. Stormwater Samples – Caffeine Detections - Comal Springs Complex

| Location | Date Collected | Caffeine (ng/L) |
|----------------|----------------|-----------------|
| HCS250 Lead | 1/22/2015 | 55 |
| | 10/23/2015 | 850 |
| HCS250 Peak | 1/22/2015 | 38 |
| | 10/23/2015 | 120 |
| HCS250 Trail | 1/23/2015 | 11 |
| | 10/23/2015 | 81 |
| HCS260 Lead | 1/22/2015 | 57 |
| | 10/23/2015 | 160 |
| HCS260 Peak | 1/22/2015 | 130 |
| | 10/23/2015 | 120 |
| HCS260 Trail | 1/23/2015 | 110 |
| | 10/23/2015 | 80 |
| FDHCS260 Trail | 1/23/2015 | 67 |
| | 10/23/2015 | 56 |
| HCS270 Lead | 1/22/2015 | 95 |
| | 10/23/2015 | 220 |
| HCS270 Peak | 1/22/2015 | 130 |
| | 10/23/2015 | 210 |
| HCS270 Trail | 1/23/2015 | 100 |
| | 10/23/2015 | 450 |
| FDHCS270 Trail | 1/23/2015 | 65 |
| | 10/23/2015 | 130 |

ng/L – nanograms per liter

7.1.4 Comal Springs Surface Water Passive Sampling

PDSs were installed in the Comal Springs system in February, April, June, August, October and December 2015. The PDS was not recovered from site HCS460 in April or December 2015. Any changes to deployment locations or non-recovered samplers are discussed in Appendix C. There is not a suitable set of regulatory standards to compare PDS results to, but rather the data are a qualitative tool for evaluating the presence of trace concentrations of organic compounds. PDSs were analyzed for a suite of SVOCs, VOCs, and organochlorine pesticides. Few compounds were detected, the most notable are relatively consistent detections of tetrachlorethene and TPH in several samples. Positive detections are shown in Table 11.

Table 11. Passive Diffusion – Volatile Organic Compounds (VOCs) - Comal Springs Complex

| Location | Month 2015 | BTEX (µg) | Chloroform (µg) | Tetrachloroethene (µg) | p/m-Xylene (µg) | o-Xylene (µg) | Toluene (µg) | TPH (µg) | Undecane (µg) |
|----------|------------|--------------|--------------------|---------------------------|--------------------|------------------|-----------------|-------------|------------------|
| HCS410 | February | <0.02 | <0.02 | 0.05 | <0.02 | <0.02 | <0.02 | 0.76 | <0.05 |
| | April | <0.02 | <0.02 | 0.02 | <0.02 | <0.02 | <0.02 | 1.82 | <0.05 |
| | June | <0.02 | 0.03 | 0.14 | <0.02 | <0.02 | <0.02 | 0.66 | <0.05 |
| | August | <0.02 | <0.02 | 0.09 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| | October | <0.02 | 0.02 | 0.07 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| | December | <0.02 | 0.03 | 0.11 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| HCS420 | February | <0.02 | <0.02 | 0.28 | <0.02 | <0.02 | <0.02 | 0.52 | <0.05 |
| | April | <0.02 | <0.02 | 0.34 | <0.02 | <0.02 | <0.02 | 1.02 | 0.16 |
| | June | <0.02 | <0.02 | 0.20 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| | August | <0.02 | <0.02 | 0.13 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| | October | <0.02 | <0.02 | 0.21 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| | December | 0.11 | <0.02 | 0.38 | <0.02 | <0.02 | 0.11 | <0.50 | <0.05 |
| HCS430 | February | <0.02 | <0.02 | 0.29 | <0.02 | <0.02 | <0.02 | 1.42 | <0.05 |
| | April | <0.02 | <0.02 | 0.42 | <0.02 | <0.02 | <0.02 | 1.26 | <0.05 |
| | June | <0.02 | <0.02 | 0.60 | <0.02 | <0.02 | <0.02 | 0.54 | <0.05 |
| | August | <0.02 | <0.02 | 0.36 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| | October | <0.02 | <0.02 | 0.55 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| | December | <0.02 | <0.02 | 0.58 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |

Table 11. Passive Diffusion – Volatile Organic Compounds (VOCs) - Comal Springs Complex

| Location | Month 2015 | BTEX (µg) | Chloroform (µg) | Tetrachloroethene (µg) | p/m-Xylene (µg) | o-Xylene (µg) | Toluene (µg) | TPH (µg) | Undecane (µg) |
|----------|------------|--------------|--------------------|---------------------------|--------------------|------------------|-----------------|-------------|------------------|
| HCS440 | February | <0.02 | <0.02 | 0.28 | <0.02 | <0.02 | <0.02 | 0.61 | <0.05 |
| | April | <0.02 | <0.02 | 0.41 | <0.02 | <0.02 | <0.02 | 0.65 | <0.05 |
| | June | <0.02 | <0.02 | 0.40 | <0.02 | <0.02 | <0.02 | 0.72 | <0.05 |
| | August | <0.02 | <0.02 | 0.39 | <0.02 | <0.02 | <0.02 | 1.23 | <0.05 |
| | October | <0.02 | <0.02 | 0.47 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| | December | <0.02 | <0.02 | 0.38 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| FDHCS440 | February | <0.02 | <0.02 | 0.28 | <0.02 | <0.02 | <0.02 | 0.76 | <0.05 |
| | April | <0.02 | <0.02 | 0.37 | <0.02 | <0.02 | <0.02 | 1.05 | <0.05 |
| | June | <0.02 | <0.02 | 0.42 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| | August | <0.02 | <0.02 | 0.44 | <0.02 | <0.02 | <0.02 | 1.44 | <0.05 |
| | October | <0.02 | <0.02 | 0.46 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| | December | <0.02 | <0.02 | 0.39 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| HCS460 | February | <0.02 | <0.02 | 0.18 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| | April | NA | NA | NA | NA | NA | NA | NA | NA |
| | June | 0.03 | <0.02 | 0.26 | 0.03 | <0.02 | <0.02 | 1.90 | <0.05 |
| | August | 0.11 | <0.02 | 0.31 | 0.08 | 0.04 | <0.02 | 2.59 | <0.05 |
| | October | <0.02 | <0.02 | 0.30 | <0.02 | <0.02 | <0.02 | <0.50 | <0.05 |
| | December | NA | NA | NA | NA | NA | NA | NA | NA |

BTEX – benzene, toluene, ethylbenzene, and xylenes

NA – Not analyzed

µg – micrograms

TPH – total petroleum hydrocarbons

7.2 San Marcos Springs Sample Results

The surface waters associated with the San Marcos Springs complex were sampled for surface water base flow conditions in March and September 2015. In general, few detections were noted. As discussed previously, surface water samples are compared to the drinking water standards for water quality in this report.

Sediments at the San Marcos Springs complex were sampled in June 2015. Sediment results were compared to the standards developed by McDonald et al. (2000) and TCEQ (2014a). These standards are based on the probability of a detected compound having a toxic effect on sediment dwelling organisms and are referred to as the TEC and PEC. Detections below the TEC are not considered to be toxic, while detections above the PEC are considered to be toxic to sediment-dwelling organisms. Detections above the TEC but less than the PEC are considered to be equally likely to be toxic or non-toxic.

Stormwater events were sampled at the San Marcos Springs complex in May and October 2015. Stormwater results did not indicate a significant number of detections of concern.

PDS sampling events were conducted at the San Marcos Springs complex in February, April, June, August, October, and December 2015. Generally speaking, various VOCs and TPH were detected at various sample locations, but only tetrachlorethene and TPH were relatively consistently detected.

7.2.1 San Marcos Springs Surface Water / Base Flow Sampling

The San Marcos Springs complex was sampled on March 25 and September 17, 2015, for surface water (base flow) events.

7.2.1.1 Surface Water / Base flow - Bacteria

Bacteria results for surface water (base flow) associated with the San Marcos Springs complex ranged from 9 MPN/100 mL through 200 MPN/100 mL for *E. coli*. Because of the presence of various fauna in surface water collection sites, positive detections are common. The 2014 Texas Surface Water Quality Standard for *E. coli* in primary recreation waters is a geometric mean of 126 MPN/100 mL with no individual sample exceeding 399 MPN/100 mL (30 TAC 307.7). The geometric mean for surface water samples collected from the San Marcos Springs complex during 2015 was 31.7 MPN/100 mL. No surface water samples collected from the San Marcos Springs in 2015 exceeded the individual sample limit of 399 MPN/100 mL. Surface water (base flow) bacteria counts are summarized in Table 12.

Table 12. Surface Water Samples – Bacteria Counts - San Marcos Springs Complex

| Location | Date | Count (MPN/100 mL) |
|----------|-----------|--------------------|
| HSM110 | 3/25/2015 | <1 |
| | 9/17/2015 | 45 |
| FDHSM110 | 3/25/2015 | 9 |
| | 9/17/2015 | 75 |
| HSM120 | 3/25/2015 | 13 |
| | 9/17/2015 | 20 |
| HSM130 | 3/25/2015 | 12 |
| | 9/17/2015 | 30 |
| HSM140 | 3/25/2015 | 200 |
| | 9/17/2015 | 37 |
| HSM150 | 3/25/2015 | 66 |
| | 9/17/2015 | 22 |
| HSM160 | 3/25/2015 | 40 |
| | 9/17/2015 | 28 |
| HSM170 | 3/25/2015 | 34 |
| | 9/17/2015 | 28 |

MPN/100 mL – most probable number per 100 milliliters

7.2.1.2 Surface Water / Base Flow - Volatile Organic Compounds (VOCs)

No VOCs were detected in the San Marcos Springs surface water samples in March or September 2015.

7.2.1.3 Surface Water / Base Flow - Semi-volatile Organic Compounds (SVOCs)

Surface water samples were analyzed for SVOCs because their detection can indicate the presence of chemicals originating from anthropogenic sources and therefore be used to evaluate potential impacts on water quality. No SVOCs were detected in any of the surface water samples from the San Marcos Springs complex during March and September 2015.

7.2.1.4 Surface Water / Base Flow - Pesticides

Surface water samples were analyzed for pesticides because their detection can indicate the presence of chemicals originating from anthropogenic sources and therefore be used to evaluate potential impacts on water quality. No pesticides were detected in any of the San Marcos Springs complex surface water samples during March and September 2015.

7.2.1.5 Surface Water / Base flow - Herbicides

Surface water samples were analyzed for herbicides because their detection can indicate the presence of chemicals originating from anthropogenic sources and therefore be used to evaluate potential impacts on

water quality. Herbicides were not detected for the March or September 2015 sampling events at all seven sites for the San Marcos Springs complex.

7.2.1.6 Surface Water / Base Flow - Polychlorinated Biphenyls (PCBs)

Surface water samples were analyzed for the various Aroclor compounds that are generally referred to collectively as PCBs. PCBs are sampled because their detection can indicate the presence of chemicals originating from anthropogenic sources and therefore help in the evaluation of potential impacts on water quality. PCBs were not detected for the March or September 2015 sampling events at all seven sites for the San Marcos Springs complex.

7.2.1.7 Surface Water / Base Flow - Metals

Surface water samples were analyzed for metals because their detection can indicate the presence of chemicals originating from anthropogenic sources and therefore be used to evaluate potential impacts on water quality. Although metals were detected for both the March and September 2015 sampling events at all seven sites for the San Marcos Springs complex, no metals of concern were detected at concentrations in excess of the drinking water standards. Barium, copper, iron, manganese, nickel, selenium, and zinc were detected; however, none of their concentrations exceeded a regulatory standard. These detections are listed in Table 13. Note many detections are “J” flagged, indicating the detected concentration is below the laboratory reporting limit, but above the method detection limit.

Table 13. Surface Water Samples – Metal Detections - San Marcos Springs Complex

| Location | Date Collected | Barium (mg/L) | Copper (mg/L) | Nickel (mg/L) | Selenium (mg/L) | Zinc (mg/L) | Iron (mg/L) | Manganese (mg/L) |
|----------|----------------|---------------|---------------|---------------|-----------------|-------------|-------------|------------------|
| HSM110 | 3/25/2015 | 0.0409 | 0.000788 J | 0.00193 | 0.00031 J | 0.00287 J | 0.0411 J | 0.0222 |
| | 9/17/2015 | 0.0407 | 0.00175 J | 0.00176 J | <0.000841 | <0.00239 | <0.0463 | 0.283 |
| FDHSM110 | 3/25/2015 | 0.0429 | 0.000673 J | 0.00203 | <0.000168 | <0.000479 | 0.0485 J | 0.022 |
| | 9/17/2015 | 0.0394 | 0.00216 J | 0.00168 J | <0.000841 | <0.00239 | <0.0463 | 0.282 |
| HSM120 | 3/25/2015 | 0.0382 | 0.000682 J | 0.00185 | 0.000351 J | <0.000479 | 0.0344 J | 0.00022 J |
| | 9/17/2015 | 0.0377 | 0.00629 | 0.0018 J | <0.000841 | <0.00239 | <0.0463 | 0.0017 J |
| HSM130 | 3/25/2015 | 0.0418 | 0.000662 J | 0.00194 | 0.000422 J | 0.00447 J | 0.0295 J | 0.00148 |
| | 9/17/2015 | 0.0414 | 0.00204 J | 0.00162 J | <0.000841 | <0.00239 | <0.0463 | 0.00154 J |
| HSM140 | 3/25/2015 | 0.0392 | 0.000951 J | 0.00208 | 0.000314 J | 0.000711 J | 0.0338 J | 0.000545 J |
| | 9/17/2015 | 0.0371 | 0.00218 J | 0.0016 J | <0.000841 | <0.00239 | <0.0463 | 0.00148 J |
| HSM150 | 3/25/2015 | 0.0388 | 0.000728 J | 0.00189 | 0.000282 J | 0.00135 J | 0.0288 J | 0.000616 J |
| | 9/17/2015 | 0.0369 | 0.00165 J | 0.00157 J | <0.000841 | <0.00239 | <0.0463 | 0.00182 J |
| HSM160 | 3/25/2015 | 0.0379 | 0.000978 J | 0.00207 | 0.000314 J | 0.0052 | 0.0342 J | 0.00128 |
| | 9/17/2015 | 0.0372 | 0.00691 | 0.00164 J | <0.000841 | 0.0559 | <0.0463 | 0.0032 J |
| HSM170 | 3/25/2015 | 0.0398 | 0.000824 J | 0.0018 | 0.000334 J | <0.000479 | 0.0348 J | 0.00167 |
| | 9/17/2015 | 0.0368 | 0.00856 | 0.00171 J | <0.000841 | <0.00239 | <0.0463 | 0.00325 J |
| MCL | | 2 | NE | NE | 0.05 | NE | NE | NE |
| PCL | | -- | 1.3 | 0.49 | -- | 7.3 | NE | 3.4 |

-- -- Not applicable

J – Detection is above the method detection limit, but below the reporting limit.

mg/L – milligrams per liter

MCL – maximum contaminant level

NE – None Established

PCL – protective concentration levels

7.2.1.8 Surface Water / Base Flow - Nitrates

Surface water samples were analyzed for nitrate-nitrite as nitrogen. Laboratory analyses indicated a limited range of nitrate-nitrite as nitrogen in surface water samples. Of the 16 surface water samples (14 environmental samples and two field duplicates) collected for the two sample events, concentrations ranged from 0.36 mg/L to 2.0 mg/L. None of the nitrate concentrations exceeded the MCL of 10 mg/L for drinking water. The highest nitrate concentration in surface water at the San Marcos Springs complex was 2.0 mg/L from HSM130 sampled on March 25, 2015. Nitrate-nitrogen results are summarized in Table 14.

Table 14. Surface Water Samples – Nitrate Detections - San Marcos Springs Complex

| Location | Date | Concentration (mg/L) |
|----------|-----------|----------------------|
| HSM110 | 3/25/2015 | 0.36 |
| | 9/17/2015 | 0.56 |
| FDHSM110 | 3/25/2015 | 0.36 |
| | 9/17/2015 | 0.59 |
| HSM120 | 3/25/2015 | 1.2 |
| | 9/17/2015 | 1.3 |
| HSM130 | 3/25/2015 | 2 |
| | 9/17/2015 | 1.7 |
| HSM140 | 3/25/2015 | 1.1 |
| | 9/17/2015 | 1.3 |
| HSM150 | 3/25/2015 | 1.1 |
| | 9/17/2015 | 1.3 |
| HSM160 | 3/25/2015 | 1.1 |
| | 9/17/2015 | 1.3 |
| HSM170 | 3/25/2015 | 1.1 |
| | 9/17/2015 | 1.3 |

mg/L – milligrams per liter

7.2.1.9 Surface water / base flow – Caffeine

Surface water base flows were analyzed for caffeine, which can indicate the possible presence of human wastewater discharge. Caffeine may enter surface water from leaking sewer or septic systems or it may be present in the aquifer from similar sources in the recharge zone (EPA, 2012). Potential ecological effects are currently unknown but could include reduced reproductive success in aquatic species (EPA, 2012). Caffeine detections in surface water samples from San Marcos Springs in 2015 ranged from 2.3 ng/L to 19 ng/L. There is no regulatory standard or expected value for comparison. Results are shown in Table 15.

**Table 15. Surface Water Samples – Caffeine
Detections - San Marcos Springs Complex**

| Location | Date Collected | Caffeine (ng/L) |
|----------|-------------------|--------------------|
| HSM110 | 3/25/2015 | 16 |
| | 9/17/2015 | 19 |
| FDHSM110 | 3/25/2015 | 10 |
| | 9/17/2015 | 13 |
| HSM120 | 3/25/2015 | <3.6 |
| | 9/17/2015 | 2.3 |
| HSM150 | 3/25/2015 | 14 |
| | 9/17/2015 | 3.8 |
| HSM160 | 3/25/2015 | <3.6 |
| | 9/17/2015 | 5.1 |

ng/L – nanograms per liter

7.2.2 San Marcos Springs Sediment Sampling

7.2.2.1 Sediment - Volatile Organic Compounds (VOCs)

VOCs were detected in sediment samples collected at three of the seven sample sites in the San Marcos Springs complex in 2015. All detections were below the laboratory reporting limits, TEC, and PEC values. The detections are summarized below in Table 16.

**Table 16. Sediment Samples – Volatile Organic Compound Detections - San Marcos Springs
Complex**

| Location | Date Collected | Acetone (mg/kg) | Carbon disulfide (mg/kg) | p-Isopropyltoluene (mg/kg) | Toluene (mg/kg) |
|----------|-------------------|--------------------|-----------------------------|-------------------------------|--------------------|
| HSM310 | 6/5/2015 | <0.022 | 0.0022J | <0.0023 | <0.0018 |
| HSM340 | 6/5/2015 | 0.011J | <0.00038 | <0.00078 | <0.00064 |
| HSM370 | 6/5/2015 | 0.014J | <0.00053 | <0.0011 | <0.00089 |
| FDHSM370 | 6/5/2015 | <0.012 | <0.00053 | 0.0014J | 0.0022J |
| TEC | | 60.030 | 0.120 | NE | 2.880 |
| PEC | | 360.180 | 0.780 | NE | 17.290 |

J – Detection is above the method detection limit, but below the reporting limit.

Mg/kg – milligrams per kilogram

NE – Not established

PEC – probable effect concentration

TEC – threshold effect concentration

7.2.2.2 Sediment - Semi-volatile Organic Compounds (SVOCs)

Generally, sediments would be expected to contain some SVOCs if the sediments have been exposed to these compounds. In summarizing the sediment SVOC detections, any compounds suspected as false positives are listed here in the text. Compounds of interest (primarily polycyclic aromatic hydrocarbons) are discussed in detail.

Non-Polycyclic Aromatic Hydrocarbon (PAH) Detections

Based on analysis of 2013 laboratory data, the EAA concluded that three compounds may have been laboratory artifacts. The compounds were bis(2-ethylhexyl) phthalate (DEHP), di-n-octyl phthalate, and di-n-butyl phthalate. The EAA noted in the *2013 Edwards Aquifer Habitat Conservation Plan Expanded Water Quality Report* (EAA, 2013) that as the data set grows, additional conclusions could be drawn. The 2014 laboratory analyses of sediment samples did not detect di-n-octyl phthalate or di-n-butyl phthalate. However, DEHP was detected in three of the sediment samples (HSM320, HSM330, and HSM350) in 2014 leading SWCA to conclude it is possible DEHP is present within the sediment and not just a laboratory artifact. DEHP was detected in three samples in 2015 as well, HSM330, HSM340 and HSM350. The 2015 detections were all below the laboratory reporting limit. The concentrations of DEHP in 2015 ranged from 0.12 J mg/kg to 0.74 J mg/kg. TEC and PEC values have not been established for DEHP. The detection of DEHP over the past two sampling events in samples only from the middle reaches of the San Marcos Spring system suggest DEHP is present within the sediment in that portion of the river.

One other non-PAH compound, 2-methylnaphthalene, was detected at a concentration of 0.15 J mg/kg at HSM370 in 2015. No TECs or PECs have been established for any of the non-PAH SVOCs detected.

PAH Detections

The remaining SVOC detections are all PAH compounds and are listed in Table 17. PAH detections are further shown in Figure 15, where the total PAH concentrations (sum of all detected concentrations for each sample point) are compared to the TEC and PEC values for total PAH concentration established by MacDonald et al. (2000) and TCEQ (2014a). Sample locations HSM320, HSM330, and HSM350 exceed the TEC for total PAH concentrations. HSM340 exceeded the PEC for total PAH concentrations.

Table 17. Sediment Samples – Semi-volatile Organic Compound Detections - San Marcos Springs Complex

| | | Acenaphthene | Anthracene | Benzo (a) Anthracene | Benzo (a) Pyrene | Benzo (b) Fluoranthene | Benzo (g,h,i) Perylene | Benzo (k) Fluoranthene | Chrysene | Dibenz (a,h) anthracene | Fluoranthene | Fluorene | Indeno(1,2,3-c,d) Pyrene | Phenanthrene | Pyrene | Total PAH | Bis(2-Ethylhexyl) Phthalate | Di-n-octyl phthalate | 2-Methylnaphthalene |
|----------|----------------|--------------|------------|----------------------|------------------|------------------------|------------------------|------------------------|----------|-------------------------|--------------|----------|--------------------------|--------------|---------|-----------|-----------------------------|----------------------|---------------------|
| Location | Date Collected | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) |
| HSM320 | 6/5/2015 | <0.32 | <0.32 | 0.51 J | 0.4 J | <0.33 | 0.33 J | <0.33 | 0.59 J | <0.24 | 0.84 J | <0.32 | 0.3 J | <0.35 | 0.79 J | 3.76 | <0.27 | <0.51 | <0.31 |
| HSM330 | 6/5/2015 | <0.079 | <0.079 | 0.65 | 0.57 J | 0.72 | 0.49 J | 0.61 J | 0.88 | 0.15 J | 1.5 | <0.079 | 0.43 J | 0.59 J | 1.4 | 7.99 | 0.12 J | <0.13 | <0.075 |
| HSM340 | 6/5/2015 | 0.27 J | 0.68 J | 4.7 | 3.8 | 4.8 | 2.9 | 3.8 | 6.2 | 0.88 J | 13 | 0.31 J | 2.7 | 6.6 | 12 | 62.64 | 0.74 J | <0.25 | <0.15 |
| HSM350 | 6/5/2015 | <0.14 | <0.14 | 0.15 J | 0.15 J | 0.21 J | 0.13 J | 0.15 J | 0.21 J | <0.1 | 0.3 J | <0.14 | 0.12 J | <0.15 | 0.66 J | 2.08 | 0.24 J | <0.22 | <0.13 |
| HSM360 | 6/5/2015 | <0.084 | <0.084 | <0.076 | <0.072 | 0.09 J | <0.073 | <0.086 | 0.087 J | <0.062 | 0.11 J | <0.084 | <0.071 | <0.091 | 0.11 J | 0.397 | <0.071 | <0.13 | <0.08 |
| HSM370 | 6/5/2015 | <0.11 | <0.11 | <0.097 | <0.093 | <0.11 | <0.094 | <0.11 | <0.11 | <0.079 | 0.11 J | <0.11 | <0.091 | <0.12 | <0.13 | 0.11 | <0.091 | 0.48 J | 0.15 J |
| TEC | | 0.0067 | 0.0572 | 0.108 | 0.150 | NE | NE | NE | 0.166 | 0.033 | 0.423 | 0.0774 | NE | 0.204 | 0.195 | 1.61 | NE | NE | NE |
| PEC | | 0.089 | 0.845 | 1.05 | 1.45 | NE | NE | NE | 1.290 | 0.140 | 2.23 | 0.536 | NE | 1.17 | 1.52 | 22.8 | NE | NE | NE |

J – Detection is above the method detection limit, but below the reporting limit.

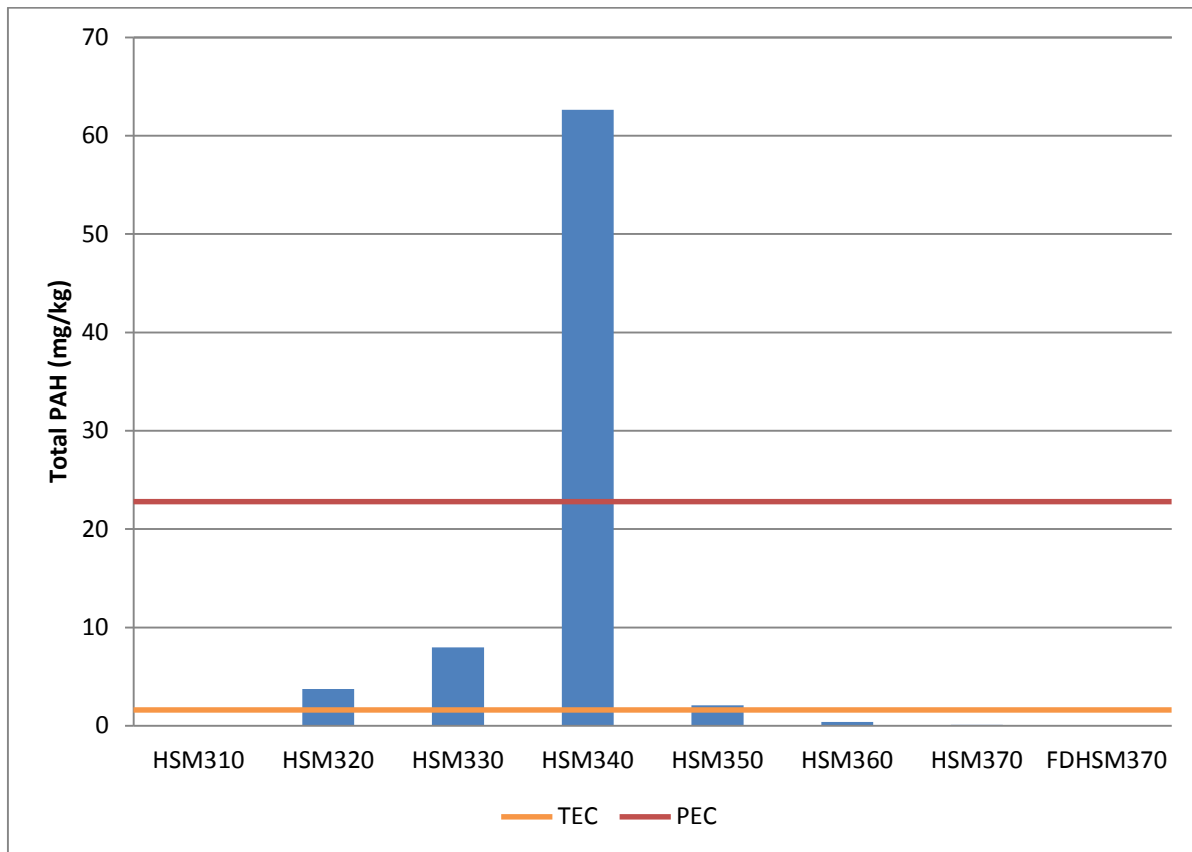
Mg/kg – milligrams per kilogram

NE – Not established

PEC – probable effect concentration

TEC – threshold effect concentration

Figure 15. San Marcos Springs Sediment Polycyclic Aromatic Hydrocarbons Detections Compared to Threshold Effect Concentration (TEC) and Probable Effect Concentration (PEC) values



7.2.2.3 Sediment - Pesticides

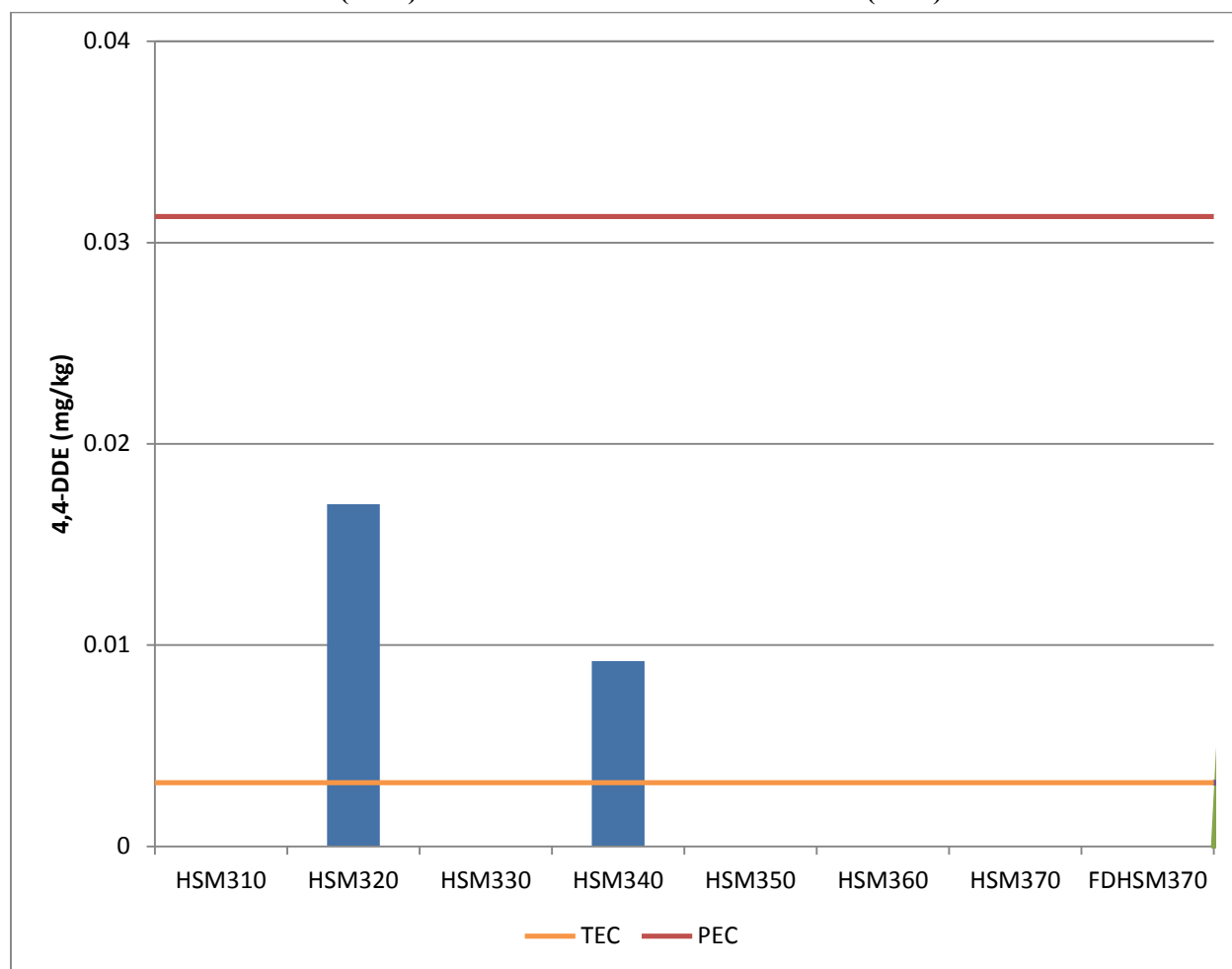
Sediment samples were analyzed for both organochlorine and organophosphorous pesticides. One pesticide, 4,4-DDE, was detected at HSM320 and HSM340. Both detections exceeded the established TEC value. The same compound was detected at HSM320 in 2013 and at HSM340 in 2014. Concentrations are listed in Table 18 and shown in Figure 16.

**Table 18. Sediment Samples – Pesticide Detections -
San Marcos Springs Complex**

| Location | Date Collected | 4,4-DDE ($\mu\text{g/kg}$) |
|----------|----------------|---------------------------------|
| HSM320 | 6/5/2015 | 17J |
| HSM340 | 6/5/2015 | 9.2 |
| TEC | | 3.16 |
| PEC | | 31.3 |

$\mu\text{g/kg}$ – micrograms per kilogram

Figure 16. San Marcos Springs Sediment 4,4-DDE Detections Compared to Threshold Effect Concentration (TEC) and Probable Effect Concentration (PEC) values



7.2.2.4 Sediment - Herbicides

Sediments were analyzed for herbicide compounds to further assess sediment quality at the San Marcos Springs complex. No herbicides were detected in any of the sediment samples from the seven sites in the San Marcos Springs complex.

7.2.2.5 Sediment - Polychlorinated Biphenyls

Sediments were analyzed for PCB compounds to further assess sediment quality at the San Marcos Springs complex. Two PCB compounds were detected in the sediment sample collected from HSM340. Aroclor-1260 was detected above the established TEC value, and the total PCBs detected were above the TEC for total PCB concentrations. The detections are listed in Table 19 and shown in comparison to the TEC and PEC values for total PCB concentrations in Figure 17.

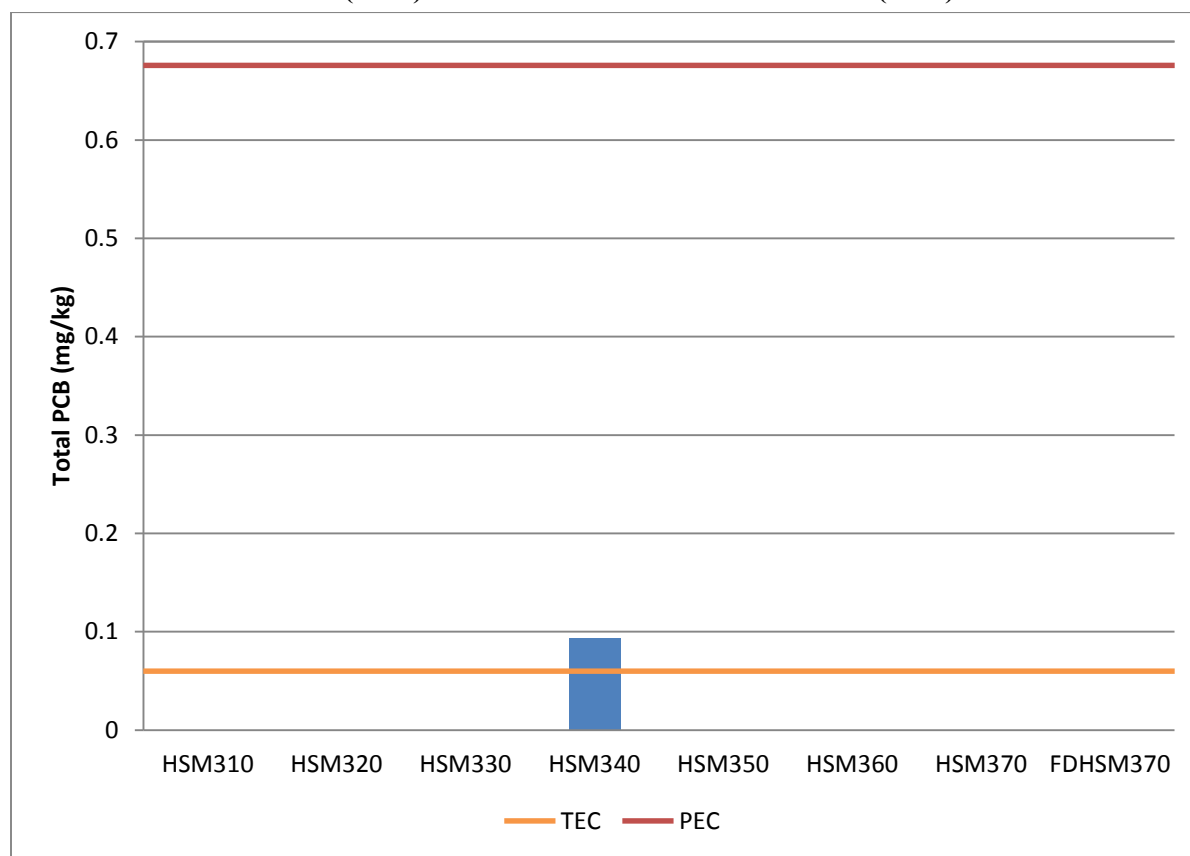
Table 19. Sediment Samples – Polychlorinated Biphenyls Detections - San Marcos Springs Complex

| Location | Date Collected | Aroclor-1254 (mg/kg) | Aroclor-1260 (mg/kg) | Total PCBs (mg/kg) |
|----------|----------------|-------------------------|-------------------------|-----------------------|
| HSM340 | 6/5/2015 | 0.041J | 0.052J | 0.093 |
| TEC | | 0.060 | 0.005 | 0.0598 |
| PEC | | 0.340 | 0.240 | 0.676 |

J – Detection is above the method detection limit, but below the reporting limit.

Mg/kg – milligrams per kilogram

Figure 17. San Marcos Springs Sediment 4,4-DDE Detections Compared to Threshold Effect Concentration (TEC) and Probable Effect Concentration (PEC) values



7.2.2.6 Sediment - Metals

Many metals are naturally occurring within soil, rock, and sediment. Sediment sample results for metals at the San Marcos Springs complex tested positive for several metals, generally at low concentrations. Metals detected above the method detection limit and subsequently evaluated in this report for potential toxic effects using the TEC and PEC standards are: antimony, arsenic, cadmium, chromium, copper, iron, lead, manganese, mercury, nickel, silver, and zinc. Other metals detected that do not have a TEC or PEC value available are aluminum, barium, beryllium, and selenium which were compared to TSBC (TCEQb, 2014). All selenium detections were below the laboratory reporting limit but above the TSBC level of 0.3 mg/kg. Sediment studies of selenium concentrations have shown that levels below four mg/kg are not likely to bioaccumulate in the food chain or have adverse impacts on the reproduction of fish or aquatic birds (Lemly, 1995; Moore et al., 1990; Van Derveer and Canton, 1996). Selenium detections did exceed this amount at HSM320 with a detected concentration of 4.20 J mg/kg.

Metal detections are listed in Table 20. Metals with detections above an established TEC value are displayed graphically in Figures 18–24, for arsenic, cadmium, copper, iron, lead, manganese, and zinc, respectively. Selenium detections are shown in Figure 25 in comparison to the TSBC and suggested toxic level of four mg/kg.

Table 20. Sediment Samples – Metal Detections - San Marcos Springs Complex

| Location | Date Collected | Antimony (mg/kg) | Arsenic (mg/kg) | Barium (mg/kg) | Beryllium (mg/kg) | Cadmium (mg/kg) | Chromium (mg/kg) | Copper (mg/kg) | Lead (mg/kg) | Nickel (mg/kg) | Selenium (mg/kg) | Silver (mg/kg) | Zinc (mg/kg) | Aluminum (mg/kg) | Iron (mg/kg) | Manganese (mg/kg) |
|-----------------------|----------------|---------------------|--------------------|-------------------|----------------------|--------------------|---------------------|-------------------|-----------------|-------------------|---------------------|-------------------|-----------------|---------------------|-----------------|----------------------|
| HSM310 | 6/5/2015 | <0.454 | 7.22 | 70.1 | 0.823 J | 0.64 J | 23.4 | 13.5 | 21.6 | 15.8 | 2.15 J | 0.506 J | 95.1 | 9790 | 16300 | 401 |
| HSM320 | 6/5/2015 | <0.665 | 7.16 | 87.3 | 1.16 J | 1.17 J | 26.6 | 58.3 | 65.4 | 20.5 | 4.2 J | 0.822 J | 253 | 12400 | 24500 | 390 |
| HSM330 | 6/5/2015 | <0.164 | 5.99 | 38.4 | <0.258 | 0.292 J | 8.38 | 3.84 | 16 | 10.8 | <0.332 | <0.179 | 60.7 | 1200 | 9350 | 334 |
| HSM340 | 6/5/2015 | 0.225 J | 2.98 | 31 | <0.256 | 0.441 J | 17 | 13.6 | 63.5 | 9.6 | <0.329 | 0.221 J | 61.2 | 2840 | 7640 | 152 |
| HSM350 | 6/5/2015 | <0.284 | 5.69 | 45.4 | 0.469 J | 0.596 J | 15.3 | 17.8 | 35.5 | 14.5 | 1.18 J | 0.384 J | 84.6 | 4890 | 12100 | 343 |
| HSM360 | 6/5/2015 | <0.175 | 4.11 | 34.7 | 0.343 J | 0.427 J | 10 | 5.74 | 54.4 | 9.41 | 0.386 J | 0.259 J | 33 | 4930 | 9170 | 240 |
| HSM370 | 6/5/2015 | <0.223 | 11 | 42.7 | 0.436 J | 0.612 J | 11.6 | 8.3 | 17.9 | 18 | <0.451 | 0.282 J | 42.6 | 5760 | 15300 | 439 |
| FDHSM370 | 6/5/2015 | <0.24 | 8.3 | 55.7 | 0.54 J | 0.56 J | 14 | 8.41 | 19.6 | 14.9 | 0.67 J | <0.262 | 43.5 | 8180 | 13800 | 471 |
| TEC | | 2 | 9.79 | NE | NE | 0.99 | 43.4 | 31.6 | 35.8 | 22.7 | NE | 1 | 121 | NE | 20000 | 460 |
| PEC | | 25 | 33.0 | NE | NE | 4.98 | 111 | 149 | 128 | 48.6 | NE | 2.2 | 459 | NE | 40000 | 1100 |
| Soil background level | | 1.0 | 5.9 | 300 | 1.5 | NE | 30 | 15 | 15 | 10 | 0.3 | NE | 30 | 30000 | 15000 | 300 |

J – Detection is above the method detection limit, but below the reporting limit.
Mg/kg – milligrams per kilogram
NE – Not established
PEC – probable effect concentration
TEC – threshold effect concentration

Figure 18. San Marcos Springs Sediment Arsenic Detections Compared to Threshold Effect Concentration (TEC) and Probable Effect Concentration (PEC) Values

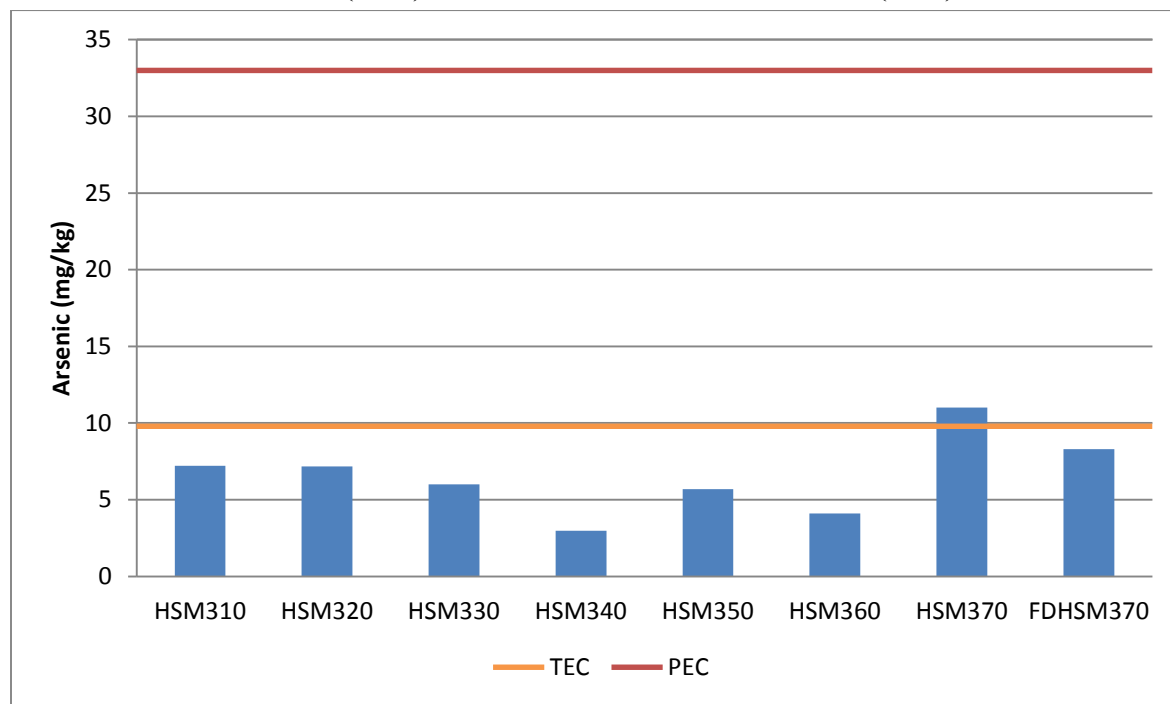


Figure 19. San Marcos Springs Sediment Cadmium Detections Compared to Threshold Effect Concentration (TEC) and Probable Effect Concentration (PEC) Values

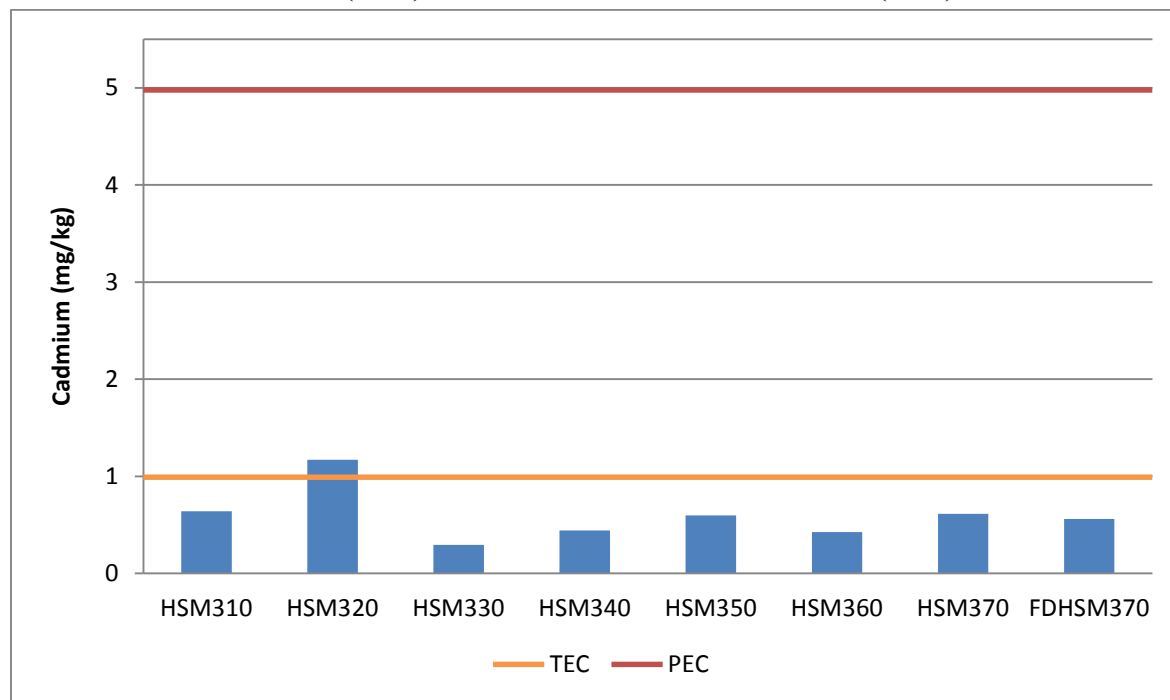


Figure 20. San Marcos Springs Sediment Copper Detections Compared to Threshold Effect Concentration (TEC) and Probable Effect Concentration (PEC) Values

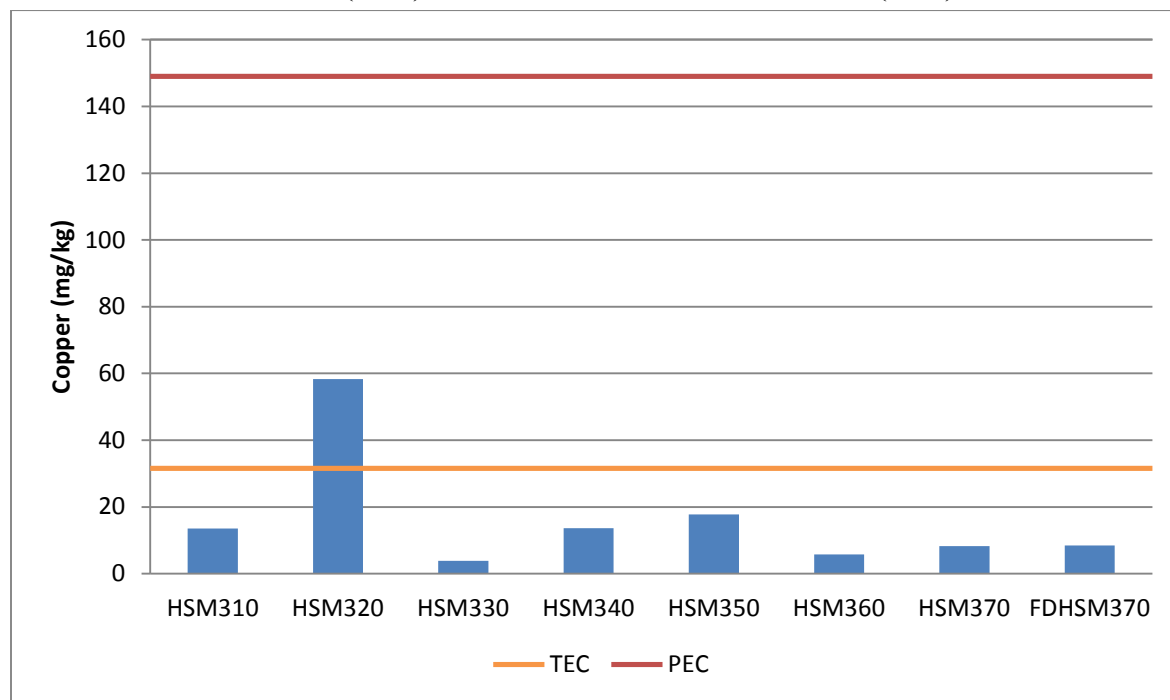


Figure 21. San Marcos Springs Sediment Iron Detections Compared to Threshold Effect Concentration (TEC) and Probable Effect Concentration (PEC) Values

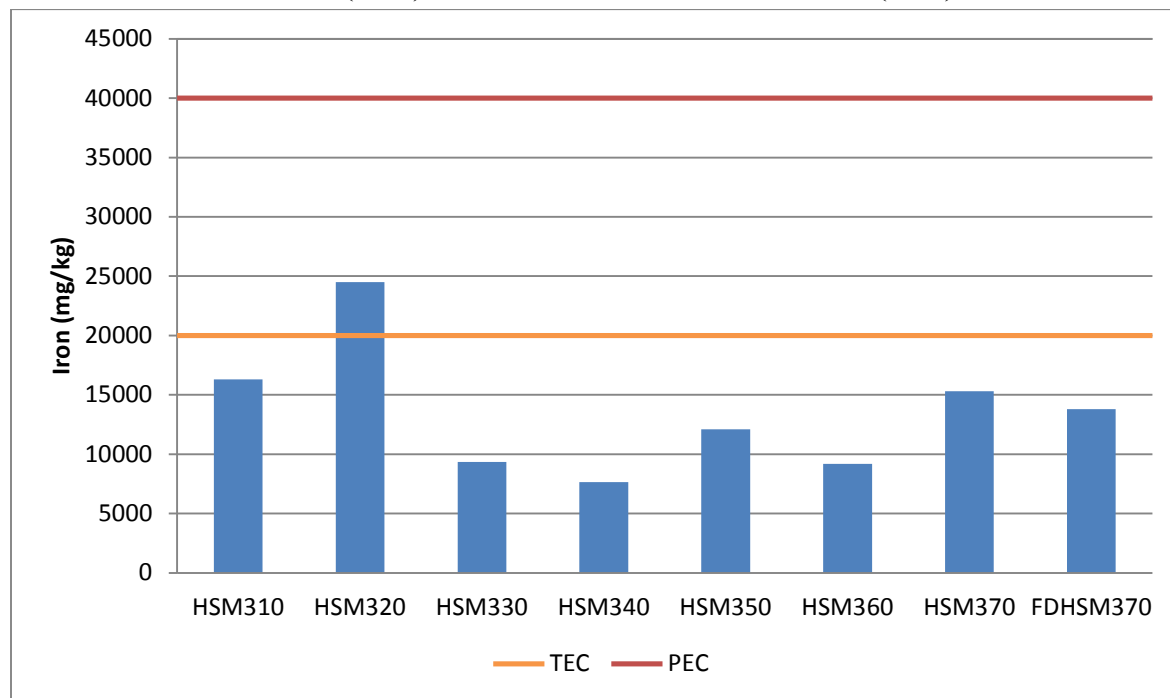


Figure 22. San Marcos Springs Sediment Lead Detections Compared to Threshold Effect Concentration (TEC) and Probable Effect Concentration (PEC) Values

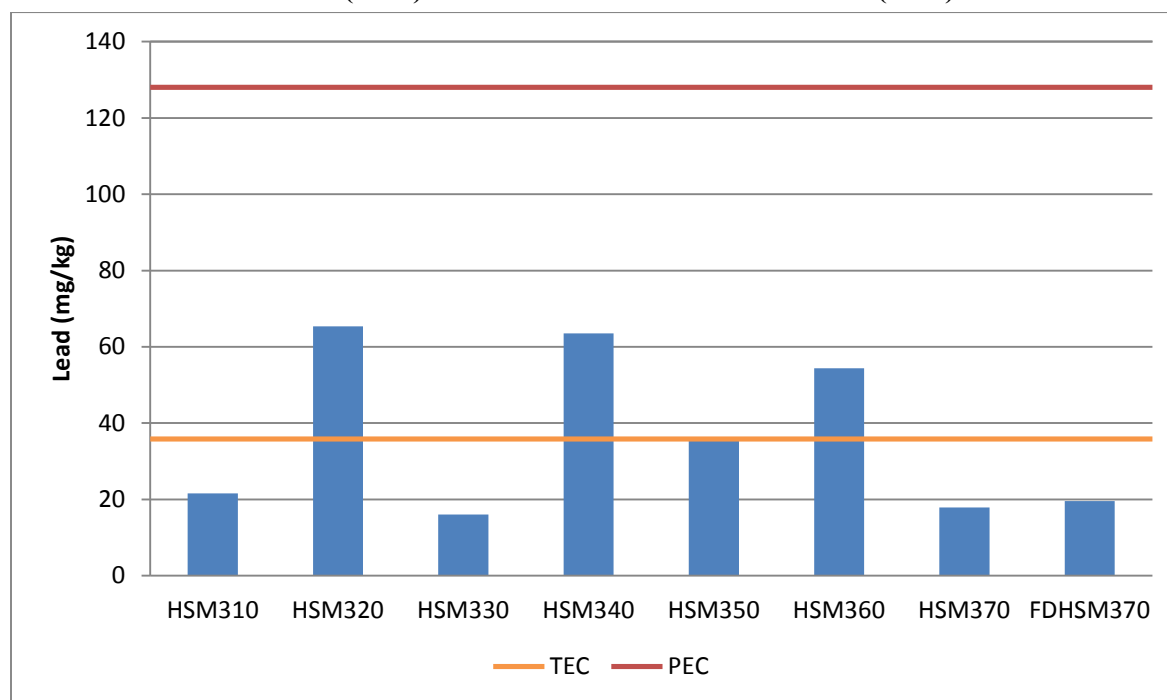


Figure 23. San Marcos Springs Sediment Manganese Detections Compared to Threshold Effect Concentration (TEC) and Probable Effect Concentration (PEC) Values

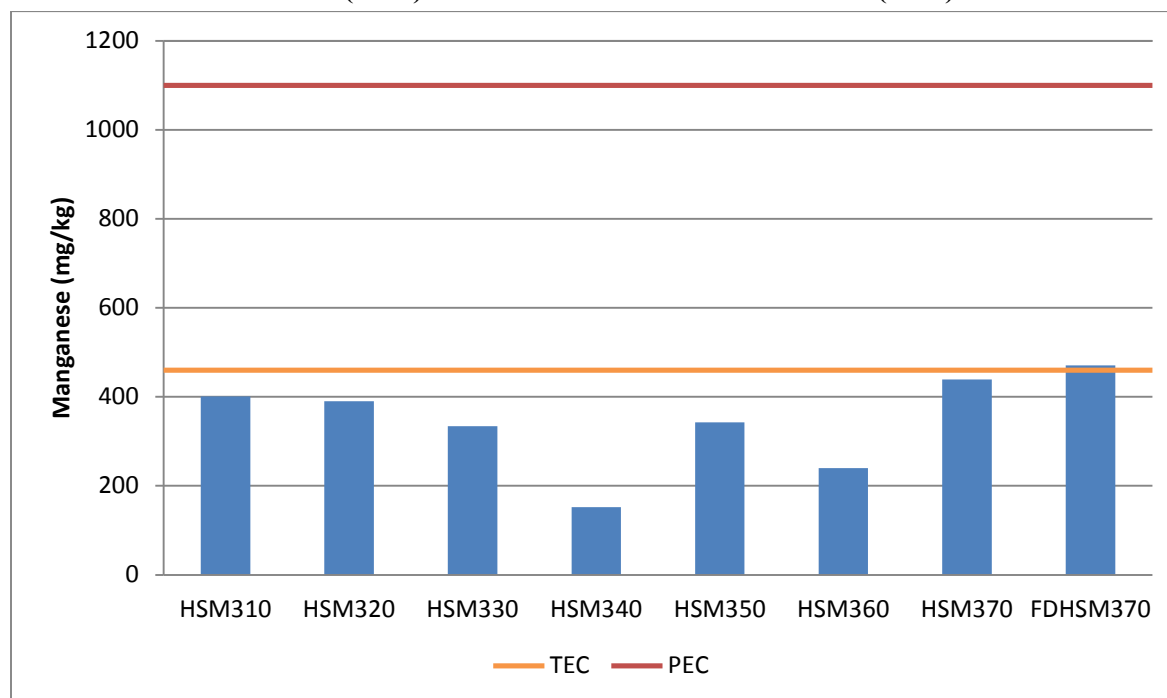


Figure 24. San Marcos Springs Sediment Zinc Detections Compared to Threshold Effect Concentration (TEC) and Probable Effect Concentration (PEC) Values

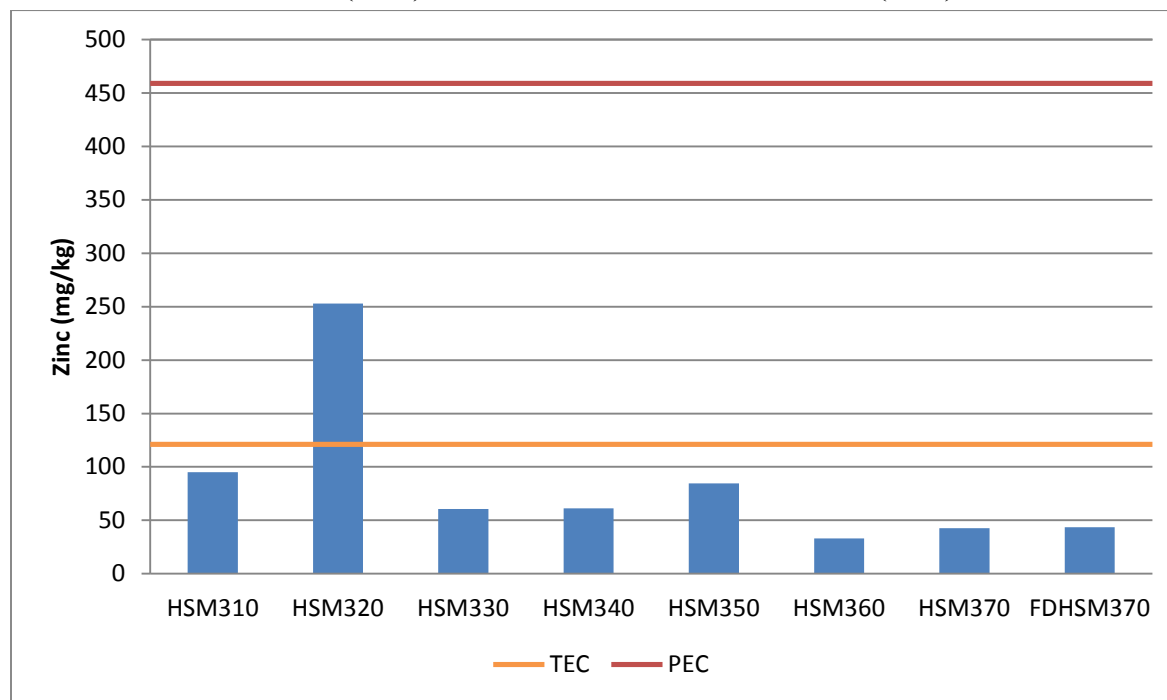
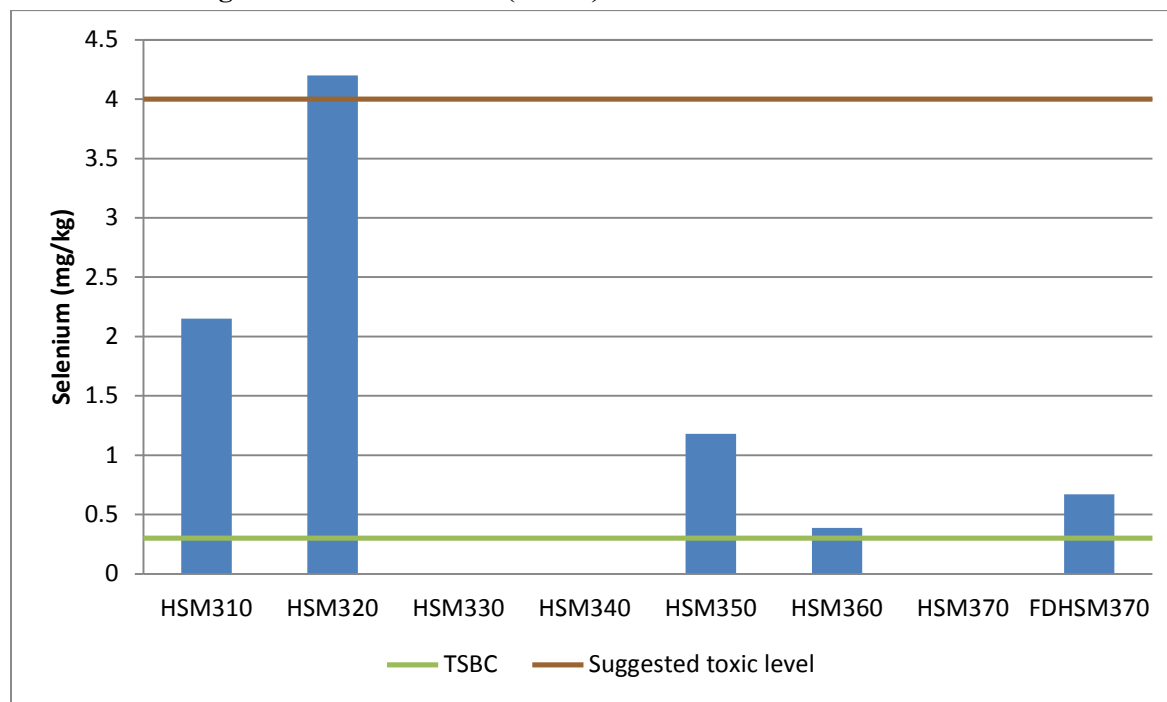


Figure 25. San Marcos Springs Sediment Selenium Detections Compared to Texas-Specific Background Concentration (TSBC) and Possible Bioaccumulation Toxic Values



7.2.3 San Marcos Springs Stormwater Sampling

Stormwater samples were collected during two storm events at the San Marcos Springs complex. Each event was sampled according to the guidelines in the EAHCP workplan. Event 1 occurred on May 5–6, 2015. Total rainfall for the first event was approximately 0.25 to 0.50 inches (NOAA, 2015) causing streamflow measured at USGS Gauge 08170500 to increase from approximately 180 cfs to a peak of 205 cfs during the event (USGS, 2015). The second event occurred on October 23, 2015, with approximately 2.50 to 2.99 inches of precipitation falling during the 24-hour period beginning at 7:00 AM on October 23, 2015 (NOAA, 2015). Sampling efforts occurred during the first pulse of the storm when precipitation was estimated to be approximately 1 inch.

7.2.3.1 Stormwater - Bacteria Detections

Stormwater samples collected and analyzed for bacteria analyses generally tested positive for high levels of bacteria. The 2014 Texas Surface Water Quality Standard for *E. coli* in primary recreation waters is a geometric mean of 126 MPN/100 mL with no individual sample exceeding 399 MPN/100 mL (30 TAC 307.7). The geometric mean for stormwater samples collected from the San Marcos Springs complex during May 2015 was 1264.1 MPN/100 mL. Bacteria counts ranged from 66 MPN/100 mL to 24,000 MPN/100 mL with several samples exceeding the individual sample limit during the May 2015 event. In October 2015, *E. coli* counts ranged from 41 MPN/100 mL to 87,000 MPN/100 mL with a geometric mean of 3693.4 MPN/100 mL. Individual detections are listed below in Table 21 and shown in relation to stream discharge and specific conductivity in Figures 26 and 27. Due to the timing of storm events and laboratory working hours, it was not possible to deliver all samples to the laboratory within sample holding times, as discussed in Appendix C. Samples exceeding hold times were included in the range and geometric mean calculations.

Table 21. Stormwater Samples – Bacteria Counts - San Marcos Springs Complex

| Location | Date | Concentration (MPN/100 mL) |
|----------------|------------|----------------------------|
| HSM210 Lead | 5/5/2015 | 77 H |
| | 10/23/2015 | 41 H |
| HSM210 Peak | 5/5/2015 | 66 H |
| | 10/23/2015 | 41 H |
| HSM210 Trail | 5/6/2015 | 270 H |
| | 10/23/2015 | 450 H |
| FDHSM210 Trail | 5/6/2015 | 330 H |
| | 10/23/2015 | 180 H |
| HSM230 Lead | 5/5/2015 | 20000 H |
| | 10/23/2015 | 87000 H |
| HSM230 Peak | 5/5/2015 | 14000 H |
| | 10/23/2015 | 11000 H |
| HSM230 Trail | 5/6/2015 | 5200 H |
| | 10/23/2015 | 73000 H |

Table 21. Stormwater Samples – Bacteria Counts - San Marcos Springs Complex

| Location | Date | Concentration (MPN/100 mL) |
|----------------|------------|----------------------------|
| FDHSM230 Trail | 5/6/2015 | 4600 H |
| | 10/23/2015 | 65000 H |
| HSM231 Lead | 5/5/2015 | 870 H |
| | 10/23/2015 | 11000 H |
| HSM231 Peak | 5/5/2015 | 770 H |
| | 10/23/2015 | 2200 H |
| HSM231 Trail | 5/6/2015 | 280 H |
| | 10/23/2015 | 1700 H |
| FDHSM231 Trail | 5/6/2015 | 330 H |
| | 10/23/2015 | 1700 H |
| HSM240 Lead | 5/5/2015 | 24000 H |
| | 10/23/2015 | 28000 H |
| HSM240 Peak | 5/6/2015 | 290 H |
| | 10/23/2015 | 1000 H |
| HSM240 Trail | 5/6/2015 | 210 H |
| | 10/23/2015 | 750 H |
| HSM250 Lead | 5/5/2015 | 2400 H |
| | 10/23/2015 | 9200 H |
| HSM250 Peak | 5/5/2015 | 2200 H |
| | 10/23/2015 | 6500 H |
| HSM250 Trail | 5/6/2015 | 2500 H |
| | 10/23/2015 | 1900 H |
| HSM260 Lead | 5/5/2015 | 2400 H |
| | 10/23/2015 | 7300 H |
| HSM260 Peak | 5/5/2015 | 2100 H |
| | 10/23/2015 | 24000 H |
| HSM260 Trail | 5/6/2015 | 1100 H |
| | 10/23/2015 | 4600 H |
| HSM270 Lead | 5/5/2015 | 1400 H |
| | 10/23/2015 | 14000 H |
| HSM270 Peak | 5/6/2015 | 5800 H |
| | 10/23/2015 | 8700 H |
| HSM270 Trail | 5/6/2015 | 1700 H |
| | 10/23/2015 | 4100 H |

H – Analyzed outside hold time, result included for comparison but not considered valid

MPN/100 mL – Most probable number per 100 milliliters of water

Figure 26. Stormwater Samples – May 2015 Bacteria Counts in Relation to Stream Discharge and Specific Conductivity – San Marcos Springs Complex

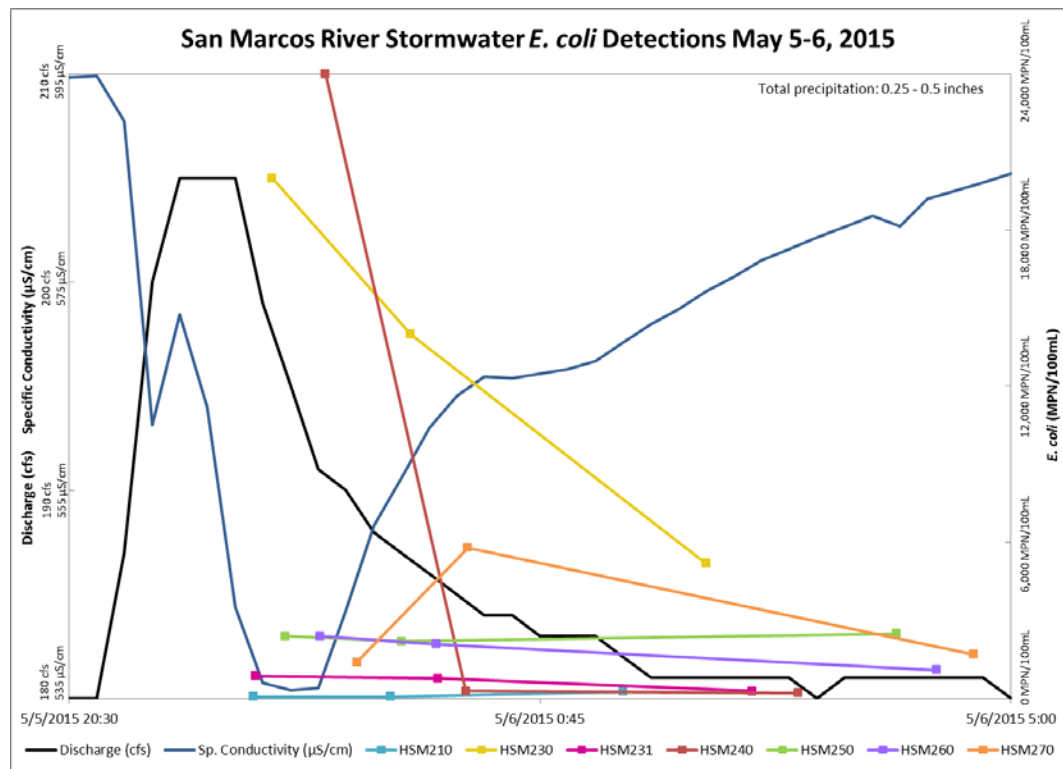
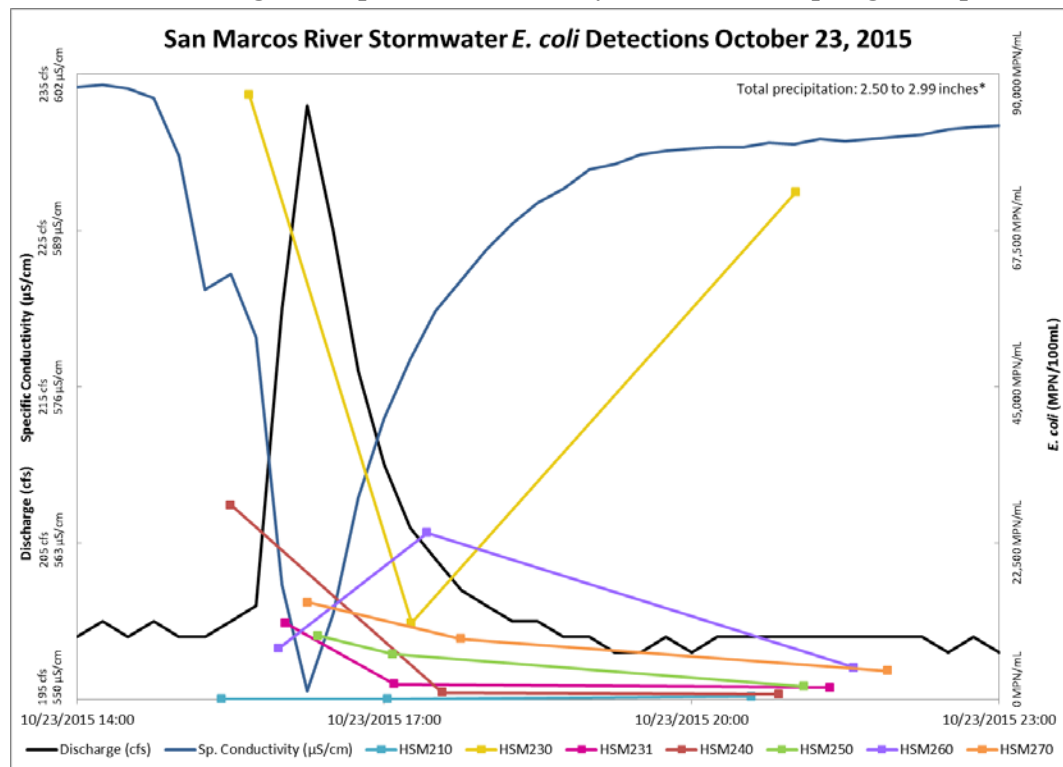


Figure 27. Stormwater Samples – October 2015 Bacteria Counts in Relation to Stream Discharge and Specific Conductivity – San Marcos Springs Complex



7.2.3.2 Stormwater - Volatile Organic Compounds (VOCs)

One VOC compound, p-Isopropyltoluene, was detected in HSM210 Trail during the May Sampling event. Trichloroethene was detected in two samples during the October sampling event. Toluene was also detected in one sample from the October event. All of these detections were below established MCLs or PCLs, and only one exceeded the laboratory reporting limit. The detections are summarized in Table 22.

Table 22. Stormwater Samples – Volatile Organic Compound Detections - San Marcos Springs Complex

| Location | Date Collected | p-Isopropyltoluene (µg/L) | Toluene (µg/L) | Trichloroethene (µg/L) |
|--------------|----------------|---------------------------|----------------|------------------------|
| HSM210 Trail | 5/6/2015 | 0.18 J | <0.24 | <0.37 |
| HSM230 Lead | 10/23/2015 | <0.16 | <0.24 | 1.1 |
| HSM240 Lead | 10/23/2015 | <0.16 | <0.24 | 0.38 J |
| HSM240 Peak | 10/23/2015 | <0.16 | 0.26 J | <0.37 |
| MCL | | NE | 1000 | NE |
| PCL | | 2400 | -- | NE |

J – Detection is above the method detection limit, but below the reporting limit

µg/L – milligrams per Liter

NE – Not established

MCL – maximum contaminant level

PCL – protective concentration levels

-- - not applicable

7.2.3.3 Stormwater - Semi-volatile Organic Compounds (SVOCs)

Stormwater samples were collected and analyzed for SVOCs. No SVOCs were detected in any of the samples collected from the San Marcos Springs complex during the May or October 2015 sampling events.

7.2.3.4 Stormwater - Herbicides and Pesticides

Stormwater samples were collected and analyzed for organochlorine pesticides, organophosphorous pesticides, and herbicides. There were very few detections of pesticides or herbicides in any of the stormwater samples collected from the San Marcos Springs complex during 2015. Endosulfan I was detected in the lead samples collected at HSM230 and HSM240 during the May 2015 event and in HSM230 Lead during the October 2015 event. 4,4-DDD was detected in the trail sample collected at HSM230 in May 2015. All of these detections were below the laboratory reporting limits and established PCLs for the compounds. These detections are shown in Table 23.

Table 23. Stormwater Samples – Pesticide Detections - San Marcos Springs Complex

| Location | Date Collected | 4,4-DDD (µg/L) | Endosulfan I (µg/L) |
|--------------|----------------|-------------------|------------------------|
| HSM230 Lead | 5/5/2015 | <0.027 | 0.048 J |
| | 10/23/2015 | <0.027 | 0.090 J |
| HSM230 Trail | 5/6/2015 | 0.038 J | <0.028 |
| HSM240 Lead | 5/5/2015 | <0.027 | 0.045 J |
| MCL | | NE | NE |
| PCL | | 3.8 | 49 |

J – Detection is above the method detection limit, but below the reporting limit

µg/L – milligrams per Liter

MCL – maximum contaminant level

NE – Not established

PCL – protective concentration levels

7.2.3.5 Stormwater - Polychlorinated Biphenyls

Stormwater samples were analyzed for the various Aroclor compounds that are generally referred to collectively as PCBs. None of the stormwater samples from the San Marcos Springs complex indicated positive detections of PCB compounds during the May 2015 sampling event.

7.2.3.6 Stormwater - Metals

Stormwater samples were analyzed for metals in accordance with the EAHCP workplan. Several positive metal detections were noted in the sample set; however, no samples contained a metal at a concentration in excess of the drinking water MCL, and most detections were below laboratory reporting limits during the May 2015 sampling event.

7.2.3.7 Stormwater - Nitrates

Stormwater samples were analyzed for nitrate-nitrite as nitrogen in accordance with the EAHCP workplan. All samples contained nitrate results below the MCL of 10 mg/L. During the May 2015 event, the range of nitrate results was 0.38 mg/L to 1.5 mg/L, with an average concentration of 0.99 mg/L. During the October 2015 event, nitrate concentrations ranged from 0.60 mg/L to 1.6 mg/L, with an average of 1.1 mg/L. For comparison, the average nitrate in spring water samples at San Marcos Springs for calendar year 2014 was 1.42 mg/L (EAA, 2015). Nitrate detections are summarized in Table 24.

**Table 24. Stormwater Samples – Nitrate Detections
- San Marcos Springs Complex**

| Location | Date | Concentration (mg/L) |
|----------------|------------|----------------------|
| HSM210 Lead | 5/5/2015 | 0.38 |
| | 10/23/2015 | 0.62 H |
| HSM210 Peak | 5/5/2015 | 0.38 |
| | 10/23/2015 | 0.60 H |
| HSM210 Trail | 5/6/2015 | 0.41 |
| | 10/23/2015 | 0.61 H |
| FDHSM210 Trail | 5/6/2015 | 0.41 |
| | 10/23/2015 | 0.62 H |
| HSM230 Lead | 5/5/2015 | 0.79 |
| | 10/23/2015 | 0.94 H |
| HSM230 Peak | 5/5/2015 | 1.4 |
| | 10/23/2015 | 1.6 H |
| HSM230 Trail | 5/6/2015 | 1.5 |
| | 10/23/2015 | 1.5 H |
| FDHSM230 Trail | 5/6/2015 | 1.5 |
| | 10/23/2015 | 1.6 H |
| HSM231 Lead | 5/5/2015 | 1 |
| | 10/23/2015 | 1.2 H |
| HSM231 Peak | 5/5/2015 | 1.1 |
| | 10/23/2015 | 1.2 H |
| HSM231 Trail | 5/6/2015 | 1.2 |
| | 10/23/2015 | 1.3 H |
| FDHSM231 Trail | 5/6/2015 | 1.2 |
| | 10/23/2015 | 1.2 H |
| HSM240 Lead | 5/5/2015 | 0.68 |
| | 10/23/2015 | 1.1 H |
| HSM240 Peak | 5/6/2015 | 1.1 |
| | 10/23/2015 | 1.2 H |
| HSM240 Trail | 5/6/2015 | 1.2 |
| | 10/23/2015 | 1.2 H |
| HSM250 Lead | 5/5/2015 | 1 |
| | 10/23/2015 | 1.2 H |
| HSM250 Peak | 5/5/2015 | 1 |
| | 10/23/2015 | NA |
| HSM250 Trail | 5/6/2015 | 1.1 |
| | 10/23/2015 | 1.2 H |
| HSM260 Lead | 5/5/2015 | 1.1 |
| | 10/23/2015 | 1.2 H |
| HSM260 Peak | 5/5/2015 | 1.1 |
| | 10/23/2015 | 1.2 H |
| HSM260 Trail | 5/6/2015 | 1.1 |
| | 10/23/2015 | 1.2 H |

**Table 24. Stormwater Samples – Nitrate Detections
- San Marcos Springs Complex**

| Location | Date | Concentration (mg/L) |
|--------------|------------|-------------------------|
| HSM270 Lead | 5/5/2015 | 1.1 |
| | 10/23/2015 | 1.1 H |
| HSM270 Peak | 5/6/2015 | 1 |
| | 10/23/2015 | 1.2 H |
| HSM270 Trail | 5/6/2015 | 1.1 |
| | 10/23/2015 | 1.2 H |

H – Analyzed outside hold time, result included for comparison but not considered valid

mg/L – milligrams per liter

NA – Not analyzed

7.2.3.8 Stormwater – Caffeine

Stormwater was analyzed for caffeine, which can indicate an anthropogenic source. Caffeine may enter surface water from leaking sewer or septic systems or it may be present in the aquifer from similar sources in the recharge zone (EPA, 2012). Potential ecological effects are currently unknown but could include reduced reproductive success in aquatic species (EPA, 2012). Caffeine detections in stormwater samples from San Marcos Springs in May 2015 ranged from 10 ng/L to 4800 ng/L and were detected in every sample. Caffeine detections in stormwater samples from San Marcos Springs in October 2015 ranged from 2.4 ng/L to 2200 ng/L and were detected in every sample. There is no regulatory standard or expected value for comparison. Results are shown in Table 25.

**Table 25. Stormwater Samples – Caffeine
Detections - San Marcos Springs Complex**

| Location | Date Collected | Caffeine (ng/L) |
|----------------|-------------------|--------------------|
| HCS210 Lead | 5/5/2015 | 10 |
| | 10/23/2015 | 5.7 |
| HCS210 Peak | 5/5/2015 | 17 |
| | 10/23/2015 | 5.8 |
| HCS210 Trail | 5/6/2015 | 13 |
| | 10/23/2015 | 2.4 |
| FDHSM210 Trail | 5/6/2015 | 14 |
| | 10/23/2015 | 5.7 |
| HSM230 Lead | 5/5/2015 | 4800 |
| | 10/23/2015 | 2200 |
| HSM230 Peak | 5/5/2015 | 830 |
| | 10/23/2015 | 260 |

**Table 25. Stormwater Samples – Caffeine
Detections - San Marcos Springs Complex**

| Location | Date Collected | Caffeine (ng/L) |
|----------------|-------------------|--------------------|
| HSM230 Trail | 5/6/2015 | 290 |
| | 10/23/2015 | 470 |
| FDHSM230 Trail | 5/6/2015 | 340 |
| | 10/23/2015 | 450 |
| HSM231 Lead | 5/5/2015 | 370 |
| | 10/23/2015 | 210 |
| HSM231 Peak | 5/5/2015 | 83 |
| | 10/23/2015 | 61 |
| HSM231 Trail | 5/6/2015 | 25 |
| | 10/23/2015 | 19 |
| FDHSM231 Trail | 5/6/2015 | 25 |
| | 10/23/2015 | 15 |
| HCS240 Lead | 5/5/2015 | 3100 |
| | 10/23/2015 | 510 |
| HCS240 Peak | 5/6/2015 | 44 |
| | 10/23/2015 | 60 |
| HCS240 Trail | 5/6/2015 | 25 |
| | 10/23/2015 | 23 |
| HCS250 Lead | 5/5/2015 | 350 |
| | 10/23/2015 | 260 |
| HCS250 Peak | 5/5/2015 | 230 |
| | 10/23/2015 | NA |
| HCS250 Trail | 5/6/2015 | 130 |
| | 10/23/2015 | 24 |
| HCS260 Lead | 5/5/2015 | 230 |
| | 10/23/2015 | 88 |
| HCS260 Peak | 5/5/2015 | 870 |
| | 10/23/2015 | 330 |
| HCS260 Trail | 5/6/2015 | 110 |
| | 10/23/2015 | 42 |
| HCS270 Lead | 5/5/2015 | 180 |
| | 10/23/2015 | 130 |
| HCS270 Peak | 5/6/2015 | 450 |
| | 10/23/2015 | 200 |
| HCS270 Trail | 5/6/2015 | 190 |
| | 10/23/2015 | 56 |

NA – Not analyzed
ng/L – nanograms per liter

7.2.4 San Marcos Springs Surface Water Passive Sampling

PDSs were installed in the San Marcos Springs system in February, April, June, August, October and December 2015. Any changes to deployment locations are discussed in Appendix C. There are no regulatory standards for comparison to PDS results. PDSs were analyzed for VOCs, SVOCs, and organochlorine pesticides. Few compounds were detected, only tetrachloroethene and TPH were consistently detected. Positive detections are shown below in Table 26.

Table 26. Passive Diffusion – San Marcos Springs Complex

| Location | Month 2015 | Chloroform (µg) | Ethylbenzene (µg) | Naphthalene (µg) | p/m-Xylene (µg) | Tetrachloroethene (µg) | Toluene (µg) | BTEX (µg) | TPH (µg) |
|----------|------------|--------------------|----------------------|---------------------|--------------------|---------------------------|-----------------|--------------|-------------|
| HSM410 | February | <0.02 | <0.02 | <0.05 | <0.02 | <0.02 | <0.02 | <0.02 | 0.90 |
| | April | <0.02 | <0.02 | <0.05 | <0.02 | <0.02 | <0.02 | <0.02 | 0.93 |
| | June | <0.02 | <0.02 | <0.05 | <0.02 | <0.02 | <0.02 | <0.02 | 0.60 |
| | August | <0.02 | <0.02 | <0.05 | <0.02 | <0.02 | <0.02 | <0.02 | <0.50 |
| | October | <0.02 | <0.02 | <0.05 | <0.02 | <0.02 | <0.02 | <0.02 | <0.50 |
| | December | <0.02 | <0.02 | <0.05 | <0.02 | <0.02 | <0.02 | <0.02 | <0.50 |
| HSM420 | February | <0.02 | <0.02 | <0.05 | <0.02 | 0.08 | <0.02 | <0.02 | <0.50 |
| | April | <0.02 | 0.03 | <0.05 | 0.09 | 0.10 | 0.02 | 0.14 | 0.54 |
| | June | <0.02 | <0.02 | <0.05 | <0.02 | 0.14 | <0.02 | <0.02 | <0.50 |
| | August | <0.02 | <0.02 | <0.05 | <0.02 | 0.06 | <0.02 | <0.02 | 0.67 |
| | October | <0.02 | <0.02 | <0.05 | <0.02 | 0.08 | <0.02 | <0.02 | <0.50 |
| | December | <0.02 | <0.02 | <0.05 | <0.02 | 0.10 | <0.02 | <0.02 | <0.50 |
| HSM430 | February | <0.02 | <0.02 | <0.05 | <0.02 | 0.13 | <0.02 | <0.02 | 1.27 |
| | April | <0.02 | <0.02 | <0.05 | <0.02 | 0.80 | <0.02 | <0.02 | <0.50 |
| | June | <0.02 | <0.02 | <0.05 | <0.02 | 1.98 | <0.02 | <0.02 | 0.96 |
| | August | <0.02 | <0.02 | <0.05 | <0.02 | 0.68 | <0.02 | <0.02 | <0.50 |
| | October | <0.02 | <0.02 | <0.05 | <0.02 | 0.94 | <0.02 | <0.02 | <0.50 |
| | December | 0.02 | <0.02 | <0.05 | <0.02 | 0.81 | <0.02 | <0.02 | <0.50 |
| FDHSM430 | February | <0.02 | <0.02 | <0.05 | <0.02 | 0.16 | <0.02 | <0.02 | 0.99 |
| | April | <0.02 | <0.02 | <0.05 | <0.02 | 0.72 | <0.02 | <0.02 | 0.54 |
| | June | <0.02 | <0.02 | <0.05 | <0.02 | 2.35 | <0.02 | <0.02 | 0.54 |
| | August | <0.02 | <0.02 | <0.05 | <0.02 | 0.89 | <0.02 | <0.02 | <0.50 |
| | October | <0.02 | <0.02 | <0.05 | <0.02 | 1.01 | <0.02 | <0.02 | <0.50 |
| | December | 0.02 | <0.02 | <0.05 | <0.02 | 0.76 | <0.02 | <0.02 | <0.50 |

Table 26. Passive Diffusion – San Marcos Springs Complex

| Location | Month 2015 | Chloroform (µg) | Ethylbenzene (µg) | Naphthalene (µg) | p/m-Xylene (µg) | Tetrachloroethene (µg) | Toluene (µg) | BTEX (µg) | TPH (µg) |
|----------|------------|--------------------|----------------------|---------------------|--------------------|---------------------------|-----------------|--------------|-------------|
| HSM440 | February | <0.02 | <0.02 | <0.05 | <0.02 | 0.13 | <0.02 | <0.02 | <0.50 |
| | April | <0.02 | <0.02 | <0.05 | <0.02 | 0.13 | <0.02 | <0.02 | <0.50 |
| | June | <0.02 | <0.02 | <0.05 | <0.02 | 0.19 | <0.02 | <0.02 | 0.68 |
| | August | <0.02 | <0.02 | <0.05 | <0.02 | 0.11 | <0.02 | <0.02 | 1.18 |
| | October | <0.02 | <0.02 | <0.05 | <0.02 | 0.11 | <0.02 | <0.02 | <0.50 |
| | December | <0.02 | <0.02 | <0.05 | <0.02 | 0.12 | <0.02 | <0.02 | <0.50 |
| HSM450 | February | <0.02 | <0.02 | <0.05 | <0.02 | 0.05 | <0.02 | <0.02 | <0.50 |
| | April | <0.02 | <0.02 | <0.05 | <0.02 | 0.08 | <0.02 | <0.02 | <0.50 |
| | June | <0.02 | <0.02 | <0.05 | <0.02 | 0.09 | <0.02 | <0.02 | <0.50 |
| | August | NA | NA | NA | NA | NA | NA | NA | NA |
| | October | <0.02 | <0.02 | <0.05 | <0.02 | 0.04 | <0.02 | <0.02 | <0.50 |
| | December | <0.02 | <0.02 | <0.05 | <0.02 | 0.05 | <0.02 | <0.02 | <0.50 |
| HSM460 | February | <0.02 | <0.02 | <0.05 | <0.02 | 0.09 | <0.02 | <0.02 | <0.50 |
| | April | <0.02 | <0.02 | <0.05 | <0.02 | 0.11 | <0.02 | <0.02 | <0.50 |
| | June | <0.02 | <0.02 | <0.05 | <0.02 | 0.15 | <0.02 | <0.02 | 0.67 |
| | August | <0.02 | <0.02 | <0.05 | <0.02 | 0.05 | <0.02 | <0.02 | <0.50 |
| | October | <0.02 | <0.02 | <0.05 | <0.02 | 0.08 | <0.02 | <0.02 | <0.50 |
| | December | <0.02 | <0.02 | <0.05 | <0.02 | 0.10 | <0.02 | <0.02 | <0.50 |
| HSM470 | February | <0.02 | <0.02 | <0.05 | <0.02 | 0.10 | <0.02 | <0.02 | 0.54 |
| | April | <0.02 | <0.02 | <0.05 | <0.02 | 0.09 | <0.02 | <0.02 | 0.68 |
| | June | <0.02 | <0.02 | <0.05 | 0.03 | 0.18 | <0.02 | 0.03 | 0.60 |
| | August | <0.02 | <0.02 | <0.05 | <0.02 | 0.05 | <0.02 | <0.02 | <0.50 |
| | October | <0.02 | <0.02 | 0.05 | <0.02 | 0.09 | <0.02 | <0.02 | <0.50 |
| | December | <0.02 | <0.02 | <0.05 | <0.02 | 0.08 | <0.02 | <0.02 | <0.50 |

BTEX – benzene, toluene, ethylbenzene, and xylenes

NA – Not analyzed

µg – micrograms

TPH – total petroleum hydrocarbons

8.0 SUMMARY OF RESULTS

SWCA staff collected surface water (base flow), stormwater, sediment, and passive diffusion samples from Comal and San Marcos springs complexes. The sampling events met the requirements of the EAHCP and provided background data for these two systems. The limited number of detections above comparative standards is indicative of generally high water quality. However, the total PAH and selenium results that exceeded comparative standards were of concern. Continued sampling will be helpful in assessing if these detected compounds are repeated and have the potential for any long-term impact on listed species. Specific detections of interest, such as compounds detected above an MCL (for water) or PEC (for sediment), are listed below.

San Marcos Sediment:

| | | |
|--------|-----------|---|
| HSM340 | Total PAH | 62.64 mg/kg (PEC = 22.8) |
| HSM320 | Selenium | 4.20 J mg/kg (Bioaccumulation Toxicity = 4.0) |

PAHs in Sediment

PAHs are a group of SVOCs common in urban runoff (Mahler et al., 2005) that can have adverse effects on aquatic life including plants, invertebrates, and fish. The effects of exposure vary but can include organ damage, reproductive harm, or immune system weakening (Mahler et al., 2005). Coal-tar parking lot sealants have been identified as a significant source of PAHs in urban waterways and were banned from use in areas surrounding the recharge zone of the Edwards Aquifer within Comal and Hays counties by the EAA in 2012. In each sample year thus far, levels of total PAH in sediment samples have exceeded TECs and PECs at several sites especially in the San Marcos Springs complex. Further investigation may be warranted to identify the extent of PAH presence in the sediment and to identify potential sources of PAHs.

Selenium in Sediment

The selenium concentration in sediment sample HSM320 of 4.20 J mg/kg is below the laboratory reporting and quantification limit. The concentration in sample HSM320 is above the TSBC level of 0.3 mg/kg. Sediment studies of selenium concentrations have shown that levels of 4 mg/kg or less are not likely to bioaccumulate in the food chain or have adverse impacts on the reproduction of fish or aquatic birds (Lemly, 1995; Moore et al., 1990; Van Derveer and Canton, 1996). Selenium detections did exceed this amount at HSM320 with a detected concentration of 4.20 J mg/kg.

DEHP in Sediment

DEHP was detected in the majority of sediment samples from the Comal and San Marcos springs complexes in 2013, but was considered by EAA to likely be a laboratory or sampling equipment artifact. DEHP was again detected in some samples from the San Marcos Springs complex in 2014 at HSM320, HSM330, and HSM350. Because other plasticizers were not present and different laboratories were used, it was a conclusion in 2014 that DEHP may be present in sediments in the San Marcos Springs complex. In 2015, equipment used to collect sediment samples did not contain plasticizers with the exception of one sample collected in the Comal River (HCS330). DEHP was again detected in some of the same San Marcos Springs complex samples (HSM330, HSM340, and HSM350). The continued detections of DEHP indicate that it may be present in the middle reaches of the San Marcos Springs complex.

Recommendations

Based on the data collected for this report combined with completion of several sample events, a few points worth consideration in future years are suggested. Although this data set adds to an initial baseline established by the 2013 and 2014 sampling programs, currently the data set is not adequate to make significant assessments regarding water or sediment quality. The current data set is extremely important for use in developing a baseline such that in the future, additional information can be added to it, allowing interested parties to better understand the surface water quality, stormwater impacts, and sediment quality trends of each system. In addition, the RTIs are very valuable for gathering basic water quality data and tracking any changes to basic water quality over time.

Based on this third year of work, the following changes should be considered for the 2016 EAHCP Workplan.

- A more detailed study of the presence of PAH compounds in the middle reaches of the San Marcos Springs complex should be considered to evaluate potential sources of the PAH compounds and the vertical and lateral extents of impacted areas.
- If the concentrations of DEHP detected in sediments are determined to be harmful to the endangered species, then additional sampling and analysis should be conducted to evaluate the vertical and lateral extents of impacted areas.

9.0 DATA QUALITY OBJECTIVES

SWCA evaluated each sampling event to determine if procedures should be modified to improve data collection to ensure data quality objectives were met. A discussion of problems encountered, deviations to the workplan, and resolutions to these circumstances are provided in Appendix C. The only ongoing challenge recognized is the inability to consistently deliver *E. coli* samples to a laboratory within hold times during stormwater sampling events. This inability is inherent to stormwater sampling events due to the occurrence of storms during non-working hours. Special runners are utilized to deliver samples to the laboratory as early as possible to minimize hold time exceedances.

Based on procedures implemented to correct or improve data collection methods and the relatively low significance of the deviations, the circumstances described in Appendix C do not compromise the integrity of the study or this report.

10.0 DEFINITIONS

| | |
|------------------|---|
| Alkalinity | The capacity of water to neutralize acids, a property imparted by the water's content of carbonate, bicarbonate, hydroxide, and on occasion borate, silicate, and phosphate. It is expressed in milligrams per liter of equivalent calcium carbonate (mg/l CaCO ₃). |
| Aquifer | Underground geological formation or group of formations containing water; source of groundwater for wells and springs. |
| ASTM | Abbreviation for American Society for Testing and Materials. A nonprofit organization that develops and publishes approximately 12,000 technical standards, covering the procedures for testing and classification of materials of every sort. |
| Bacteria | Microscopic living organisms that can aid in pollution control by metabolizing organic matter in sewage, oil spills, or other pollutants. However, certain bacteria in soil, water, or air can also cause human, animal, and plant health problems. |
| Basin | Any area draining to a point of interest. |
| Baseline data | Initial data generated by consistent monitoring of the same sites over time. |
| Caffeine | A stimulant drug found naturally in coffee, tea, and chocolate, and also within soft drinks and other foods. If detected, it might indicate an anthropogenic source of water impacts. |
| Channel | A long, narrow excavation or surface feature that conveys surface water and is open to the air. |
| Deionized water | Water with all ions removed. |
| Detection limit | The lowest concentration of a given pollutant that an analytical method or equipment can detect and still report as greater than zero. Generally, as readings approach the detection limit, they become less and less reliable quantitatively. |
| Dissolved solids | The total amount of dissolved material, organic, and inorganic, contained in water or wastewater. Measurements are expressed as ppm or mg/L. |
| DO | Abbreviation for dissolved oxygen. Oxygen molecules that are dissolved in water and available for living organisms to use for respiration. Usually expressed in milligrams per liter or percent of saturation. The concentration of DO is an important environmental parameter contributing to water quality. |
| DOC | Abbreviation for dissolved organic carbon, a broad classification of organic molecules of varied origin and composition within aquatic systems. Organic |

carbon compounds are a result of decomposition processes from dead organic matter, such as plants.

| | |
|-----------------|--|
| DQO | Abbreviation for data quality objectives, a process used to develop performance and acceptance criteria or data quality objectives that clarify study objectives, define the appropriate type of data, and specify tolerable levels of data needed to support decisions. |
| Drainage | The collection, conveyance, containment, and/or discharge of surface and stormwater runoff. |
| EARIPHCP | Abbreviation for Edwards Aquifer Recovery Implementation Program Habitat Conservation Plan. |
| Endpoint | That state in titration at which an effect, such as a color change, occurs, indicating that a desired point in the titration has been reached. |
| Equipment blank | Sample used to assess the effectiveness of the decontamination process on sampling equipment. The equipment blank is prepared by pouring reagent-grade water over/through sampling equipment and analyzing for parameters of concern (to match the sampling routine applicable to the site). |
| Field duplicate | Second sample collected simultaneously from the same source as the parent sample, but which is submitted and analyzed as a separate sample. This sample should generally be identified such that the laboratory is unaware that it is a field duplicate. |
| Filtration | The process of separating solids from a liquid by means of a porous substance (filter) through which only the liquid can pass. |
| Groundwater | Water found beneath Earth's surface that fills pores between materials, such as sand, soil, or gravel. |
| Habitat | The specific area of environment in which a particular type of plant or animal lives and grows. |
| HCP | Abbreviation for Habitat Conservation Plan. A planning document that is required by the United States Fish and Wildlife Service as part of their enforcement of the Endangered Species Act. |
| LCS/LCSD | Abbreviation for Laboratory control samples and laboratory control sample duplicate. LCS/LCSD are evaluated to assess overall method performance and are the primary indicators of laboratory performance. In general, laboratory control samples are similar in composition as the environmental samples, contain known concentrations of all the analytes of interest, and undergo the same preparatory and determinative procedures as the environmental samples. A |

LCS/LCSD may be analyzed to provide information on the precision of the analytical method.

| | |
|---------------|--|
| MS/MSD | Abbreviation for matrix spike/matrix spike duplicate. MS/MSD results are examined to evaluate the impact of matrix effects on overall analytical performance and potential usability of the data. A matrix spike is a representative environmental sample that is spiked with target analytes of interest prior to being taken through the entire analytical process in order to evaluate analytical bias for an actual matrix. A matrix duplicate is a collected (e.g., a VOC soil sample) or a homogenized sample that is processed through the entire analytical procedure in order to evaluate overall precision for an actual matrix. |
| MDL | Abbreviation for method detection limit, minimum concentration of a substance that can be measured and reported with 99 percent confidence that the analyte concentration is greater than zero, as determined from analysis of a sample containing the analyte in a given matrix. |
| MPN | Abbreviation for most probable number. An analytical method used to detect the presence of coliforms in a water sample and estimate their numbers. |
| PCBs | Abbreviation for polychlorinated biphenyls. Group of more than 200 chlorinated toxic hydrocarbon compounds that can be biomagnified. |
| PCL | Abbreviation for protective concentration levels which is established to protect human health. |
| Peak | Maximum instantaneous flow at a specific location resulting from a given storm condition. |
| pH | A measure of the alkalinity or acidity of a substance. Also defined as “the negative logarithm of the hydrogen ion concentration ($-\log_{10}[\text{H}^+]$) where H^+ is the hydrogen ion concentration in moles per liter. The pH of a substance is neutral at 7.0, acidic below 7.0, and alkaline above 7.0. |
| PQL | Abbreviation for practical quantitation limit, which is the smallest concentration of the analyte that can be reported with a specific degree of confidence. |
| Precipitation | The discharge of water, in liquid or solid state, out of the atmosphere, generally upon a land or water surface. Precipitation includes rainfall, snow, hail, and sleet. |
| Precision | The ability of a measurement to be consistently reproduced. |
| QA/QC | Abbreviation for quality assurance/quality control. The total integrated program for assuring reliability of monitoring and measurement data. |

| | |
|----------------|--|
| Recession | End of runoff event, which is defined as the point in time when the recession limb of the hydrograph is <2 percent of the peak or is within 10 percent of the prestorm base flow, whichever is greater. |
| RPD | Abbreviation for relative percent difference. The RPD provide a measure of precision. |
| Representative | Said of samples collected that are similar to those of groundwater in its in situ condition. |
| RL | Abbreviation for reporting limit, the smallest concentration of an analyte reported by the laboratory to a customer. The RL is never less than the PQL and is generally twice the MDL. |
| Runoff | Precipitation, snow melt, or irrigation water that runs off the land into surface water. Runoff can carry pollutants from the air and land into the receiving waters. |
| Sediment | Fragmental material that originates from weathering of rocks and is transported by, suspended in, or deposited by water or air. |
| Shelby Sampler | A thin-walled tube with a cutting edge at the toe. A sampler head attaches the tube to the drill rod and pressure vents. Generally used in cohesive soils. Soil or sediment sampled from this sampler is considered undisturbed. |
| Spring | Water coming naturally out of the ground. |
| Stormwater | Stormwater is the water that runs off surfaces such as rooftops, paved streets, highways, and parking lots. It can also come from hard, grassy surfaces such as lawns, play fields, graveled roads, and parking lots. |
| Surface water | That which forms and remains above ground, such as lakes, ponds, rivers, streams, bays, and oceans. |
| SVOC | Abbreviation for semi-volatile organic compounds, which is a group of chemicals composed primarily of carbon and hydrogen that have a relatively low tendency to evaporate (volatilize) into the air from water or soil. Some of the compounds that make up asphalt are examples of SVOCs. |
| TDS | Abbreviation for total dissolved solids, or the total amount of all inorganic and organic substances, including minerals, salts, metal, cations, or anions that are dispersed within a volume of water. |
| Temporal | Over a period of time. |
| TKN | Abbreviation for total Kjeldahl nitrogen, which is the total concentration of organic and ammonia nitrogen in wastewater. |

| | |
|--------------|--|
| TOC | Abbreviation for total organic carbon, which is the gross amount of organic matter found in natural water. Suspended-particulate, colloidal, and dissolved organic matter are part of the TOC measurement. Settable solids consisting of inorganic sediments and some organic particulate are not transferred from the sample by the lab analyst and are not part of the TOC measurement. |
| TSBC | Texas-specific Background Concentrations as established by the Texas Commission on Environmental Quality. |
| Turbidimeter | An instrument for measuring turbidity in which a standard suspension is used for reference. |
| Turbidity | A measure of how clear the water is; how much the suspended material in water results in the scattering and absorption of light rays. An analytical quantity is usually reported in turbidity units and determined by measurements of light diffraction. Material that can increase turbidity (reduce clarity of water) are suspended clay, silt, sand, algae, plankton, microbes, and other substances. |
| Trip blank | Sample known to be free of contamination (for target analytes) that is prepared in the laboratory and treated as an environmental sample after receipt by the sampler. Trip blank samples are applicable to VOC analysis only. |
| TSS | Abbreviation for total suspended solids, which are the nonfilterable residue retained on a glass-fiber disk filter mesh measuring 1.2 micrometers after filtration of a sample of water or wastewater. |
| USGS | Abbreviation for United States Geological Services. USGS is a federal research organization which provides impartial information on health of ecosystems and environment, natural hazards which may threaten us, natural resources, impacts of climate and land use change, and core science systems which provide timely, relevant, and useable information. |
| VOC | Abbreviation for volatile organic compounds, which are often used as solvents in industrial processes and are either known or suspected carcinogens or mutagens. The five most toxic are vinyl chloride, tetrachloroethene, trichloroethene, 1,2-dichloroethane, and carbon tetrachloride. |
| Whirl-Pak® | Sterilized, clear polyethylene bag used to collect water samples for analysis. |
| WQAL | Abbreviation for a list of parameters defined as the following: pH, conductivity, temperature, dissolved oxygen, turbidity, and alkalinity in the field. Other parameters submitted for laboratory analysis include cations, anions, nutrients, metals, VOCs, SVOCs, herbicides and pesticides, bacteria, TOC, PCBs, and phosphorous. |

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APPENDIX A

**EDWARDS AQUIFER HABITAT CONSERVATION PLAN
(EAHCP) WORKPLAN**

2015 Water Quality Monitoring Program Strategy for Comal Springs and San Marcos Springs in Support of the Edwards Aquifer Habitat Conservation Plan

March 25, 2014

Prepared by

Edwards Aquifer Authority

INTRODUCTION

This Work Plan details the sampling strategy and protocols for surface water quality monitoring in 2015 for the Edwards Aquifer Habitat Conservation Plan (EAHCP) document (Section 5.7.2) implemented by the Edwards Aquifer Authority (EAA), utilizing a third party contractor. The goal of the water quality monitoring program, first implemented in 2013, is to detect water quality impairments that may negatively impact the listed species. In the event that certain constituents of concern are detected at levels indicating the potential for adverse effects, the Implementing Committee member with jurisdictional authority will be consulted to identify sources and consider Best Management Practices (BMPs) to reduce and/or eliminate the constituents of concern. If necessary, additional testing could be included in the current or following year to assist in determining the source of contamination and the Science Committee could be consulted to assist with BMP identification and source determination.

SCOPE OF WORK

The Water Quality Monitoring Program described below includes surface water, storm water, groundwater, and sediment sampling within Comal Springs and San Marcos Springs and associated river systems. Sample collection and analyses performed by EAA staff in 2013, has been performed by a contractor since 2014.

For 2015, the contractor will use the same sampling locations used in 2014 and as shown in the attached figures. However, changes in springflow, surface water runoff, land use, site security and access may dictate minor modification to sample collection locations and schedules as sampling efforts progress. Any minor changes resulting from these factors that are necessary as a result of safety or equipment concerns will be noted in the field sample sheets and dedicated field books. Should logistics or safety issues require any significant changes to this workplan, the sampling contractor shall report those issues to the EAA. Subsequently, the EAA will present those changes to the Science and Implementing committees for review and approval as needed prior to their implementation.

COMAL SPRINGS

Comal Springs discharges an average of 291 cubic feet second (cfs) into Landa Lake, located within the city of New Braunfels, Texas. Comal Springs is considered a spring complex with multiple discharge points along the 4,500 foot reach of Landa Lake. The springs issue from the Edwards Group limestone along the 4,500-foot section of the northeast-southwest trending escarpment formed by the Comal Springs Fault. Landa Lake forms the headwaters of the Comal River which flows approximately two miles before entering the Guadalupe River.

Discharge measurements have been collected from Comal Springs since 1933 and the EAA has been collecting water quality samples for more than ten years. In recent years, the EAA has been collecting samples from Spring 1, Spring 3, and Spring 7 on a quarterly basis during normal flow conditions and on a monthly basis when the San Antonio pool critical period triggers have been reached. Spring 1, Spring 3, and Spring 7 discharge into Landa Lake and make up part of the Comal Springs complex. Figure 1 indicates these historical groundwater sampling locations. Water quality samples are collected and analyzed for: dissolved oxygen (DO), pH, conductivity, and temperature, in the field and for alkalinity¹. Samples are also submitted to the EAA contract laboratory for analysis of cations, anions, nutrients, metals, VOCs, SVOCs, herbicides and pesticides, bacteria, TOC, PCBs, and phosphorous. This list of parameters is defined as the water quality analytical list (WQAL).

Sampling Methods

All samples will be collected following the EAA's *Field Sampling Plan* or contractor established methodology upon approval by the EAA. Samples shall be analyzed by a NELAP accredited contract laboratory. To date no requests to deviate from the EAA's *Field Sampling Plan* have been received or approved.

EAHCP Surface Water Sampling Locations

To comply with sampling requirements outlined in the EAHCP, five additional surface water sampling locations (Figure 2) were identified in the 2013 EAHCP Water Quality Work Plan for intensive monitoring in the Landa Lake and Comal River area as listed below:

- Upper Springs (near Bleiders Creek);
- Upper Landa Lake - (near Spring Island);
- Lower Landa Lake - (above outfalls);
- Upper Old Channel - (Elizabeth Street); and,
- USGS Gauge - (above San Antonio Street Bridge).

¹ Alkalinity analysis will be conducted within eight hours of sample collection.

Surface Water Sampling Frequency

In 2015, water samples will be collected twice from each of the five surface water locations listed above. The interval between sampling events will be approximately six months. Water samples will be analyzed for the WQAL parameters and caffeine using a National Environmental Laboratory Accreditation Program (NELAP) laboratory. A listing of analytical parameters is provided in Appendix A.

Sediment Sampling

One sediment sample will be collected during 2015 from each of the surface water sampling locations (Figure 2). Three samples will be collected from each sample site and composited into a single sample for analysis (to minimize VOC loss, it is recommended the compositing process be performed at the laboratory). Sediment samples will be analyzed for the analytical parameters provided in Appendix B. Results of sediment sampling analysis will be used to formulate future sediment sampling at Landa Lake and the Comal River. Sediment samples will be collected from zero to three inches below the surface for calendar year 2015. Sediment sample intervals will likely vary in subsequent sample years based on the results of each year of sediment analyses.

Storm Water Sampling Program for Comal Springs

Two storm water sampling events will be performed in 2015 to evaluate storm water quality from the urban landscape. A storm water sampling event will be triggered when a local rainfall event causes a significant increase in spring flow at the historic Comal Springs gauging station or changes in two water quality parameters at any of the real time water quality monitoring stations. Three water quality samples will be collected and analyzed from each surface water sampling location during the sampling event. Sample times will be spaced to reflect changes in the stream hydrograph (initial rise or first flush, peak flow, and recession limb). Water samples will be analyzed for the WQAL parameters. A listing of analytical parameters for storm water samples is provided in Appendix A.

The following locations will be sampled for storm water as indicated on Figure 3:

- Upper Springs (near Blieders Creek);
- New Channel – (below confluence with Dry Comal Creek);
- Upper Old Channel - (at Elizabeth Street);
- Lower Old Channel - (above Hinman Island); and,
- Comal River - (above confluence with Guadalupe River).

Groundwater Sample Collection for Extreme Low Flow Scenarios for Comal Springs

In the event flow at Comal Springs drops below 30 cfs, the EAHCP (6.4.3.3) calls for weekly monitoring of three wells in the vicinity of the spring complex for DO, conductivity, pH, and temperature. Should springflow drop below 20 cfs, then additional parameters to include nutrients,

EAHCP Water Quality Sampling Plan

TDS, and TOC are to be added to the weekly sample regimen. Analytical parameters for all low flow sampling is included in Appendix A. Based on conditions during the drought of record (circa 1950s), sampling for lower flow scenario could last for up to 21 weeks. The three specific wells to be used will be determined at the time of low flow sample initiation, based on well conditions and aquifer levels.

Real Time Instrument Water Quality Data Logging Program for Comal Springs

Continuous water quality monitoring stations will continue in 2015 at the following locations indicated on Figure 4:

Spring Run 3;
Spring 7; and,
New Channel (below confluence with Dry Comal Creek).

Monitoring will be performed using a data logging sonde capable of collecting data on 15 minute intervals. The parameters measured will include temperature, dissolved oxygen, pH, turbidity, and conductivity. These data will be evaluated to identify short-term and long-term water quality variations of the spring system as well as changes in water quality related to storm water runoff.

This monitoring effort will continue to be performed by EAA staff in 2015.

Collection of Passive Diffusion Samples (Passive Samples) at Comal Springs

Passive samples are to be collected during the 2015 sampling effort using a passive diffusion type sampling device. Devices will be obtained from Amplified Geochemical Imaging LLC (AGI), or equivalent to the AGI device for functionality and analytical parameters. Sample locations for PDS samples are provided in Figure 5. The passive sampling effort shall be performed in February, April, June, August, October, and December. The devices shall be installed for a two-week interval at the same locations as the base flow surface water samples. Specifically at the sample points below.

Upper Springs (near Bleiders Creek);
Upper Landa Lake - (near Spring Island);
Lower Landa Lake - (above outfalls);
Upper Old Channel - (Elizabeth Street); and,
USGS Gauge - (above San Antonio Street Bridge).

The general parameter set for PDS samples is listed in Appendix A, under *Analytical Parameters for Passive Diffusion Samplers, Comal and San Marcos Springs*.

SAN MARCOS SPRINGS

Located in San Marcos, Texas on the campus of Texas State University, San Marcos Springs discharges an average of 176 cfs into Spring Lake. The springs issue from the Edwards Group limestone along the northeast-southwest trending escarpment formed by the San Marcos Springs Fault. Spring Lake forms the headwaters of the San Marcos River. Discharge measurements have been collected from San Marcos Springs since 1957 and the EAA has been collecting water quality samples for more than ten years.

In recent years, the EAA has been collecting samples from Deep Spring and Hotel Spring on a quarterly basis during normal flow conditions and on a monthly basis when the San Antonio pool critical period triggers have been reached. Both Deep and Hotel springs are located in the bed of Spring Lake and make up part of the San Marcos Springs complex. Figure 6 indicates these historical groundwater sample locations at San Marcos Springs. Water quality samples are collected and analyzed for: dissolved oxygen (DO), pH, conductivity, and temperature, in the field and for alkalinity². Samples are also submitted to the EAA contract laboratory for analysis of cations, anions, nutrients, metals, VOCs, SVOCs, herbicides and pesticides, bacteria, TOC, PCBs, and phosphorous. For the purposes of EAHCP related water sampling, the analyte caffeine has been added to the list of analyzed parameters. This list of WQAL parameters is an identical to the list of parameters analyzed for at Comal Springs.

Sampling Methods

All samples will be collected following the EAA's *Field Sampling Plan* or contractor established methodology upon approval by the EAA. Samples shall be analyzed by a NELAP accredited contract laboratory. To date no requests to deviate from the EAA's *Field Sampling Plan* have been received or approved.

Surface Water Sampling Locations

To comply with sampling requirements outlined in the EAHCP document, seven additional surface water sampling locations (Figure 7) were identified in the 2013 HCP work plan for intensive monitoring as listed below:

- Sink Creek;
- Spring Lake;
- Sessoms Creek;
- City Park;
- Rio Vista Dam;
- I-35 reach; and,
- Capes Dam/Willow Creek.

² Alkalinity analysis will be conducted within eight hours of sample collection.

Surface Water Sampling Frequency

In 2015, water samples will be collected twice from each of the seven surface water locations listed above. The interval between sampling events will be approximately six months. Water samples will be analyzed for the WQAL parameters and caffeine using a National Environmental Laboratory Accreditation Program (NELAP) laboratory. A listing of analytical parameters is provided in Appendix A.

Sediment Sampling

One sediment sample will be collected during 2015 from each of the surface water sampling locations (Figure 7). Three samples will be collected from each sample site and composited into a single sample for analysis (it is recommended that compositing of the sample be performed at the laboratory to minimize VOC loss). Sediment samples will be analyzed for the analytical parameters provided in Appendix B. Results of sediment sampling analysis will be used to formulate future sediment sampling at Spring Lake and the San Marcos River. Sediment samples will be collected from zero to three inches below the surface for calendar year 2015. Sediment sample intervals will vary in subsequent sample years based on the results of each year of sediment analyses.

Storm Water Sampling Program for San Marcos Springs

Two storm water sampling events will be performed in 2015 to evaluate storm water runoff from the urban landscape. A storm water sampling event will be triggered when a local rainfall event causes a significant increase in spring flow at the San Marcos Springs gauging station or changes in two water quality parameters at any of the real time water quality monitoring stations. Three water quality samples will be collected from each surface water sampling location during the sampling event. Sampling times will be spaced to reflect changes in the stream hydrograph (initial rise or first flush, peak flow, and recession limb). Water samples will be analyzed for the WQAL parameters. A listing of analytical parameters is provided in Appendix A, for storm water samples.

The following locations will be sampled for storm water as indicated on Figure 8:

- Sink Creek;
- Sessoms Creek;
- Dog Beach Outflow;
- Hopkins Street Outflow;
- Purgatory Creek (above San Marcos River);
- I-35 Reach; and
- Willow Creek (above San Marcos River).

Groundwater Sample Collection for Extreme Low Flow Scenarios for San Marcos Springs

In the event flow at San Marcos Springs drops below 50 cfs, the EAHCP (6.4.4.3) calls for weekly monitoring of three wells in the vicinity of the spring complex for DO, conductivity, pH, and temperature. Should springflow drop below 30 cfs, then additional parameters to include Nutrients, TDS, and TOC are to be added to the sample regimen. Analytical parameters for all low flow sampling is included in Appendix A. Based on conditions during the drought of record (circa 1950s), sampling for lower flow scenario could last for up to 21 weeks. The three specific wells to be used will be determined at the time of sampling abased on well conditions and aquifer levels.

Real Time Instrument Water Quality Data Logging Program for San Marcos Springs

Continuous water quality monitoring stations were established in 2013 and will continue in 2015 at the following locations indicated on Figure 9:

USGS gauging station; and,
Rio Vista Dam.

Monitoring will be performed using a data logging sonde capable of collecting data on 15 minute intervals. The parameters measured will include temperature, dissolved oxygen, pH, turbidity, and specific conductance. These data will be evaluated to identify short-term and long-term water quality variations of the spring system as well as changes in water quality related to storm water runoff. **Continuous water quality monitoring stations will be operated and maintained by EAA.**

In 2015, an additional water quality data logging point is recommended for installation at the lower (south) end of the sample area for San Marcos. The additional station will help with the timing of storm sample collection as well as improved monitoring of the IH-35 and Willow Creek runoff impacts. Costs for this are included in Appendix C of this document. The location of the proposed new monitoring point is coincident with surface water sample point HSM170 (Capes Dam/Willow Creek area).

Collection of Passive Diffusion Samples (Passive Samples) at San Marcos Springs

Passive samples are to be collected during the 2015 sampling effort using a passive diffusion type sampling device. Devices will be obtained from AGI, or equivalent to the AGI device for functionality and analytical parameters. Sample locations for PDS samples are provided in Figure 10. The passive sampling effort shall be performed in February, April, June, August, October, and December. The devices shall be installed for a two-week interval at the same locations as the base flow surface water samples. Specifically at the sample points that follow.

EAHCP Water Quality Sampling Plan

Sink Creek;
Spring Lake;
Sessoms Creek;
City Park;
Rio Vista Dam;
I-35 reach; and,
Capes Dam/Willow Creek.

The general parameter set for PDS samples is listed in Appendix A, under *Analytical Parameters for Passive Diffusion Samplers, Comal and San Marcos Springs*.

WATER QUALITY MONITORING REPORT

The contractor will compile and present sampling results in an annual report to the EAA. The report will include an evaluation of analytical data, graphs of results that exceed comparative or regulatory standards, a discussion of water and sediment quality, laboratory reports and field data sheets, photographs, sampling locations and rationale, description of sampling methods, and a description and rationale for any deviations from the Water Quality Sampling Plan due to logistics or safety issues. The report is to be submitted in hard copy and electronically and will be reviewed internally by EAA. The deadline for submittal to the EAA is December 21, 2015.

DATA COMPILATION, ANALYSES AND REPORTING

All of the data collected as a result of the 2015 HCP Water Quality Monitoring Plan will be compiled, and analyzed, and the results will be presented to the Implementing Committee by February 15, 2016; prior to inclusion in the annual EAHCP Annual Report that is required by Sections 6.2.4 and 9.3 of the EAHCP and Section 11.1c of the Implementing Agreement. The report will include an evaluation of all analytical data, including graphs, key photographs and general summary of results.

CHANGES TO WORKPLAN FROM THE 2014 SAMPLING EFFORT

In summary, the workplan has few changes from 2014. Funding is needed to add a single RTI water quality data logger at the downstream end of the sampling area in San Marcos. Funding is also requested for maintenance and replacement needs for existing RTIs, as well as data transmission and web hosting fees. Detail for the RTIs is listed in Appendix C.

A change to the sample interval for sediment samples is included in this workplan. The former sample interval is 0 – 18-inches below surface. For 2015, the sediment sample interval will be 0-3-inches below the surface.

SCIENCE COMMITTEE REVIEW

EAHCP Water Quality Sampling Plan

This 2015 Water Quality Work Plan will be reviewed by the EAHCP Science Committee prior to implementation. The Science Committee will be asked to conform the need for the following additions or changes to the Water Quality Work Plan:

- An additional Real Time Instrument for water quality data logging will be added to the downstream end of the San Marcos Springs sampling area in 2015.
- A change to the sample interval for sediment samples from 0 – 18-inches below surface to 0-3-inches below the surface.

BUDGET

Table 7.1 Budget: \$200,000

Requested workplan Budget: \$504,530

2015 EAHCP Sampling as performed by an outside contractor, annual costs \$474,430.00

Real Time Instruments (RTI): \$30,100 (see Appendix C)

Justification for Budget Adjustment

The real time water quality data logging instrumentation is in need of funding for maintenance (warranties will expire in 2015), in addition spare instrumentation is needed to prevent extended down time in the event of catastrophic failure. The instruments also require funding for calibration fluids, batteries, and other incidental costs. Funding is also needed for the addition of a downstream San Marcos instrument. Cost details are provided in Appendix C.











Figure 6
San Marcos Springs Groundwater Sample Locations



Most of the historical EAA sampling records for San Marcos Springs pertains to the locations known as Hotel and Deep (spring vents). Other locations at San Marcos Springs may have a limited sample record.

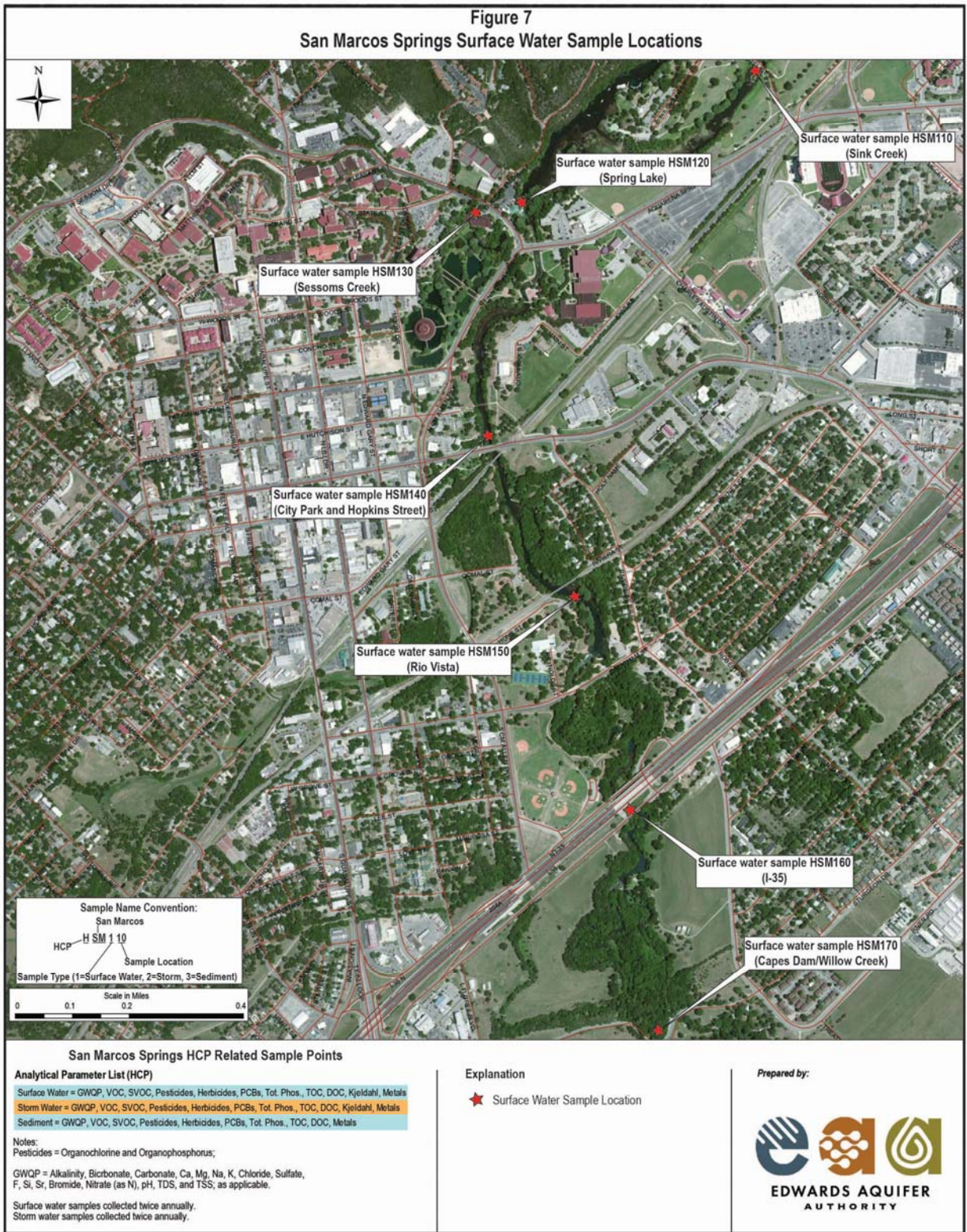
Samples are collected monthly during low flow conditions (critical period), and quarterly during normal conditions.

Explanation

★ Historical Groundwater (Spring) Sample Location

Prepared by:











EAHCP Water Quality Sampling Plan

Appendix A

Analytical Parameters for Assessing Water Quality from Storm Water and Surface Water Locations, Comal and San Marcos Springs

| Analyses | | |
|--|--|--|
| Volatile Organic Compounds (VOCs) | | |
| Semi-volatile Organic Compounds (SVOCs) | | |
| Organochlorine Pesticides | | |
| Polychlorinated Biphenyls (PCBs) | | |
| Organophosphorous Pesticides | | |
| Herbicides | | |
| Metals (Al, Sb, As, Ba, Be, Cd, Cr (total), Cu, Fe, Pb, Mn, Hg, Ni, Se, Ag, Tl, and Zn) | | |
| General Chemistry (GWQP) Total Alkalinity (as CaCO ₃), Bicarbonate Alkalinity (as CaCO ₃), Carbonate Alkalinity (as CaCO ₃); (Cl, Br, NO ₃ , SO ₄ , F ⁻ , pH, TDS, TSS, Ca, Mg, Na, K, Si, Sr, CO ₃), and Total Suspended Solids (TSS). | | |
| Phosphorus (total) | | |
| Total Organic Carbon (TOC), | | |
| Dissolved Organic Carbon (DOC) | | |
| Kjeldahl Nitrogen | | |
| Bacteria Testing (<i>E. coli</i>) | | |
| Caffeine | | |

| Method | Method Description | Protocol |
|----------|----------------------------------|-----------------------------|
| 8260B | Volatile Organic Compounds | (GC/MS) SW846 |
| 8270C | Semivolatile Organic Compounds | (GC/MS) SW846 |
| 8081B | Organochlorine Pesticides | (GC) SW846 |
| 8082A | Polychlorinated Biphenyls (PCBs) | by Gas Chromatography SW846 |
| 8141A | Organophosphorous Pesticides | (GC) SW846 |
| 8151A | Herbicides | (GC) SW846 |
| 6010B | Metals | (ICP) SW846 |
| 6020 | Metals | (ICP/MS) SW846 |
| 7470A | Mercury | (CVAA) SW846 |
| 300.0 | Anions, | Ion Chromatography |
| 340.2 | Fluoride | MCAWW |
| 365.4 | Phosphorus, | Total EPA |
| 9040C | pH | SW846 |
| 9060 | Organic Carbon, | Total (TOC) SW846 |
| SM 2320B | Alkalinity | SM |
| SM 2540C | Solids, | Total Dissolved (TDS) SM |
| SM 2540D | Solids, Total Suspended (TSS) | SM |
| 351.2 | Nitrogen, Total Kjeldahl | MCAWW |
| 1694 | Caffeine | |

Protocol References:

EPA = US Environmental Protection Agency

MCAWW = "Methods For Chemical Analysis Of Water And Wastes", EPA-600/4-79-020, March 1983 And Subsequent Revisions.

SM = "Standard Methods For The Examination Of Water And Wastewater",

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

EAHCP Water Quality Sampling Plan

Number of required QA/QC Samples for Surface (Base Flow) Sampling, Storm Sampling, and Sediment Sampling

| QA/QC Samples (Duples/EQ Blanks) | Equip. Blanks | Duples | Total |
|----------------------------------|---------------|-----------|-----------|
| Comal Surface Water= | 2 | 2 | 4 |
| San Marcos Surface Water= | 2 | 2 | 4 |
| Comal Storm Water= | 2 | 4 | 6 |
| San Marcos Storm Water= | 2 | 6 | 8 |
| Comal Sediments= | 1 | 1 | 2 |
| San Marcos Sediments= | 1 | 1 | 2 |
| Total Costs QA/QC Samples | 10 | 16 | 26 |

Analytical Parameters for Critical Period Related (Low Flow) Sampling of Water Wells, Comal and San Marcos Springs

| |
|--|
| Analyses |
| General Chemistry (GWQP) Total Alkalinity (as CaCO ₃), Bicarbonate Alkalinity (as CaCO ₃), Carbonate Alkalinity (as CaCO ₃); (Cl, Br, NO ₃ , SO ₄ , F, pH, TDS, TSS, Ca, Mg, Na, K, Si, Sr, CO ₃ ,) |
| Total Organic Carbon (TOC) |
| Total Dissolved Solids (TDS) |
| |

Analytical Parameters for Passive Diffusion Samplers, Comal and San Marcos Springs

| |
|---|
| PDS devices are to be placed at the locations listed Figures 5 and 10, for a two-week time period in the months of February, April, June, August, October, and December. |
| PDS devices will be from Amplified Geochemical Imaging, LLC, or equivalent and shall provide analyses for the following: TPH, BTEX, 1,3,5 and 1,2,4-trimethylbenzene, MTBE, phenanthrene, naphthalene1-methyl naphthalene, octane, cis and trans-1,2,-dichloroethene, 1,1-dichloroethane, chloroform, 1,1,1-trichloroethane, 1,2-dichloroethane, carbon tetrachloride, trichloroethene, tetrachloroethene, chlorobenzene, 1,4-dichlorobenzene, 1,1,2-trichloroethane, 1,1,1,2-tetrachloroethane, 1,1,2,2-tetrachloroethane, 1,3-dichlorobenzene, and 1,2-dichlorobenzene. |

EAHCP Water Quality Sampling Plan

Appendix B

Analytical Parameters for Assessing Water Quality from Sediment Sample Locations, Comal and San Marcos Springs

| Analyses |
|---|
| Volatile Organic Compounds (VOCs) |
| Semi-volatile Organic Compounds (SVOCs) |
| Organochlorine Pesticides |
| Polychlorinated Biphenyls (PCBs) |
| Organophosphorous Pesticides |
| Herbicides |
| Metals (Al, Sb, As, Ba, Be, Cd, Cr (total), Cu, Fe, Pb, Mn, Hg, Ni, Se, Ag, Tl, and Zn) |
| General Chemistry Total Alkalinity (as CaCO ₃), Bicarbonate Alkalinity (as CaCO ₃), Carbonate Alkalinity (as CaCO ₃); Ca, Mg, Na, K, Chloride, Sulfate, |
| Phosphorus (total) |
| Total Organic Carbon (TOC), |
| Dissolved Organic Carbon (DOC) |
| Bacteria Testing (<i>E coli</i>) |

| Method | Method Description | Protocol |
|----------|----------------------------------|-----------------------------|
| 8260B | Volatile Organic Compounds | (GC/MS) SW846 |
| 8270C | Semivolatile Organic Compounds | (GC/MS) SW846 |
| 8081B | Organochlorine Pesticides | (GC) SW846 |
| 8082A | Polychlorinated Biphenyls (PCBs) | by Gas Chromatography SW846 |
| 8141A | Organophosphorous Pesticides | (GC) SW846 |
| 8151A | Herbicides | (GC) SW846 |
| 6010B | Metals | (ICP) SW846 |
| 6020 | Metals | (ICP/MS) SW846 |
| 7470A | Mercury | (CVAA) SW846 |
| 300.0 | Anions, | Ion Chromatography |
| 340.2 | Fluoride | MCAWW |
| 365.4 | Phosphorus, | Total EPA |
| 9040C | pH | SW846 |
| 9060 | Organic Carbon, | Total (TOC) SW846 |
| SM 2320B | Alkalinity | SM |
| SM 2540C | Solids, | Total Dissolved (TDS) SM |
| SM 2540D | Solids, Total Suspended (TSS) | SM |

Protocol References:

EPA = US Environmental Protection Agency

MCAWW = "Methods For Chemical Analysis Of Water And Wastes", EPA-600/4-79-020, March 1983 And Subsequent Revisions.

SM = "Standard Methods For The Examination Of Water And Wastewater",

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

Appendix C

Estimated Costs for Addition of a Real Time Water Quality Monitoring Instrument at San Marcos Springs, and Ongoing Costs for Operation and Maintenance.

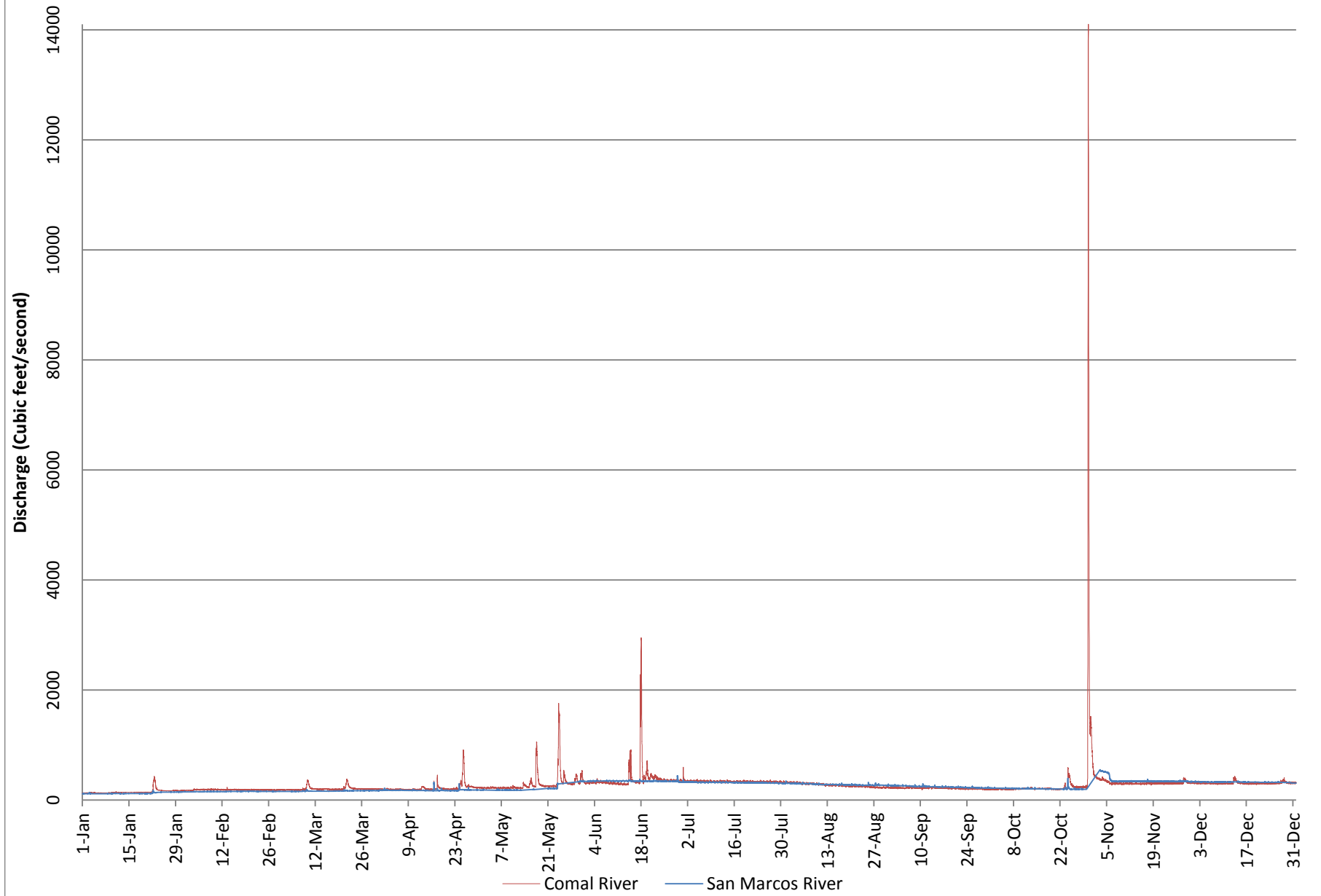
| | |
|---|---|
| Three new Eueka, Manta 2 Probe, equipped to monitor: DO, pH, Temperature, Conductivity, and Turbidity with Associated Netronix Telemetry System. | \$6,000.00 each for a total of \$18,000 |
| Annual maintenance costs for equipment, to include batteries (as needed), repairs, and calibration standards (estimated costs are for six total instruments, which includes the proposed new addition in San Marcos)* | \$6,000.00 |
| Annual data contract to include cellular data fees, and web hosting at Netronix site (estimated costs are for six total instruments, which includes the proposed new addition in San Marcos) | \$5,100.00 |
| Installation costs for proposed new unit to be located on the San Marcos system | \$1,000.00 |
| Total Estimated Costs for Real Time Water Quality Instrumentation calendar year 2015 | \$30,100.00 |

**Instrument warranties expires in 2015, extra funding needs to be earmarked for potential repairs.*

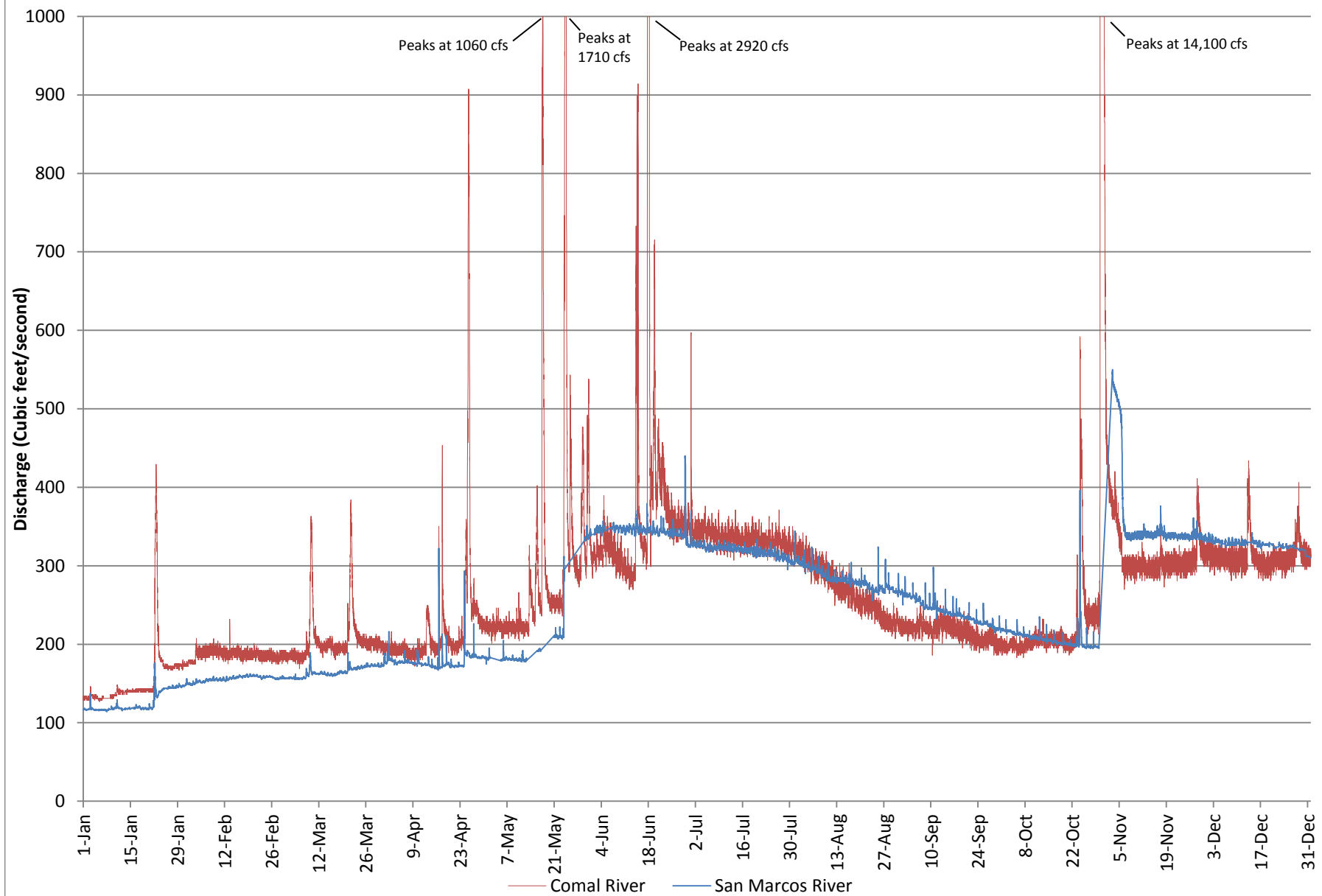
APPENDIX B

SPRING HYDROGRAPHS AND STORM WATER QUALITY GRAPHS

Discharge (cfs) Raw Data Plot 2015

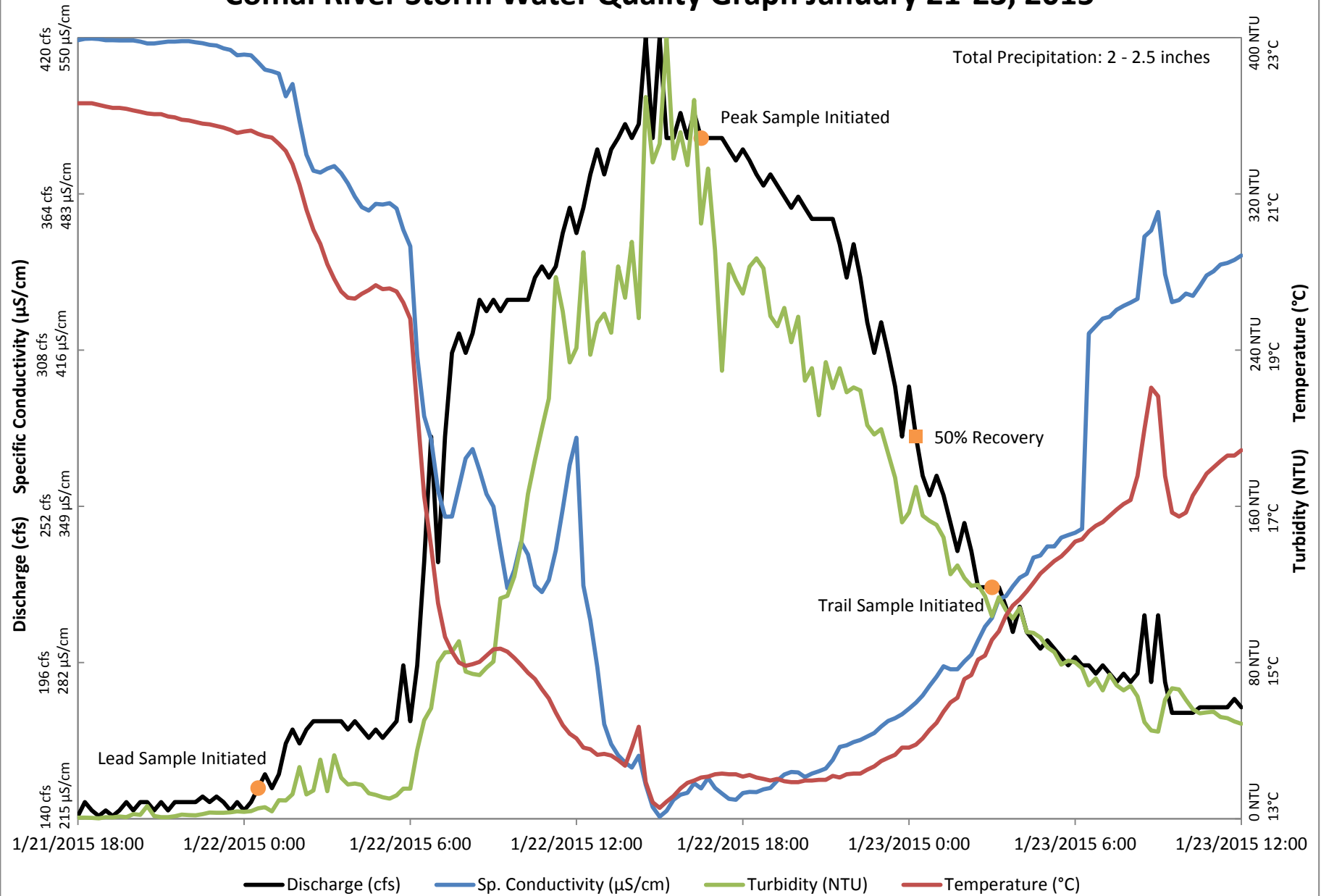


Discharge (cfs) Raw Data Plot 2015

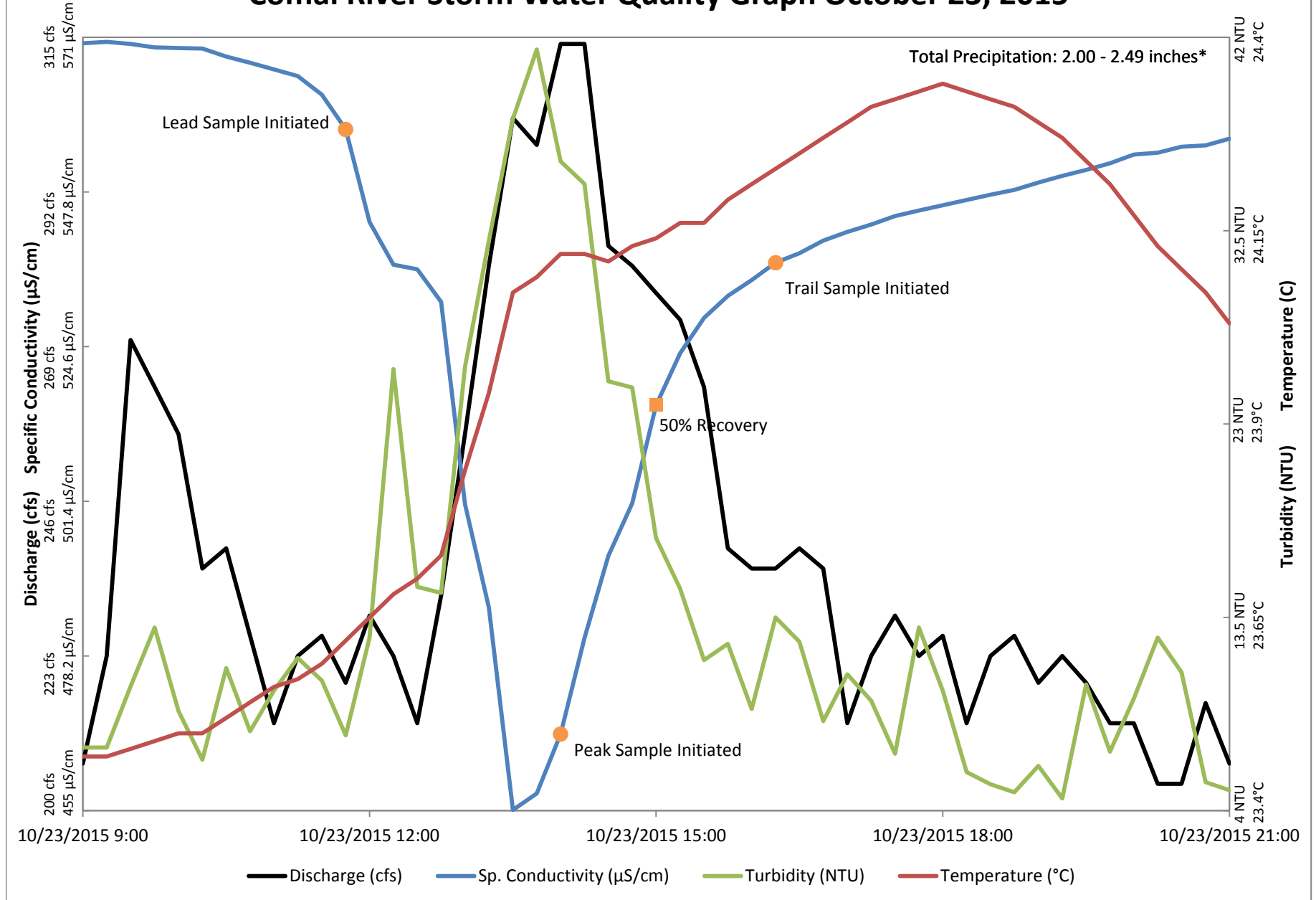


Comal River Storm Water Quality Graph January 21-23, 2015

Total Precipitation: 2 - 2.5 inches



Comal River Storm Water Quality Graph October 23, 2015

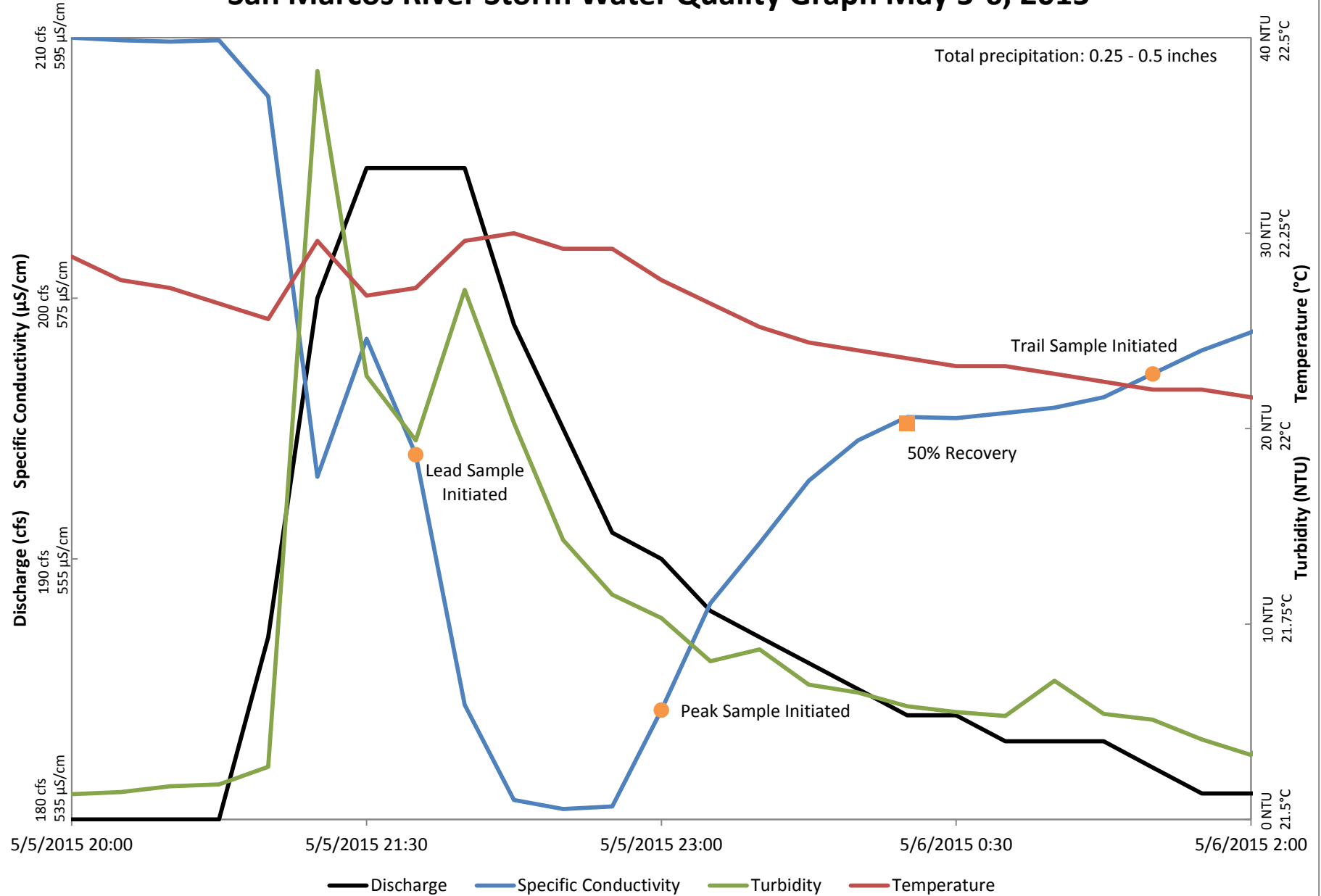


*2.00-2.49 inches total precipitation over 24 hour period according to NOAA. Sampling occurred during first pulse of storm, which might have totaled approximately 1 inch.

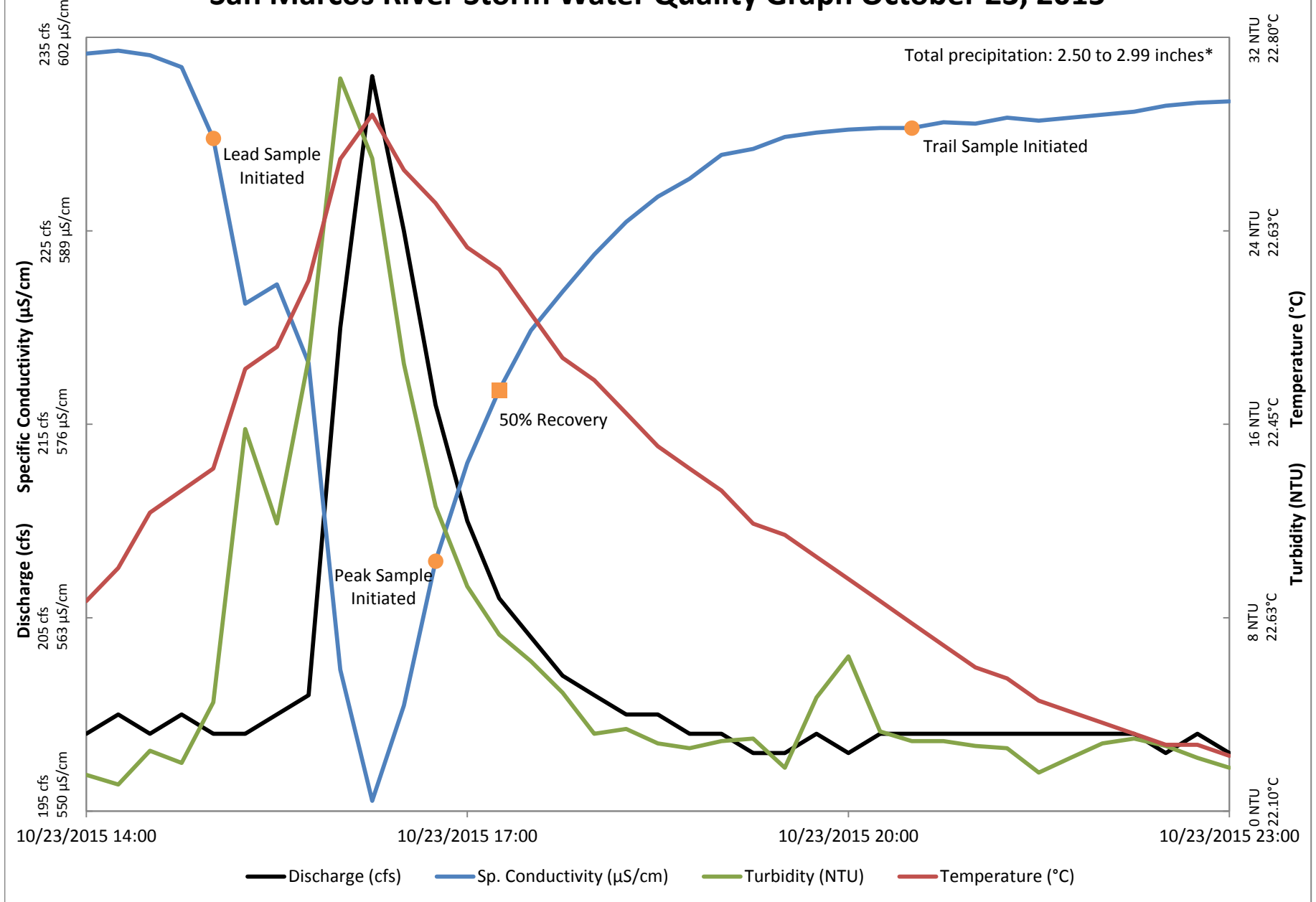
Data Source: Environet 2015, USGS 2015

San Marcos River Storm Water Quality Graph May 5-6, 2015

Total precipitation: 0.25 - 0.5 inches



San Marcos River Storm Water Quality Graph October 23, 2015



*2.50-2.99 inches total precipitation over 24 hour period according to NOAA. Sampling occurred during first pulse of storm, which might have totaled approximately 1 inch.

Data Source: Environet 2015, USGS 2015

APPENDIX C

DISCUSSION OF DEVIATIONS

Appendix C Discussion of Deviations

Comal Springs

Surface Water

Surface water was collected in the Comal Springs complex in March and September 2015. Surface water sampling locations did not deviate from the EAHCP Workplan.

March 16, 2015 and September 9, 2015, Collection

In accordance with the Workplan, filtration for metals and alkalinity samples was performed in the field using disposable bailers. Alkalinity analysis was performed at the SWCA office in San Antonio. Because the hold-time for pH analysis is only 15 minutes, laboratory analysis of pH was performed outside of hold time, but pH was measured in the field as part of the water quality parameter data set.

Stormwater

Stormwater events were sampled January 22–23 and October 23, 2015, in the Comal Springs complex. The ongoing drought made storm water sampling exceedingly difficult to perform. Rain events were generally scattered in nature and often too small in magnitude to generate sufficient runoff to sample. Storm water sampling locations did not deviate from those proposed in the EAHCP Workplan.

January 22–23, 2015, Event

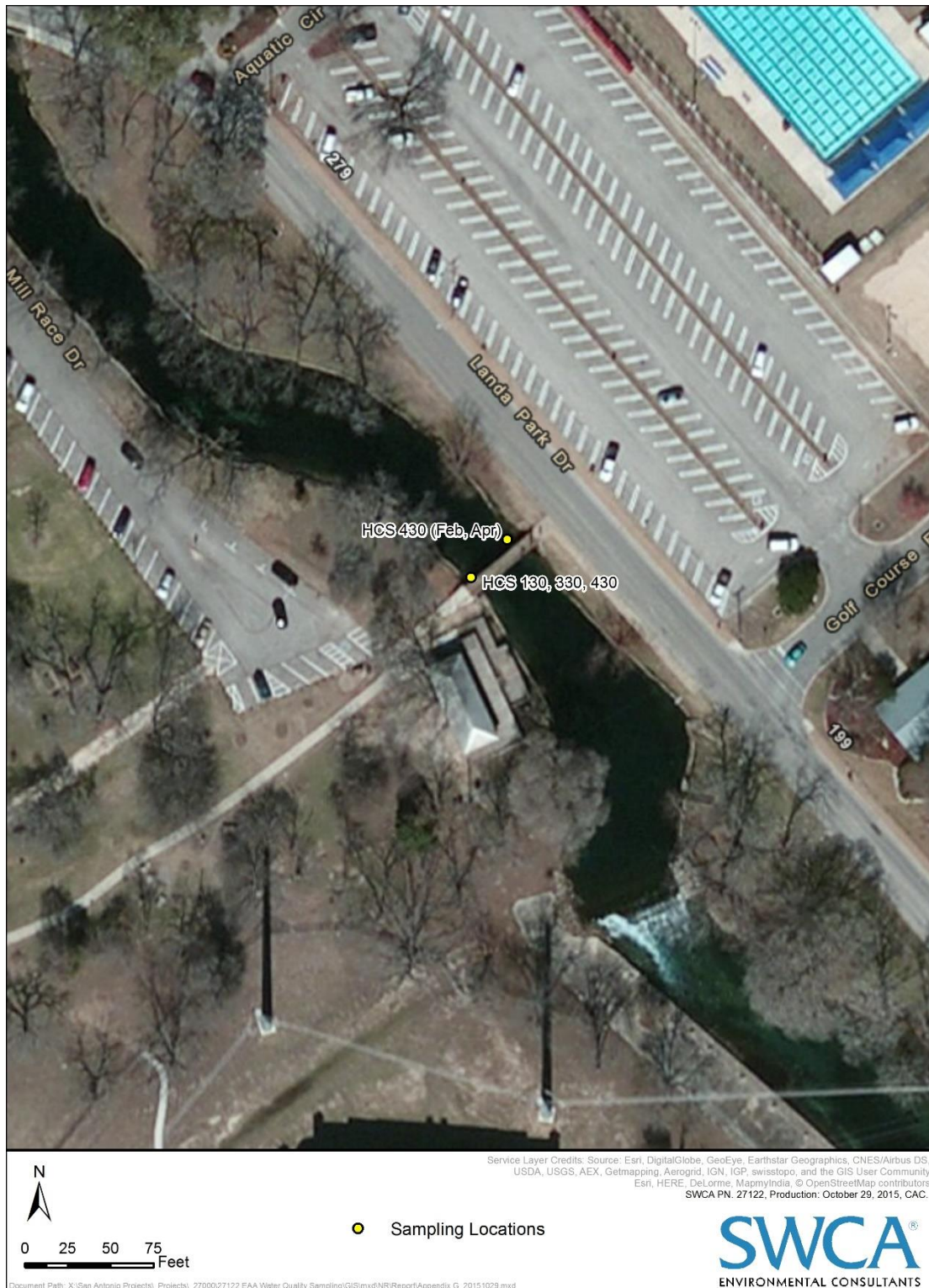
Lead sample collection began at approximately 0:30 on January 22, 2015 making it impossible for *E. coli* samples to reach the laboratory within the 6-hour hold time. Lead samples were submitted when the San Antonio River Authority (SARA) laboratory opened on the morning of January 22, 2015. The peak occurred at approximately 16:30 on January 22, 2015, also making it impossible for *E. coli* samples to reach the laboratory within the 6-hour hold time. Peak samples were submitted when the SARA laboratory opened on the morning of January 23, 2015. Trail samples were collected and delivered to the laboratory on January 23, 2015, within hold time. Because the hold-time for pH analysis is only 15 minutes, laboratory analysis of pH was performed outside of hold time, but pH was measured in the field as part of the water quality parameter data set.

Sediments

Sediment samples were collected from the Comal Springs complex on June 4, 2015. The EAHCP Workplan designated sediment collection sites to coincide with surface water collection points but allowed for some deviation from these locations as field conditions dictated. Sediment samples were collected more than 30 feet from surface water collection points due to limited amounts of available sediment at the surface water locations at HCS340 and HCS360. Sediment collection location deviations are discussed below.

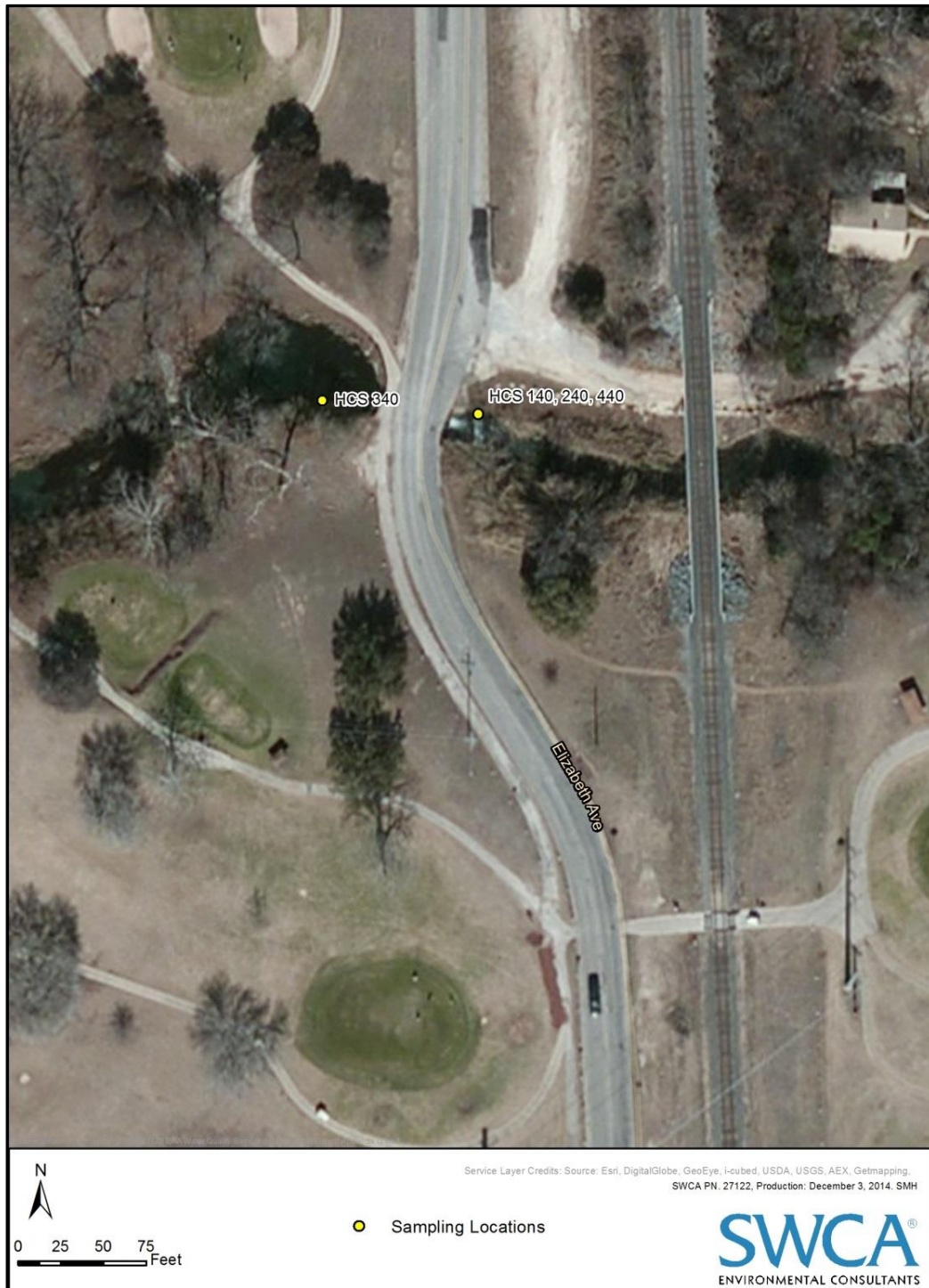
HCS330

This sediment sample was collected at the surface water location. Due to river depth, this sample was collected utilizing the core sampler.



HCS340

Surface water samples are collected from the eastern side of Elizabeth Avenue at this sample site, however the substrate in this area is very rocky with a strong current and little to no sediment. Sediment samples were collected upstream on the western side of Elizabeth Avenue, approximately 30 feet from the road. Sediment was collected approximately 10 feet from the southern shore.



HCS360

Surface water collection occurred on the upstream side of the bridge near the western bank at this location. Sediment was collected on the upstream side of the bridge approximately 12 feet from the eastern bank due to high river depth and limited available sediment near the surface water location. A field duplicate was also collected in the same location.



Passive Diffusion Samplers

Passive diffusion samplers (PDSs) were deployed at each of the surface water sample collection sites. When at all possible, deployment locations coincided with the surface water locations. Some adjustments had to be made to account for river depth, accessibility by SWCA staff for installation and retrieval, and interference by the public. PDSs were deployed for two-week periods during the months of February, April, June, August, and October 2015.

Storm events did occur during some PDS deployment periods, which may affect the analytes present in PDS results. Rain events large enough to cause an increase in discharge at the USGS Streamflow gauge in the Comal River occurred during PDS deployments on February 13, April 16, April 17, and June 14–15. Heavy rains and flooding occurred the week prior to the June 2015 deployment, discharge rates were higher throughout the deployment period than 2015 levels prior to June.

Any alterations to sample locations or lost PDS are discussed below.

HCS430

February 2015 – Bank stabilization construction continued from 2014 at site HCS430. PDS device deployed from Eastern bank of river.

April 2015 – Construction finished, however Western bank still inaccessible due to erosion protection measures, PDS device deployed from Eastern bank of river.

June 2015 – All construction and bank access restrictions had been cleared. PDS device was deployed from the Western bank of river, underneath a bridge to limit visibility to the public. This remained the normal deployment location for the remainder of 2015.

HCS460

April 2015 – PDS not recovered. The sample area was searched thoroughly but no sign of the deployment device or sampler was observed.

San Marcos

Surface water

Surface water was collected in the San Marcos Springs complex in March and September 2015. Surface water sampling locations did not deviate from the EAHCP Workplan. Because the hold time for pH analysis is only 15 minutes, laboratory analysis of pH was performed outside of the hold time. However, pH was measured in the field as part of the water quality parameter data set. In accordance with the Workplan, filtration for metals and alkalinity samples was performed in the field using disposable bailers. Alkalinity analysis was performed at the SWCA office in San Antonio.

Stormwater

Stormwater events were sampled May 5–6 and October 23, 2015, in the San Marcos Springs complex. Stormwater sampling locations did not deviate from those proposed in the EAHCP Workplan.

May 5–6, 2015, Event

All sample collection occurred outside of SARA laboratory operating hours making it impossible for the *E. coli* samples to reach the laboratory within the 6-hour hold time. All samples were submitted when the San Antonio River Authority (SARA) laboratory opened on the morning of May 6, 2015. SVOC analysis was not performed for FDHSM230 Trail, the bottle holding the sample for this analysis was broken in transit after sample collection. Because the hold-time for pH analysis is only 15 minutes, laboratory analysis of pH was performed outside of hold time. However, pH was measured in the field as part of the water quality parameter data set.

October 23, 2015, Event

Information will be added for the October event in the final draft of this report.

Sediment

Sediment samples were collected from the San Marcos Springs complex June 5, 2015. The EAHCP Workplan designated sediment collection sites to coincide with surface water collection points but allowed for some deviation from these locations as field conditions dictated. Sediment samples were collected more than 30 feet from surface water collection points due to limited amounts of available sediment and river depth at HSM310, HSM340, HSM350 and HSM370. Sediment collection location deviations are discussed below.

HSM310

Surface water samples were collected from the center of the downstream side of the West Laurel Street bridge. The bottom of this area has many large rocks and little exposed sediment. Sediment samples were collected approximately 6 feet upstream of the bridge, approximately 5 feet from the eastern bank.



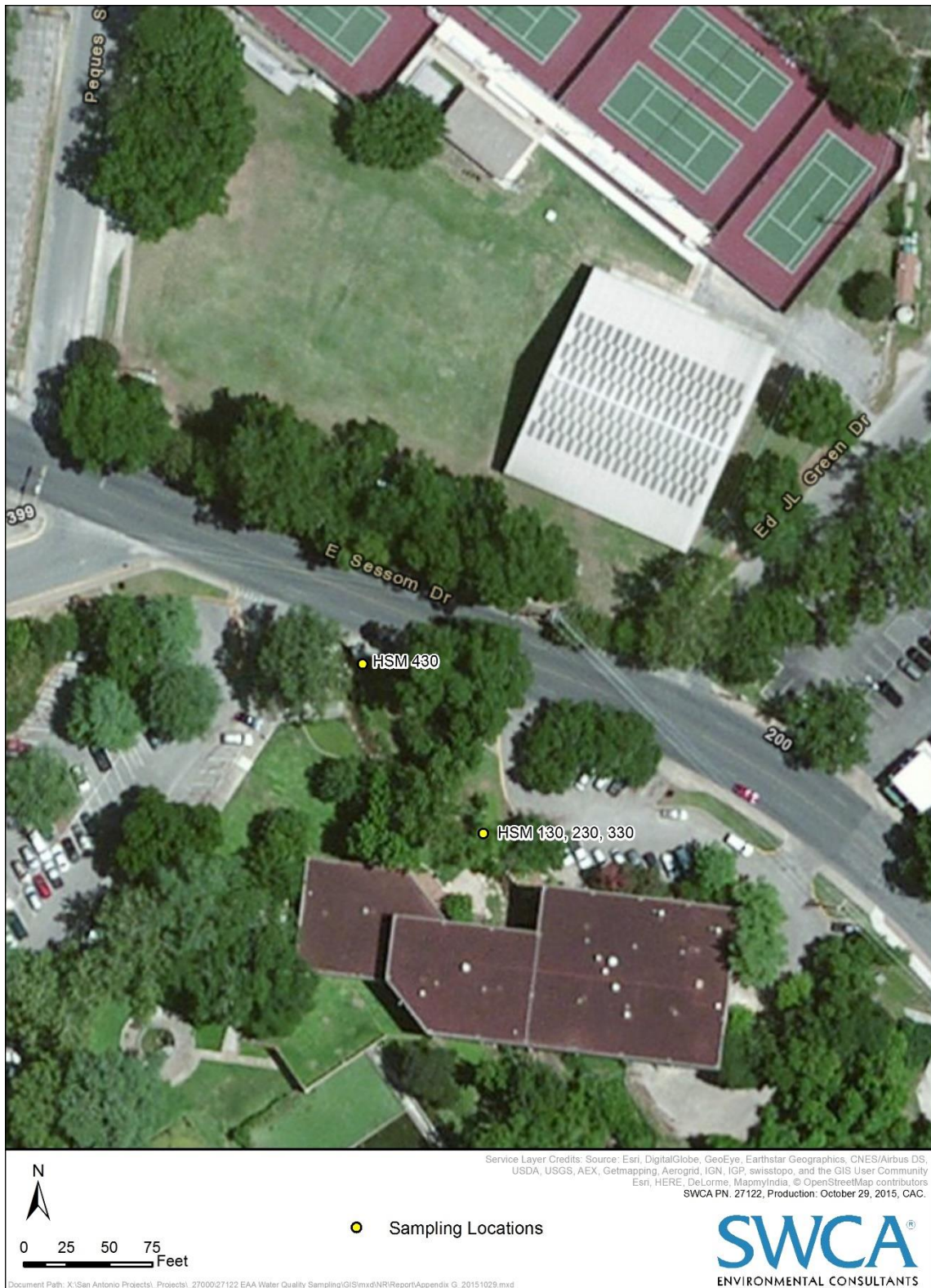
HSM320

No sediment location deviation.



HSM330

No sediment location deviation.



HSM340

Surface water samples are collected near the middle of the river channel at this site. Sediment samples were collected along the western bank and approximately 30 feet upstream of the surface collection point due to high river depth in this location. The samples were collected in a vegetated area approximately 10 feet from the western bank.



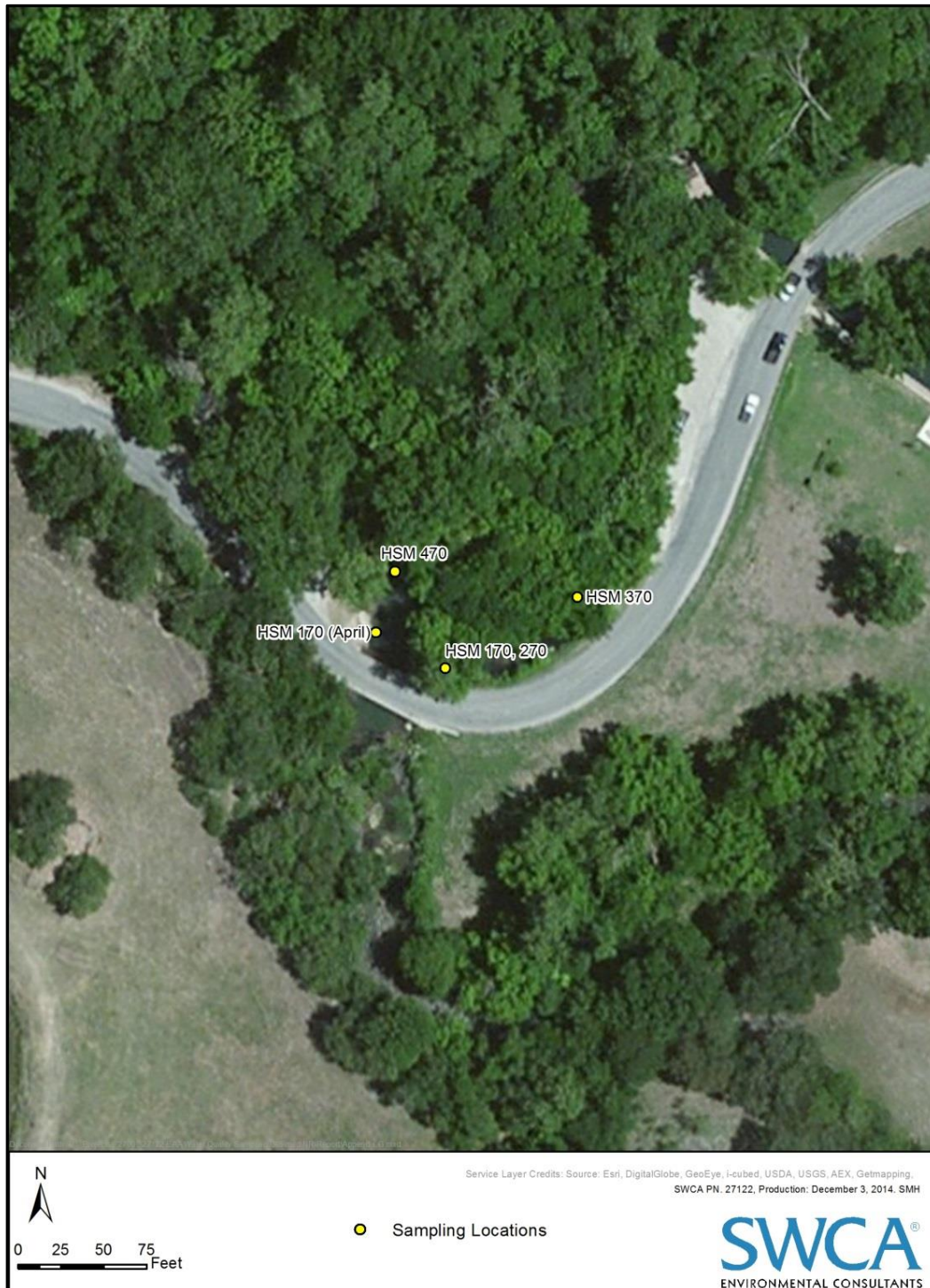
HSM350

Water depth and access along the Rio Vista Park footbridge limited the possibilities for sediment collection. Sediment samples were collected on the east side of the island approximately 15 feet to the east of the island.



HSM370

Surface water samples are collected near the middle of the channel on the upstream side of the Cape Street Bridge. High water depth, strong current, and rocky bottom limited sediment collection possibilities. Sediment samples were collected approximately 10 feet from the western bank of the Cape Street Bridge. A field duplicate was collected at the same location.



Passive Diffusion Sampling

Passive diffusion samplers were deployed at each of the surface water sample collection sites. When at all possible, deployment locations coincided with the surface water locations, some adjustments had to be made to account for river depth, accessibility by SWCA staff for installation and retrieval, and interference by the public. PDSs were deployed for two-week periods during the months of February, April, June, August, and October 2015. In 2014, SWCA staff designed and constructed a concrete and stainless steel deployment device to hold the PDS. Use of the devices continued throughout 2015.

Storm events did occur during some PDS deployment periods which may affect the analytes present in PDS results. Rain events large enough to cause a change in discharge at the USGS Streamflow gauge in the San Marcos River occurred during PDS deployments on April 10, 13, 14, 16, 17 and 18, and June 14–15. Heavy rains and flooding occurred the week prior to the June 2015 deployment, discharge levels were higher throughout the deployment period than 2015 levels prior to June.

Any alterations to sample locations or lost PDS are discussed below.

HSM420

In 2014, the PDS location for HSM420 was moved downstream from the surface water collection site to an area with easier and safer access for SWCA staff. This location was used for all 2015 deployments.

HSM430

In 2014, the PDS location for HSM430 was moved upstream from the surface water collection site. This area was chosen because a children's education program accesses the river near the main sampling site. The upstream location protects the PDS from interference by the children and any sediment disturbed by their activity. This location was utilized throughout 2015.

APPENDIX D

GROUNDWATER QUALITY MONITORING PLAN

Groundwater Quality Monitoring Plan

EDWARDS AQUIFER AUTHORITY

900 E. Quincy Street
San Antonio, Texas, 78215

Version 1.3
Revised July 2013



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ACRONYMS AND ABBREVIATIONS

| | |
|----------|---|
| ASTM | American Society of Testing and Materials |
| bgs | below ground surface |
| COC | chain of custody |
| DQO | data quality objective |
| EAA | Edwards Aquifer Authority |
| EAHCP | Edwards Aquifer Habitat Conservation Plan |
| e-line | electronic water level measurement device |
| GW | groundwater |
| MSL | mean sea level |
| NAWQA | national water quality assessment |
| PPCP | pharmaceutical and personal care products |
| psi | pounds per square inch |
| QA | quality assurance |
| QC | quality control |
| SOP | standard operating procedure |
| TWDB | Texas Water Development Board |
| USGS | United States Geological Survey |
| U.S. EPA | United States Environmental Protection Agency |
| VOA | volatile organic analysis |
| VOC | volatile organic compound |

SECTION 1

PURPOSE AND OBJECTIVES

1.1 PURPOSE AND OBJECTIVES OF THE PLAN

Data derived from water quality sampling and analysis provide the primary indicator of the state of water quality in the Edwards Aquifer. These data are also a key component of assessing water quality changes over time. Water quality data also compose the primary source of information for our understanding and monitoring of contaminant loading and migration in the Edwards Aquifer. As such, analytical samples collected for assessing water quality must be collected under a set of standard operating procedures (SOPs), which are outlined in this plan. Included herein are sections on data quality objectives (DQOs), sampling programs, analytical methods, field procedures, and guidelines for plan review.

The purpose of this plan is to provide an SOP document ensuring that useful, consistent, and defensible water quality data are produced by implementation of appropriate procedures and methods when water quality samples are being collected and analyzed. Water quality samples are currently collected under various sampling programs at the Edwards Aquifer Authority (EAA). Data quality requirements vary by program and are discussed in Sections 2 and 3.

Section 2 of this plan provides a description of DQOs in general, as well as DQOs for this program. Section 3 provides detailed information for each of the sampling programs. Section 4 provides a listing of analytical methods used by the EAA, as well as data-flagging requirements, information for sample containers, hold times, and sample preservation. Section 5 outlines field procedures; Section 6 discusses staff training and field audits. Section 7 provides information regarding annual plan review, and Section 8 provides a list of references cited in the document. The appendices (A–G) provide maps of sample locations, a glossary of terms, instrument operation and calibration information, field forms, information on regulatory limits for various compounds, stormwater sample-collection details, and equipment-decontamination procedures.

The purpose of this plan can be achieved by implementation of the objectives listed below and discussed in detail in Sections 2–7 of the plan. Each EAA staff member charged with the responsibility of collecting water quality or other analytical samples is required to be familiar with this plan, along with the objectives and procedures outlined in it. The objectives of this plan are to

- Obtain quality data that are defensible for their intended purpose,

- Analyze field samples in an appropriate and consistent manner such that the results are accurate and repeatable (see calibration procedures in Appendix C),
- Collect samples for laboratory analysis in an appropriate and consistent manner that will ensure accurate and reliable analytical results with a minimal number of anomalous data,
- Select sample sites and time periods that will provide representative water quality data for a range of aquifer conditions, and
- Review the plan annually and revise as needed.

SECTION 2

DATA QUALITY OBJECTIVES

The U.S. Environmental Protection Agency (U.S. EPA) has developed criteria for data quality objectives utilizing a seven-step process that optimizes sample collection and analysis on the basis of data uses, fiscal budget, sample quantity, and other parameters (U.S. EPA, 2000). The process is iterative and may be modified by the planning team to incorporate changes as required:

- 1. State the Problem**
Define the problem, identify the planning team, and examine the budget and schedule.
- 2. Identify the Decision**
State the decision, identify study questions, and define alternative actions.
- 3. Identify Inputs to the Decision**
Identify information needed for the decision, such as information sources, bases for action level, and sampling and analysis methods.
- 4. Define the Boundaries of Study**
Specify sample characteristics, and define spatial/temporal limits and units of decision making.
- 5. Develop a Decision Rule**
Define parameters for decision rules, specify action levels, and develop logic for action.
- 6. Specify Tolerable Limits on Decision Errors**
Set acceptable limits for decision errors relative to consequences (health effects, costs, other impacts).
- 7. Optimize the Design for Obtaining Data**
Select a resource-effective sampling and analysis plan that meets performance criteria.

2.1 U.S. EPA DQO Process as Applied to EAA Analytical Programs

2.1.1 DQO—State the Problem

Collect and analyze groundwater, spring water, and surface water samples that are contained in, issue from, or provide recharge to the Edwards Aquifer. In addition, collect stormwater and sediment samples as needed to satisfy program requirements. Sampling activities are to be conducted such that sufficient funding is held in reserve to collect confirmation samples if needed. In addition, the program must be flexible enough to collect samples in the event of a contingency (spill or other event) that affects or could potentially affect water quality of the Edwards Aquifer. The planning team includes the

Chief Technical Officer (CTO) and supervisory staff of the Aquifer Science Team of the EAA. Budget is proposed by the team and presented for board approval annually. The schedule is annual, with a general goal of collecting a minimum of 80 samples from wells, sampling all major springs (monthly or quarterly, depending on hydrologic conditions), and sampling surface waters twice annually while maintaining a budget reserve sufficient to address other needs (confirmation and contingency sampling).

Under a separate budget, the same team is charged with collecting surface water, stormwater, and sediment samples in support of the Edwards Aquifer Habitat Conservation Plan (EAHCP). Surface water, stormwater, and sediment samples are collected upstream, within, and downstream of Comal and San Marcos springs. Comal Springs has five designated sample locations, whereas San Marcos Springs has seven. Surface water and stormwater samples are to be collected twice annually, whereas sediment samples are collected once annually for the first year (to obtain baseline sediment quality information). Subsequent years may vary depending on results. See *Water Quality Monitoring Program Strategy for Comal Springs and San Marcos Springs in Support of the Edwards Aquifer Habitat Conservation Plan* (EAHCP Workplan).

2.1.2 DQO—Identify the Decision

The decision is to collect the samples as described earlier under the sampling programs and protocols outlined in detail in this document. Study questions are:

- Can the quality of water entering into, residing in, and issuing forth from the Edwards Aquifer be representatively monitored?
- For the allowed budget, how many analytical parameters can be collected?
- What analytical parameters are the most informative with regard to water quality?
- Can a relevant data set that provides historical and current water quality information as relates to the Edwards Aquifer, be developed and maintained?
- Can the data indicate trends in water quality over time?
- Can contingency sampling functionally define contaminant flowpaths and ultimately help in the prevention of public exposure to contaminants in the event of a spill?
- How does the EAA functionally share the information collected with stakeholders and the public?

Alternative actions are to

- Modify the analytical parameter list to accommodate budget constraints,
- Reduce the number of sample points and sample frequency if needed to accommodate budget constraints, and

- Continually review results to assess the need for, and feasibility of, modifying the parameter list such that analytical parameters collected provide the most information for the program, as well as cost-effective information.

2.1.3 DQO—Identify Inputs to the Decision

Sample frequency, sample type, and analytical program are all based on many inputs. The EAA strategic plan dictates minimum sample numbers, for example. Other inputs of importance include findings from karst researchers worldwide regarding the varying nuances of sampling in karst environments (i.e., multiple samples from a single location are generally more valuable than single samples from multiple locations). Assimilating and incorporating information gleaned from EAA sample results annually provide significant inputs to the process as well.

Action levels as defined for this study are not directly comparable to action levels for hazardous waste cleanup. In this program, action levels generally depend on sample type and program: for example, stormwater samples are triggered by specific stormwater events. Action levels may also be related to contingencies. If a contaminant of concern is detected in relation to a contingency, then additional sampling may be triggered. In other cases, an action level may be reached if an anthropogenic compound is detected above a regulatory limit. The resulting action will generally be to utilize additional sampling so as to delineate a possible source if a “contaminant” is the trigger.

Sampling and analysis methods are specific to each sampling program and are designed to provide data on water quality and changes to water quality that may occur over time. Results of each program are reviewed regularly, and changes to the parameters for each program may be made on the basis of these reviews or other needs. All programs are generally analyzed for field parameters (conductivity, dissolved oxygen [DO], turbidity, pH, and temperature) at the time the sample is collected. Other laboratory analytical parameters are then designated on the basis of the program.

2.1.4 DQO—Define Boundaries of the Study

Spatially the study is limited to the Edwards Aquifer Region, which includes contributing area, recharge zone, and artesian zone of the aquifer, as well as contiguous areas that may be pertinent to data collection. Temporal limits are defined by sample program and hydrologic condition. Temporal parameters are described in more detail under sample programs.

2.1.5 DQO—Develop a Decision Rule

Decision rules are defined by multiple factors:

- Strategic plan,
- Board directives,
- Approved budget,

- Data analyses and results,
- Historical data for a particular site, and
- EAHCP requirements.

2.1.6 DQO—Specify Tolerable Limits on Decision Errors

Decision-error limits are dictated by sample program. Whereas all results are considered important, contingency samples have an elevated priority because of the potential to provide a warning to the public in the event water quality is impacted. As such, in the event of a major contingency that requires long-term sampling and analysis, the budget impact would be significant. In some scenarios, additional laboratory funding would be requested from the board to cover these costs. Other sample programs are expected to be well planned and orchestrated such that no budget overruns occur.

The goal of the program in general is to collect a number of samples adequate to monitor the health of the Edwards Aquifer with high confidence that results are representative and accurate. These samples are collected through various sampling programs, as outlined in the next section.

2.1.7 DQO—Optimize the Design for Obtaining Data

The sampling plan as designed provides a resource-effective plan that meets performance criteria through data review, data assessment, and program requirements. The design is optimized by the data needs of each sample program, in which analytical parameters are specific to a program and designed to provide a maximum number of data cost-effectively.

2.2 Additional Inputs for DQO Process

Another definition of DQOs is provided by the Air Force Center for Environmental Excellence (AFCEE) in its *Quality Assurance Project Plan* (QAPP), which states that “DQOs specify the data type, quality, quantity, and uses needed to make decisions and are the basis for designing data collection activities” (AFCEE, 2001). The U.S. EPA and the AFCEE both generally utilize DQOs for hazardous waste clean-up sites, which often represent a threat to public health and the environment. However, sampling programs at the EAA differ in that most samples taken are “clean” and are not used to assess the success of a clean-up action.

Therefore, for the purposes of this plan, DQOs are met by assigning a level of precision and procedural techniques and parameter suites that are appropriate for the sample type and monitoring program. Whereas it is the purpose of this plan for all data produced to be representative and fully defensible, all data do not necessarily need to be analyzed by reference methods in the analytical laboratory utilizing a full suite of QA/QC samples. Most water quality samples collected are intended for monitoring the general status of water quality within the Edwards Aquifer, with one potential exception. In some cases, contingency sampling may be used to assess the impact of an event (i.e., a spill) to the Edwards Aquifer that has the potential for public health implications.

Therefore, DQOs developed for this document are designed to provide data of quality and quantity adequate to reflect the needs of the sample program under which a particular sample is collected. Most analytical data collected are designed to assess

- The presence or absence of anthropogenic compounds in the sample.
- Changes to chemical quality of the sample point when compared with prior data,
- Development of data adequate to establish a record of water quality such that future changes to water quality can be measured,
- Measurement of changes to water quality against changes in hydrologic conditions, and
- In the case of confirmation samples, assessment with a high degree of confidence the presence or absence of a compound of interest.

SECTION 3

SAMPLING PROGRAMS AND OBJECTIVES

Water quality samples are collected under one of the EAA sample programs described in detail in this section. Sample parameters vary with the sample program. For a better understanding of the sampling programs and sample distribution, typical water quality sample locations, see Appendix A, which is a listing of sample type and program. EAHCP sample locations are also provided.

3.1 SAMPLE TYPES AND SAMPLE PROGRAMS

Sample type is simply defined by source and media. The EAA collects samples from wells, springs, surface water, and, at times, groundwater in caves. Samples of soil or sediment may also be collected under some circumstances. As such, sample types are:

- Wells (applies to groundwater samples and includes water collected in caves),
- Springs,
- Surface water,
- Soil or sediment, and
- Stormwater.

Sample programs exist for each sample type, driving the DQO process for a given sample. Each sample program has a defined sample frequency and analytical parameter list. However, the analytical parameter list is always subject to future revision to accommodate changing circumstances. Table 3-1 summarizes current sample types and individual sample programs conducted by the EAA.

Table 3-1. Sample Types and Sample Programs

| Sample Type | Sample Program | Sample Frequency | Analytical Parameters |
|---------------|----------------|---|--|
| Wells | Passive | Quarterly | FP, GWQP, VOC, TPH, TOC, PAH, metals, bacteria |
| | NAWQA | Annually | FP, GWQP, VOC, 8081B, 8141A, 8151A, 8082A, TOC, PAH, metals, bacteria |
| | Routine | Annually | FP, GWQP, VOC, 8081B, 8141A, 8151A, TOC, PAH, metals, bacteria |
| | TWDB | Annually | FP, GWQP, VOC, 8081B, 8141A, 8151A, TOC, PAH, metals, bacteria |
| | PPCP | Annual | FP, PPCP (limited to nine wells annually) |
| | Contingency | As needed | Defined by contingency event |
| | Confirmation | As needed | Defined by detection needing confirmation |
| | QA/QC | Per QA needs | Defined by QA program |
| | EAHCP | Water level dependant | FP, GWQP, TOC, TDS |
| Springs | Primary | Quarterly (noncritical period) Monthly (critical period) | FP, GWQP, SVOC, VOC, 8081B, 8141A, 8151A, 8082A, TOC, metals, total phosphorous, bacteria, orthophosphate as P |
| | Secondary | Annually | FP, GWQP, SVOC, VOC, 8081B, 8141A, 8151A, 8082A, TOC, metals, total phosphorous, bacteria |
| | PPCP | Annually | FP, PPCP (limited to six spring samples annually) |
| | Contingency | As needed | Defined by contingency event |
| | Confirmation | As needed | Defined by detection needing confirmation |
| | QA/QC | Per QA needs | Defined by QA program |
| Surface water | Primary | Twice annually | FP, GWQP, 8081B, 8141A, 8151A, 8082A, TOC, PAH, metals, total phosphorous, bacteria |
| | Secondary | Annual | FP, GWQP, 8081B, 8141A, 8151A, 8082A, TOC, PAH, metals, total phosphorous, bacteria |
| | EAHCP | Twice annually | FP, GWQP, VOC, SVOC, 8081B, 8141A, 8151A, 8082A, TOC, metals, total phosphorous, bacteria, TKN, DOC |
| | PPCP | Annually | FP, PPCP (limited to two surface water samples annually) |
| | Contingency | As needed | Defined by contingency event |
| | Confirmation | As needed | Defined by detection requiring confirmation |
| | QA/QC | Per QA needs | Defined by QA program |
| Soil/sediment | EAHCP | Annually | FP, GWQP, VOC, SVOC, 8081B, 8141A, 8151A, 8082A, TOC, metals, total phosphorous |
| | Contingency | As needed | Defined by contingency event |
| | Confirmation | As needed | Defined by detection requiring confirmation |
| | QA/QC | Per QA needs | Defined by QA program |
| Stormwater | EAHCP | Twice annually | FP, GWQP, VOC, SVOC, 8081B, 8141A, 8151A, 8082A, TOC, metals, total phosphorous, bacteria, TKN |
| | Confirmation | As needed | Defined by detection requiring confirmation |
| | QA/QC | Per QA needs | Defined by QA program |

FP=field parameter, GWQP=general water quality parameters, SVOC=semivolatile organic compound, VOC=volatile organic compound, TOC=total organic carbon, TKN=total kjeldahl nitrogen, PPCP=personal care and pharmaceutical products., PAH=polynuclear aromatic hydrocarbons, TPH=total petroleum hydrocarbons, DOC=dissolved organic compounds

3.2 SAMPLE PROGRAM DETAIL

The sample types and programs summarized in Table 3-1 comprise the various analytical samples collected and analyzed by the EAA. Specific details of each program are provided in this section.

Sample Programs for Well Sample Types

1. Passive Sampling Program

The passive sampling program is a program to provide continuous monitoring of particular wells (referred to as sentinel wells) through the use of a passive sampling device. The device currently used is the Amplified Geochemical Imaging (AGI), LLC passive diffuse sample module (aka, Gore Module). This device utilizes a sorbent material encased in GoreTex® fabric that is capable of detecting certain analytes for volatile and semivolatile compounds, as well as petroleum hydrocarbon compounds. The Gore Modules are hung at specific intervals continuously in a sentinel well and replaced each month. The module is then shipped to AGI, LLC. for analysis (which is included as part of the module cost). Currently six wells designated as sentinel wells are located in Medina, Bexar, and Hays counties. These wells are sampled via grab sample quarterly. Sample parameter selection for this sample type is generally based on collecting parameters that are also detectable by the Gore Module, plus some additional parameters of value to an understanding of long-term trends in water quality. Sample frequency is also selected to detect temporal changes in water quality at a single sample point.

2. National Water Quality Assessment (NAWQA) Program

The NAWQA wells are a series of thirty wells installed by the USGS for long-term assessment of water quality on a regional and national scale. Ten of these wells (all in the recharge zone of Bexar County) are sampled annually. The sample parameter list is selected on the basis of the NAWQA program and is used to contribute data to that study, as well as to build a historical record of water quality for the EAA data set. Ten out of 30 NAWQA wells are sampled annually, and every well must be sampled within a three-year period.

3. Routine Water Quality Monitoring

Routine water quality samples are collected from a variety of well types (monitoring, domestic, agricultural, industrial, and municipal) to provide a data set for water quality regionwide for different well types. Sample parameters are broad in spectrum and designed to detect the most common anthropogenic compounds, as well as to document changes in concentrations of common cations and anions. These wells are generally sampled annually or less frequently.

4. Texas Water Development Board (TWDB)

Twenty TWDB samples are collected at designated wells using a split-sample technique, such that a sample set is sent to the TWDB contract laboratory (at no cost to the EAA). The remaining sample is sent to the EAA contract laboratory and analyzed for some of the same (TWDB) parameters, as well as additional parameters. This sample type provides a cost-effective tool for evaluation and comparison of analytical results for certain parameters (metals and anions). These wells (or springs, in some cases) are sampled annually under this program for a wide variety of parameters and are also used to assess the health of the system and to establish potential changes or trends in quality.

5. Pharmaceuticals and Personal Care Products (PPCPs)

These parameters detect various compounds found in common personal care products, as well as medications and household items. The primary value in this sample group is the conclusiveness of the data. Because the detection limits are low and the percentage of detections (at low concentrations) to date is high, this sample program appears to provide the most conclusive evidence of anthropogenic impacts on the Edwards Aquifer. The current sample budget allows for nine wells, six springs, and two surface waters to be sampled annually for these parameters. The same locations are sampled each year (with some exceptions) to provide a temporal record of water quality changes associated with the compounds. This program is being evaluated for an increase in sample frequency at some locations.

6. Contingency Samples

Contingency samples are collected only on an as-needed basis to assess potential contamination events related to spills or similar contingencies that have a high potential for affecting water quality in the Edwards Aquifer. Sample parameters and sample frequency are determined on the basis of type of spill (or other contingency), as well as the size of the event. Sample parameters and frequency are decided on by management. EAA staff members are subsequently directed to an appropriate course of action on the basis of assessment of the event by management.

7. Confirmation Samples

Confirmation samples are samples collected in response to an unexpected detection at a site where additional confirmation is needed in order to assess the probability that detection is not a sampling artifact or otherwise false detection. Confirmation detections are method and analyte specific and are taken at the direction of management.

8. QA/QC Samples

QA/QC samples are discussed in detail in Section 3.3.

9. EAHCP Drought Contingency-Sampling of transect wells and Springs

Well samples collected for the EAHCP are collected only when certain springflow criteria are met—specifically, low-flow situations at Comal and San Marcos springs. For Comal Springs, when flows fall below 30 cubic feet per second (cfs), weekly monitoring at three wells is to be conducted for DO, conductivity, pH, and temperature. The next trigger at Comal Springs is 20 cfs, and weekly monitoring is conducted using the same parameters plus nutrients, TDS, and TOC. For San Marcos Springs, the first trigger is 50 cfs, and the second trigger is 30 cfs.

Sample Programs for Spring Sample Types**1. Primary Springs**

Primary springs are Comal, Hueco, and San Marcos. They are sampled monthly during critical periods (critical period = a ten-day average when water levels at Bexar, County, index well J-17 of below 660 feet msl, and/or a ten-day average springflow rate at either Comal or San Marcos springs is less than 225 cfs for Comal Springs and less than 96 cfs for San Marcos Springs). During noncritical periods, sampling is generally conducted quarterly. Sample parameters are extensive because the springs represent a composite sample of aquifer water and are directly associated with habitat for threatened and endangered species.

2. Secondary Springs

Secondary springs generally produce a smaller volume of springflow and may or may not be located within the San Antonio Segment of the Edwards Aquifer. These springs are Las Moras (Fort Clark Springs), San Pedro, San Antonio, Government Canyon, and other springs that may be designated for infrequent sampling. Las Moras is generally sampled annually, whereas the others are sampled quarterly or annually if flowing. Sample parameters are the same as those for the primary springs, except that sample frequency differs between primary and secondary.

3. Pharmaceuticals and Personal Care Products (PPCPs)

These parameters detect various compounds found in common personal care products, as well as medications and household items. The primary value in this sample group is the conclusiveness of the data. Because the detection limits are low and the percent of detections (at low concentrations) to date are high, this sample program appears to provide the most conclusive evidence of anthropogenic impacts on the aquifer. The current sample budget allows for nine wells, six springs, and two surface waters to be sampled annually for these parameters. The same locations are sampled each year (with some exceptions)

to provide a temporal record of water quality changes associated with the compounds. This program is being evaluated for an increase in sample frequency at some locations.

4. Contingency Samples

Contingency samples are collected only on an as-needed basis to assess potential contamination events related to spills or similar contingencies that have a high potential for affecting water quality in the Edwards Aquifer. Sample parameters and sample frequency are determined on the basis of type of spill (or other contingency), as well as the size of the event. Sample parameters and frequency are decided on by management. EAA staff members are subsequently directed to an appropriate course of action on the basis of assessment of the event by management.

5. Confirmation Samples

Confirmation samples are samples collected in response to an unexpected detection at a site where additional confirmation is needed in order to assess the probability that detection is not a sampling artifact or otherwise false detection. Confirmation detections are method and analyte specific and are taken at the direction of management.

6. QA/QC Samples

QA/QC samples are discussed in detail in Section 3.3.

7. Drought Contingency-Sampling of transect wells and Springs

Sample Programs for Surface Water Sample Types

1. Primary Surface Water

Primary surface waters are collected twice annually from eight locations: Nueces River at Laguna, Dry Frio River at Reagan Wells, Frio River at Concan, Sabinal River near Sabinal, Seco Creek at Miller Ranch, Hondo Creek near Tarpley, Medina River at Bandera, and Blanco River at Wimberley. These sample locations have a significant historical sample record and provide information regarding the quality of waters that effectively provide recharge to the Edwards Aquifer. Sample parameter lists are fairly significant, but do not generally include VOCs because of the low probability of detection of these compounds in a surface water environment.

2. Secondary Surface Water

Secondary surface water sites may have varying locations and are generally sampled only annually. They are generally sites of interest because of their ability to provide recharge to the aquifer, or they may be indicators of water

quality from springs issuing forth from the Trinity Aquifer. Sample parameter lists are fairly significant but do not generally include VOCs because of the low probability of detection of these compounds in a surface water environment.

3. EAHCP Surface Water Samples

EAHCP surface water samples are collected at Comal and San Marcos springs; Comal Springs has five sample locations, whereas San Marcos has seven sample locations, which are situated upstream and downstream of the spring orifice locations. Parameters provide a broad spectrum of analyses so that water quality might be better understood in detail at these locations. The parameters list will also be used to study trends in water quality at these locations over time. Sample frequency is twice annually.

4. Pharmaceuticals and Personal Care Products (PPCPs)

These parameters detect various compounds found in common personal care products, as well as medications and household items. The primary value in this sample group is the conclusiveness of the data. Because the detection limits are low and the percent of detections (at low concentrations) to date high, this sample program appears to provide the most conclusive evidence of anthropogenic impacts on the aquifer. The current sampling budget allows for nine wells, six springs, and two surface waters to be sampled annually for these parameters. The same locations are sampled each year (with some exceptions) to provide a temporal record of water quality changes associated with the compounds. This program is being evaluated for an increase in sample frequency at some locations.

5. Contingency Samples

Contingency samples are collected only on an as-needed basis to assess potential contamination events related to spills or similar contingencies that have a high potential for affecting water quality in the Edwards Aquifer. Sample parameters and sample frequency are determined on the basis of type of spill (or other contingency), as well as the size of the event. Sample parameters and frequency are decided on by management. EAA staff members are subsequently directed to an appropriate course of action on the basis of assessment of the event by management.

6. Confirmation Samples

Confirmation samples are samples collected in response to an unexpected detection at a site where additional confirmation is needed in order to assess the probability that detection is not a sampling artifact or otherwise false detection. Confirmation detections are method and analyte specific and are taken at the direction of management.

7. QA/QC Samples

QA/QC samples are discussed in detail in Section 3.3.

Sample Programs for Sediment Sample Types**1. EAHCP Sediment Samples**

EAHCP sediment samples will be collected for a broad spectrum of parameters to establish a base-line data set for sediments in and around Comal and San Marcos springs. These sample data are important to an understanding of potential issues with disturbing sediments in these areas.

2. Contingency Samples

Contingency samples are collected only on an as-needed basis to assess potential contamination events related to spills or similar contingencies that have a high potential for affecting water quality in the Edwards Aquifer. Sample parameters and sample frequency are determined on the basis of type of spill (or other contingency), as well as the size of the event. Sample parameters and frequency are decided on by management. EAA staff members are subsequently directed to an appropriate course of action on the basis of assessment of the event by management.

3. Confirmation Samples

Confirmation samples are samples collected in response to an unexpected detection at a site where additional confirmation is needed in order to assess the probability that detection is not a sampling artifact or otherwise false detection. Confirmation detections are method and analyte specific and are taken at the direction of management.

4. QA/QC Samples

QA/QC samples are discussed in detail in Section 3.3

Sample Programs for Stormwater Sample Types**1. EAHCP Stormwater Samples**

EAHCP stormwater samples are collected twice annually for a broad spectrum of parameters to establish a base-line data set for stormwater quality in and around Comal and San Marcos springs. Stormwater samples are collected across the hydrograph at three points (rising, peak, and recession) to ascertain changes in water quality associated with storm flow.

2. Confirmation Samples

Confirmation samples are samples collected in response to an unexpected detection at a site where additional confirmation is needed in order to assess the probability that detection is not a sampling artifact or otherwise false detection. Confirmation detections are method and analyte specific and are taken at the direction of management.

3. QA/QC Samples

QA/QC samples are discussed in detail in Section 3.3

3.3 QUALITY CONTROL AND QUALITY ASSURANCE SAMPLES (QA/QC)

So that the data quality process is adhered to, additional samples for QA/QC must be taken and analyzed on occasion so that the quality of the sample collection and analysis process might be assessed. The various types of QA/QC samples applicable to this plan are outlined in the following paragraphs. Approximately ten percent of all samples will be QA/QC samples.

3.3.1 Matrix Spike and Matrix Spike Duplicate

Matrix spike and matrix spike duplicate samples (MS/MSD) are used to assess the effects of the sample matrix on the analytical process. The MS/MSD is a split (or replicate) of a parent sample collected in the field concurrently during the normal sample-collection process. Ideally, one MS/MSD is collected for each media type (soil, water, sludge, etc.) every 20 samples for each analysis being performed. For most sampling, no media changes will be encountered; i.e., most samples will be water. However, should the samples vary significantly in turbidity, collection of a specific MS/MSD for a sample with elevated turbidity may be advisable.

The MS/MSD is spiked and analyzed, and if the spiked analytes are recovered within a method-specific percentage, then matrix effects will be deemed minimal and no matrix data flag will be attached to the results. However, if spike recovery does not fall within the designated percentage, then analytical results will be flagged with an M-flag, indicating that a matrix effect is present. The sample name for MS/MSDs is identical to that of the parent sample, with the MS/MSD attached as a modifier at the end of the sample name. The MS/MSD will also be noted on the chain of custody (COC).

3.3.2 Ambient Blanks

Ambient blanks are taken to assess the possibility of site-specific atmospheric contamination of VOC samples. Ambient blanks are taken only when an area is suspected of having detectable quantities of atmospheric VOCs present (e.g., if VOC samples are being collected near a fueling operation). Ambient blanks are prepared by pouring ASTM

II, reagent-grade water directly into a 40-milliliter (mL), VOA container at the sample site during collection. The VOA is allowed to remain open and exposed to the atmosphere for the duration of the sample-collection process. The water is treated and analyzed as a sample from this point forward, with the designation *AB* on the COC. Ambient blanks are applicable to VOC samples.

3.3.3 Equipment Blanks

Equipment blanks consist of ASTM II, reagent-grade water poured over/through any sampling equipment used for collection of definitive samples. Most sample-collection equipment is disposable; however, in some cases, an equipment blank may be required. Equipment blanks are used to assess the effectiveness of decontamination procedures (for new materials provided to the EAA or from EAA decontamination processes) and are designated as *EB* on the COC. The frequency of collection of equipment blanks will depend on the sampling routine and sampling equipment in use.

3.3.4 Trip Blanks

Trip blanks are applicable only to VOC samples and are prepared and supplied by the contracted analytical laboratory. Trip blanks are to be shipped from the laboratory and maintained along with the VOC samples collected in the field. The purpose of trip blanks is to assess any potential contamination that may be introduced during shipping and sample handling. Trip blanks are designated on the COC as *TB*. Trip blanks are not to be opened in the field.

3.3.5 Duplicate or Replicate Samples

Duplicate and replicate samples are intended to assess the precision or repeatability of the analytical process. Typically one in ten samples should have a duplicate sample collected. The collection frequency of one duplicate per ten samples is generally acceptable. Note, however, that if a confirmation sampling event involves only three wells, then the duplicate (as well as other) QA/QC samples are still required. In other words, duplicates compose 10% of the sample set such that a sample population of ten would contain one duplicate. However, a sample population of 11 would contain two duplicates. The calculated number of duplicates is always rounded to the next whole number. Duplicates will generally be collected only at the 10% level for EAHCP analysis. For other programs, duplicate analysis is covered generally by the application of a TWDB sample set. Exceptions may apply and will be designated by management.

A duplicate sample is a second sample collected at the same location as that of the parent, either simultaneously or immediately following collection of the first sample (AFCEE, 2001). Both samples are collected, stored, and transported identically. A replicate sample, sometimes called a *split sample* is defined as a single sample divided into two samples (AFCEE, 2001). As with a duplicate, collection, storage, and transport of the resulting

samples must be identical. Duplicate and replicate samples each have unique identifiers (see Section 4).

3.3.6 Spike Samples

Spike samples are used as part of EAA's quality control on the contracted laboratory. EAA sampling staff members collect and subsequently spike twelve liters of water at one of the major springs, the spike containing a known percentage of a substance (contaminant). The spiked sample is then submitted to the contracted laboratory for analysis. If the contracted laboratory reports the findings within the specified amount, then EAA has confidence in their data. However, if the contracted laboratory is unable to detect or report the spikes, then EAA will pursue corrective action with the help of laboratory personnel to resolve the discrepancy. The corrective-action process will be initiated by the Hydrogeology Supervisor.

3.3.7 Recording QA/QC Samples in Analytical Workbook

Samples collected for QA/QC or spiked samples are to be recorded in chronological order in the laboratory notebook. The laboratory notebook is to be kept in the EAA Camden Building in the water quality area with the calibration notebook.

SECTION 4**ANALYTICAL METHODS, SAMPLE IDENTIFICATION, AND CUSTODY PROCEDURES**

This section will discuss analytical methods applicable to the EAA sampling program, as well as provide a summary of analytical hold times, acceptable sample containers, and preservation techniques. In addition, a discussion of proper identification and sample custody procedures is provided herein.

4.1 ANALYTICAL METHODS

A variety of analytical methods are used in the various water quality and sediment sampling programs. Table 4-1 lists standard analytical reference methods that have possible application to the various programs. Recall, too, that Table 3-1 provides a current listing of analytical methods/parameters for each sample type and program.

Table 4-1. Analytical Reference Methods

| Analysis | Method |
|--|------------------------|
| VOC | SW-8260b |
| SVOC | SW-8270c |
| Chlorinated herbicides | SW-8151a |
| Organophosphorus compounds | SW-8141a |
| Nonvolatile compounds by HPLC | SW-8321 |
| Organochlorine pesticides | SW-8081b |
| Polychlorinated biphenyls (PCBs) | SW-8082a |
| PAH | SW-8310 |
| Determination of triazine pesticides | EPA-619 |
| Organonitrogen pesticides in industrial/municipal wastewater | EPA-633 |
| Oryzalin in industrial/municipal wastewater | EPA-638 |
| TPH | TX-1005 |
| Metals (except mercury) | SW-6010b or SW-6020 |
| Mercury | SW-7470A |
| Cyanide | SW-9010B |

Table 4-1. Analytical Reference Methods (continued)

| Analysis | Method |
|-----------------------------------|-------------------------|
| Alkalinity | EPA-310.1 |
| Common anions | SW-9056 |
| Sulfate (SO ₄) | EPA 300.0 |
| pH | SW-9040B |
| Total dissolved solids (TDS) | EPA 160.1 |
| Total suspended solids (TSS) | EPA 160.2 |
| Ortho-phosphate | EPA 365.3 |
| Nitrate/nitrite (both as N) | EPA 353.2 |
| Ammonia (as N) | EPA 350.3 |
| Kjeldahl (as N) | EPA 351.3 |
| Total organic carbon (TOC) | EPA 415.1 or SW-9060 |
| Sulfide | EPA 376.2 |
| Dissolved organic compound | SM 5310C- 2000 |
| E-coli most probable number (MPN) | SM9223B-2004 |
| Dissolved orthophosphate lab | EPA 365.3- 1978 |
| Ammonia as N-nondistilled | SMA4500 NH3D-1997 |
| Bromide | EPA 300.0- 1993 |
| Chloride | EPA 300.0- 1993 |
| Nitrate as N | EPA 300.0- 1993 |
| Total phosphorous | EPA 365.3- 1978 |
| Enterococci | ENTEROLERT |
| Eshcerichia coli-colilert | SM 9223B 20Ed |
| Total coliform_colilert | SM 9223B 20Ed |
| TWDB anions | EPA 300.1 |
| TWDB cations | EPA 200 |
| TWDB nitrate | EPA 353.2 |
| Anti-bacterial agents | 1694 |
| Pharmaceuticals | 1694 |
| Steroids/hormones | 1698 |

| | |
|--|-------------------------|
| SIM analysis | MS-SIM-GX/MS |
| Nonylphenols | WS-MS-0010 |
| General water quality parameters (GWQP), general chemistry—(alkalinity, bicarbonate, carbonate, Ca, Mg, Na, K, Cl, SO ₄ , F, Si, Sr, bromide, nitrate as N, pH, TDS, and TSS) | Methods listed in table |

4.2 DATA-FLAGGING CONVENTIONS

Analytical data must be qualified by the EAA-contracted analytical laboratory, which is done summarily by the addition of data flags to the data result. Table 4-2 provides a summary of the data-flagging convention used in this plan (modified from AFCEE, 2001).

Table 4-2. Data Flags

| Flag | Description |
|-------------|--|
| J | Analyte positively identified. Quantitation is an estimation because the associated numerical value is below the reporting limit (RL). |
| U or ND | Analyte analyzed for, but not detected. Associated numerical value at or below method detection limit (MDL). |
| R | Data rejected because of deficiencies in ability to analyze sample and meet QC criteria. |
| B | Analyte found in associated blank, as well as in sample. |
| M | Matrix effect present. |
| T | Tentatively identified compound (using GC/MS). |
| No flag | Analyte detected at reported concentration. |

4.3 SAMPLE CONTAINERS AND HOLD TIMES

Samples sent to the analytical laboratory must be properly containerized, preserved, and analyzed within specified hold times for the method for the data to be of defensible quality. In addition to the requirement for samples to be chilled to 4°C, ±2°, some analytical methods require the sample to be maintained at specific pH values. As such, Table 4-3 lists acceptable container types, preservatives, and hold times for common analytical methods. The table includes all scheduled analyses for the various sampling programs. In the event an analysis is required that is not included in the table, Aquifer Science Team members listed herein (hydrogeology supervisor or hydrologic data coordinator) will communicate with the EAA contracted laboratory regarding appropriate containers, preservatives, and hold times for the methods in question.

Table 4-3. Sample Containers, Preservatives, and Hold Times

| Analyte or Method ¹ | Container | Preservation | Minimum Sample Volume | Holding Times |
|--|----------------------------|---------------------|---|--|
| Volatile organic compounds (SW8260B) | G, Teflon®-lined septum, T | 4°C, HCl to pH <two | 3× 40 mL with no head space or (1) 250 mL amber bottle with no head space | 14 days (water and soil); seven days if unpreserved by acid |
| Semivolatile organic compounds (SW8270C) | G, Teflon®-lined cap, T | 4°C | 1L or 8 ounces/soil | Seven days until extraction and 40 days after extraction (water); 14 days until extraction and 40 days after extraction (soil) |
| Chlorinated herbicides (SW8151a) | G, Teflon®-lined cap, T | 4°C | 1L or 8 ounces/soil | Seven days until extraction and 40 days after extraction (water); 14 days until extraction and 40 days after extraction (soil) |
| Organophosphorus compounds (SW8141A) | G, Teflon®-lined cap, T | 4°C | 1L or 8 ounces/soil | Seven days until extraction and 40 days after extraction (water); 14 days until extraction and 40 days after extraction (soil) |
| Organochlorine pesticides (SW8081) | G, Teflon®-lined cap, T | 4°C | 1L or 8 ounces/soil | Seven days until extraction and 40 days after extraction (water); 14 days until extraction and 40 days after extraction (soil) |
| Polychlorinated biphenyls (SW8082) | G, Teflon®-lined cap, T | 4°C | 1L or 8 ounces/soil | Seven days until extraction and 40 days after extraction (water); 14 days until extraction and 40 days after extraction (soil) |

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| Analyte or Method ¹ | Container | Preservation | Minimum Sample Volume | Holding Times |
|--|----------------------------|--|---|---|
| Polynuclear aromatic hydrocarbons (SW8310) | G, Teflon®-lined cap, T | 4°C | 1L or 8 ounces/soil | Seven days until extraction and 40 days after extraction (water); 14 days until extraction and 40 days after extraction (soil) |
| Total petroleum hydrocarbons (TX1005) | G, Teflon®-lined septum, T | 4°C, HCl to pH <2 | 3× 40 mL with no head space or (1) 250 mL amber bottle with no head space | 14 days (water); to extraction, and 14 days after extraction |
| General water quality parameters (alkalinity, bicarbonate, carbonate, Ca, Mg, Na, K, Cl, SO ₄ , F, Si, Sr, bromide, nitrate (as N), pH, TDS, and TSS) | P, G | 4°C | 250 mL | 28 days |
| Cyanide | P, B | 4°C; NaOH to pH >12 | 500 mL or four ounces /soil | 14 days (water and soil) |
| Ortho-phosphate (as P) | P, G | 4°C | 50 mL | 48 days |
| Nitrate (as N) and nitrite (as N) | P, G | 4°C | 250 mL | 48 days |
| Ammonia (as N) | P, G | 4°C | 250 mL | 28 days |
| Kjeldahl (as N) | P, G | 4°C | 250 mL | 28 days |
| Total organic carbon | P, G | 4°C, H ₂ SO ₄ to pH <2 | 250 mL | 28 days |
| Dissolved organic carbon | P, G | 4°C, H ₂ SO ₄ | 400 mL | 28 days |
| Phosphorus | P, G | 4°C, H ₂ SO ₄ | 500 mL | 28 days |
| Alkalinity E310.1 | P, G | 4°C | 50 mL | 14 days |
| Common anions SW9056 | P, G | None required | 50 mL | 28 days for Br ⁻ , F ⁻ , Cl ⁻ , and SO ₄ ²⁻ ; 48 hours for NO ₃ ⁻ , NO ₂ ⁻ , and PO ₄ ³⁻ |
| Cyanide, total and amenable to chlorination SW9010A SW9012 | P, G, T | 4°C; NaOH to pH >12, 0.6 g ascorbic acid | 500 mL or four ounces /soil | 14 days (water and soil) |
| Total dissolved solids (TDS) E160.1 | P, G | 4°C | 100 mL | Seven days |
| Total suspended solids (TSS) E160.2 | P, G | 4°C | 100 mL | Seven days |
| Biological oxygen demand (BOD), five-day | P, G | 4°C | 1L | 48 hours |
| Sulfide | P, G | 4°C | 1L | Seven days |

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| Analyte or Method ¹ | Container | Preservation | Minimum Sample Volume | Holding Times |
|---|-------------------------|--|-----------------------|--|
| Total inorganic carbon | P, G | 4°C | 250 mL | 28 days |
| Escherichia coli-colilert | P, G, WP | 4°C, dark, sodium thiosulfate, one-inch headspace | 100–250 mL | Six + two h (this holding time represents six field hours and two lab hours) |
| Enterococci | P, G, WP | 4°C, dark, sodium thiosulfate, one-inch headspace | 100–250 mL | Six + two h (this holding time represents six field hours and two lab hours) |
| Total coliform-colilert | P, G, WP | 4°C, dark, sodium thiosulfate, one-inch headspace | 100–250 mL | Six + two h (this holding time represents six field hours and two lab hours) |
| TWDB anions | P, G | 4°C, filtered on site | 500 mL | 28 days |
| TWDB cations | P, G | 4°C, HNO ₃ , filtered on site | 250 mL | 28 days |
| TWDB nitrate | P, G | 4°C, H ₂ SO ₄ , filtered on site | 500 mL | 28 days |
| 1694 Pharmaceuticals (LCMS/MS) Acetaminophen Caffeine Carbamazepine Cotinine DEET Diltiazem Fluoxetine Gemfibrozil Ibuprofen Lincomycin Naproxen Sulfamethoxazole Trimethoprim Tylosin Iopromide | G, Teflon®-lined cap, T | 4°C | 1L or 8 ounces/soil | Seven days (unpreserved), 14 (days preserved) |
| 1694 Antibacterial (LCMS/MS) Triclobarban Triclosan | G, Teflon®-lined cap, T | 4°C | 1L or 8 ounces/soil | Seven days (unpreserved), 14 (days preserved) |

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| Analyte or Method ¹ | Container | Preservation | Minimum Sample Volume | Holding Times |
|---|-------------------------|-------------------------------------|-------------------------|--|
| 1698 Steroids/hormones (LCMS/MS) 17a-Estradiol 17a-Ethynyl estradiol 17b-Estradiol Equilenin Estriol Estrone Progesterone Testosterone | G, Teflon®-lined cap, T | 4°C, H ₂ SO ₄ | 1L or 8 ounces/soil | Seven days (unpreserved), 14 (days preserved) |
| Nonylphenols/ethoxylates/bisphenol-A (GCMS) Bisphenol-A Nonylphenol diethoxylate (tech.) Nonylphenol monoethoxylate (tech.) p-Nonylphenol (tech.) p-tert-octylphenol para-n-nonylphenol | G, Teflon®-lined cap, T | 4°C, H ₂ SO ₄ | 1L or 8 ounces/soil | Seven days (unpreserved), 14 (days preserved) |
| Selected metals—6020 (Al, Sb, As, Ba, Be, Cd, Cr (total), Cu, Fe, Pb, Mn, Hg, Ni, Se, Ag, Tl, and Zn) | P, G, T | HNO ₃ to pH <2, 4°C | 500 mL or 8 ounces/soil | 180 days (water and soil) |
| Hg—Cold vapor 7470.7471 | P, G | HNO ₃ to pH <2, 4°C | 250 mL | 28 days (14 days if in plastic bottle) |
| Selected metals—(ICP unless otherwise noted) 6020/7470/7471 (Al, Sb-ICP-MS or GFAA, As, Ba, Be, Cd, Cr (total), Cu, Fe, Pb, Mn, Hg-ICP-MS or CVAA, Ni, Se-ICP-MS or GFAA, Ag, Tl-ICP-MS or GFAA, and Zn) | P, G, T | HNO ₃ to pH <2, 4°C | 500 mL or 8 ounces/soil | 180 days (water and soil) |
| Hg- ICP-MS or CVAA 7470/7471 | P, G | HNO ₃ to pH <2, 4°C | 250 mL | 28 days (14 days if in plastic bottle) |

- a. Polyethylene (P); glass (G); brass sleeves in sample barrel, sometimes called California brass (T).
- b. No pH adjustment for soil.
- c. Preservation with 0.008 percent Na₂S₂O₃ only required when residual chlorine present.

4.4 SAMPLE IDENTIFICATION

Each sample must have a unique identifier so that it can be differentiated from other samples. In addition, sample names must meet the required criteria for entry into the data base and subsequent electronic storage and retrieval of the data. Therefore, sample names must conform to the guidelines herein.

4.4.1 Sample Identification, for Non-EAHCP Samples

The primary method for non-EAHCP sample identification will be to use the state well registration number for wells (and springs as applicable) or the site name for surface water samples. When no well number is available for a spring, then an abbreviation for the spring name and orifice will be used. For example,

- The unique identifier, for use on the COC for Comal Springs, Orifice 1 is DX 68-23-301,
- The unique identifier for use on the COC for Comal Springs Orifice 3 (no state well number) is CS3,
- The unique identifier for use on the COC for the Nueces River at Laguna is Nueces@Laguna, and
- For wells that are sampled in more than one location within the borehole, the interval number is attached to the well name. For example, well LR-67-09-101 is regularly sampled at two intervals, so the COC name is LR-67-09-101-1 (interval 1 or upper interval) and LR-67-09-101-4 (interval 4, or the deepest interval).

Note that to the extent possible, custody forms and sample-container labels will be preprinted by the laboratory.

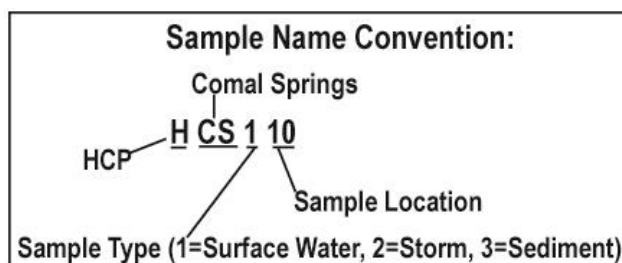
In some cases no well number or other recognized registration number will exist for the sample point. Then documentation for the sample location will require location (latitude/longitude and address if available) and name of well owner. Photographic documentation is also required. The subsequent sample name will be a *pseudo* state well number derived from the well location and owner name. For example,

The unique identifier for a sample taken from the Mary Smith residence in San Antonio, a private well with no state well registration number and located in Bexar County (abbreviation AY) at state well grid location 68-23-8, would be AY-68-23-8MS.

When wells of this type are sampled, proper documentation to include collection location, sample name, sample parameters, date, and time is extremely important and will be recorded in the field log for cross reference to the COC.

4.4.2 Sample Identification, for EAHCP Samples

For samples collected under the EAHCP, sample names are designed to provide additional data regarding sample type. Specifically the sample name will indicate the sample as an EAHCP-related sample, the spring group (Comal or San Marcos), sample type (surface water, stormwater, or sediment), and sample location. In the example below, the sample name refers to an EAHCP sample at Comal Springs, collected for surface water, at location 10. Sample locations are noted on the sample-collection maps for the EAHCP (included in Appendix A with calendar year 2013 non-EAHCP sample locations).



4.4.3 Sample Identification, QA/QC

For QA/QC samples, a modifier is added to the sample name to indicate the QA/QC type, for example, DX-68-23-301 (Comal Spring 1). If an MS/MSD sample were collected, a separate set of samples named DX-68-23-301MS/MSD would be collected. The appropriate modifier for each QA/QC sample is listed in Table 4-4.

Table 4-4. QA/QC Sample Nomenclature

| Sample Type | Modifier |
|-------------------------------------|-----------------|
| Matrix spike/matrix spike duplicate | MS/MSD* |
| Ambient blank | AB [#] |
| Equipment blank | EB [#] |
| Trip blank | TB [#] |
| Duplicate | FD* |
| Replicate | FR* |

* Requires sample, with same sample name as parent + modifier at end.

[#] Numerical suffix to be attached and referenced in laboratory notebook; suffix starts at 1 at beginning of each calendar year. Details for location, etc. included in field notebook documentation.

4.5 SAMPLE CUSTODY

All samples shipped to the analytical laboratory must have proper custody documentation. One person on each sampling team is to have primary responsibility for sample custody (generally the lead sampler). This person will be designated as the *sample custodian* for sample collection. A person has custody of a sample group if samples are (1) in his/her possession, (2) in his/her view after being in his/her possession, (3) placed in a secure area by the sample custodian.

Furthermore, the laboratory COC form is to be filled out completely by the sample custodian in the field. The form must contain all required information for proper sample identification (if not preprinted) and must contain appropriate signatures. In addition, samples must remain in control of the sample custodian. Once collected, samples must be under the supervision of the sample custodian or secured in a manner such that no reasonable chance of unauthorized access to the samples exists. Furthermore, samples shipped by a common courier (i.e., Federal Express), require that the sample custodian note on the COC when the samples were released to the courier and why. The contracted analytical laboratory will sign the COC upon receipt. A breach of sample custody can invalidate the defensibility of the sample set.

4.6 DATA VALIDATION

Analytical data require review in order to be validated prior to publication. The amount of review (or level of review) is a function of the sample type. Field-collected data results are reviewed in the field by the analyst. One of the best ways for the field analyst to assess the acceptability of field data and subsequently validate them is to compare the results with historical data. This comparison, combined with proper equipment calibration, maintenance, and analytical technique, will provide an adequate validation process for field-parameter data. In the event that the analyst finds a discrepancy in the field data, a second analysis for the parameter in question should be performed. If the analyst feels that the data may be inaccurate because of issues with the field analysis, this fact is to be noted on the sample field sheet.

Contract analytical-laboratory data will receive a 100% analyst review at the analytical laboratory prior to posting of analytical results. A subsequent analytical laboratory review by the QA/QC section is required prior to the analytical laboratory's certification of the results. A subsequent 10% review by EAA staff of the analytical data is required upon receipt of the final analytical report. The analytical report will contain numerical analytical results for the laboratory QA/QC samples (i.e., LCS, method blanks, etc.). These laboratory analytical data are to have data flags assigned by the analytical laboratory.

SECTION 5

FIELD PROCEDURES AND SAMPLE COLLECTION

Possibly the most significant part of any successful sample collection is the field procedures and documentation that occur in the field. Field procedures to include sample equipment decontamination; sample-collection procedures for well, spring, surface water, and sediment samples; a listing of potential sources of contamination; and the proper use of field notebooks are included in this section.

5.1 RESPONSIBILITIES

The CTO and hydrogeology supervisor for the data-collection program will ensure that the samples obtained represent the environment being investigated. The hydrologic data coordinator will ensure that all field crews are provided with the necessary information, equipment, and supplies to successfully schedule and complete sampling. The hydrologic data coordinator will also be the primary point of contact between the contract analytical laboratory project manager and the EAA sampling team(s). The hydrologic data coordinator will report sampling deviations to the CTO and hydrogeology supervisor. Sample-collection staff (generally, environmental science technicians) are responsible for being familiar with the instructions provided in this SOP and for collection of samples in accordance with this SOP. For most sample-collection events, a sample team of two people will be utilized. Teams will have a lead sampler (according to experience level) who is directly responsible for adherence to directives of the SOP.

5.2 EQUIPMENT DECONTAMINATION

In order to obtain samples that are reliable and defensible, all (nondisposable) sample-collection equipment must be decontaminated prior to use. When possible, sample collection from a wellhead valve directly to a sample container is best. When this kind of collection is not possible, disposable equipment is preferable.

If neither option is plausible, then nondisposable sample-collection devices (constructed of Teflon® when possible) must be used. Sampling equipment that is exposed directly to sample media (pumps, peristaltic or submersible pump tubing, reusable bailers, or other devices) will be washed in a nonphosphate, laboratory-grade detergent such as Alconox®, followed by a double rinse in potable water. A final rinse of deionized or distilled water will be applied after completion of the initial decontamination process.

Equipment that will not be used immediately must be kept clean by wrapping in aluminum foil or placed inside clean plastic bags. Such storage will prevent

contamination of the equipment prior to use. See Appendix G for additional detail regarding equipment-decontamination procedures.

5.3 SOURCES OF SAMPLE CONTAMINATION

Samples can easily become contaminated during the sample-collection process. It is the responsibility of the sampler to prevent contamination from occurring. A multitude of potential cross-contamination sources are present in the field environment. Because many of the analytical methods used can quantify various analytes in parts per billion or less, even minute sources can potentially contaminate a sample. For example, Table 5-1 summarizes some of the potential sources that can cause a false-positive reading in a sample. These should be considered when samples are collected in the field. Also note that water has a strong affinity for many anthropogenic compounds. Use of good judgment is another aspect of collecting defensible data. Steps should be taken to avoid cross-contamination of samples. If the sampler suspects the possibility of cross-contamination, he/she should note it in the field log for the sample set in question, or the site should be sampled again if necessary.

Table 5-1. Potential Sources of Cross-Contamination

| Source | Possible Contaminant |
|--|------------------------------|
| Fuels—generators, work vehicles | BTEX/TPH/VOC/SVOC |
| Exhaust fumes—generators, vehicles, heavy roadway traffic, overhead air traffic | BTEX/TPH/VOC/SVOC |
| Oil/grease residue on tools, gloves, etc. | TPH/SVOC |
| Tape | VOC |
| Insect spray | VOC/SVOC/pesticides |
| Insect repellent | SVOC/VOC/pesticides |
| Sunscreen | VOC/SVOC/PPCP |
| Soil/debris | Bacteriological/metals/SVOCs |
| Foods/drinks/medications and other personal care products such as soap, makeup, deodorant, etcetera. | PPCPs |

5.4 FIELD NOTEBOOKS

The field notebook is a legal document and should be treated as such. All pertinent site information should be in the notebook, including site name, weather information, site conditions, well condition (if applicable), equipment problems, sample-collection notes such as approximate sample times, and any other information that may be deemed valuable. The names of individuals on the sample team, as well as visitors to the site, should also be recorded in the notebook. All information recorded in the field notebook should follow the format described herein. No blank spaces are to be left on pages. All blank areas should be marked through with a single line and initialed by the author. The top of each page should have the date and sample site. The base of each page should contain the initials of the author. Mistakes are to be crossed out with a single line and initialed. Field notebooks are to be recorded in black ink only.

5.5 SAMPLE COLLECTION

Field personnel must wear clean (disposable) nitrile gloves during the sample-collection process. Generally samples for field water quality parameters are to be collected first, followed by VOC, SVOC, and metals samples. Any required information is to be recorded in the field notebook before, during, and after sampling.

5.5.1 Well Samples

Each well must be gauged and sounded (if possible). The general condition of the well will be noted in the field notebook. After the water level is gauged, the purge volume for the well will be calculated by the following equation,

$$V = H \times F,$$

where V is one well volume, H is the difference between depth of the well and depth to water in feet (i.e., length of water column in well), and F is the number of gallons per foot of water for the well size (Table 5-2).

Table 5-2. Well-Casing Volume in Gallons per Foot

| Casing Diameter (in inches) | F (gallons per foot of water in well) |
|-----------------------------|---------------------------------------|
| 2 | 0.16 |
| 4 | 0.65 |
| 6 | 1.47 |
| 8 | 2.6 |
| 10 | 4.1 |
| 12 | 5.9 |
| 16 | 10.4 |

The relationship $F = \pi (D/2)^2 \times 7.48$ gallons/ft³ can be used to calculate pipe volumes not listed in the table. Note that D = pipe diameter in feet and F = volume per foot.

A well may be sampled upon achieving *one* of the following: a minimum of three well volumes are purged from the well *or* field-parameter readings are stabilized for a minimum of three parameter measurements. Wells that go dry prior to purging the three well volumes, or the field-parameter readings have not stabilized, shall be purged to dryness (except for drinking-water supply or irrigation wells). During purging, water will be monitored for the following field parameters: *temperature, pH, DO, conductivity, and turbidity*.

Stabilization is defined as

- Temperature fluctuations limited to $\pm 1^\circ \text{C}$,
- pH fluctuations ± 0.1 unit,
- DO fluctuations ± 0.3 milligrams per liter (mg/L),
- Conductivity fluctuations $\pm 5\%$, and
- Turbidity ± 10 NTU.

In the event that these parameters do not stabilize (after purging of three well volumes), a maximum of six well volumes will be purged prior to sample collection (if the field parameters stabilize at any point, the well is considered ready to sample, and purging may cease). Once the well has stabilized or the maximum purge volume is reached, and the well has recovered to at least 80% of its initial level, it is ready to sample.

5.5.2 Spring Samples

Springwater samples should be as representative of the actual water issuing forth from the spring as possible and not be “contaminated” by surrounding surface waters. As such, various sample-collection techniques may be necessary. For spring orifices located below surface water, samplers should use a peristaltic pump to collect the springwater sample by placing the intake part of the pump tubing in the spring orifice. This placement allows for filling of sample bottles without introducing surface waters or overflowing the bottles and losing any preservatives inside. This technique is not feasible or necessary for all spring sites but should be utilized as appropriate. When a spring that can be sampled without a pump is being sampled, then a typical grab sample may be collected. In some cases (high flow volume) it may be necessary to collect samples in a clean bottle (such as a clean 1,000-mL amber glass bottle, clean Teflon beaker, or something similar) and the container used to transfer water into subsequent containers. Doing so will prevent the loss of any preservatives that may be in sample bottles. However, the action should be performed with as little agitation to the sample as possible to preserve potential VOCs in the parent sample.

Note: If preservatives in the sample container are diluted or lost because of the collection technique, a new bottle should be used. If a new bottle is unavailable, the lack of preservatives must be communicated to the laboratory to ensure that the sample remains valid by being analyzed within the appropriate hold time.

Current information and observations concerning springflow at the time of sample collection should be entered in the field notebook. For example, approximate springflow volume (can be listed as low, medium, high) is the flow representative of an extreme volume (high or low); observed water quality should be noted (clear, cloudy, or murky), along with other observations deemed appropriate by the lead sampler.

5.5.3 Surface Water Samples

Surface water samples should be collected without disturbing the sediment, if at all possible. The presence of sediment in the sample may bias the results. Samples should be collected from the flowing parts of the stream on the upstream side of the sample collector. Samples are not to be collected from stagnant areas, and they should also be taken from approximately the same location for each sample event. Sample bottles should be filled by collecting the water sample in a clean bottle or by using a peristaltic pump and transferred into the final sample bottle. Caution should be used to prevent overfilling of the sample bottle and diluting any preservatives that may be in the bottle.

Note: If preservatives in the sample container are diluted or lost because of the collection technique, a new bottle should be used. If a new bottle is unavailable, the lack of preservatives must be communicated to the laboratory to ensure that the sample remains valid by being analyzed within the appropriate hold time.

Information regarding the sample point in the stream, streamflow, and water conditions, as well as other information deemed appropriate by the sampler, should be entered into the field notebook at the time of sample collection.

5.5.4 Sediment Samples

Sediment samples are scheduled for collection by the EAHCP sampling program. Furthermore, the possibility exists that EAA staff may be required to collect samples of this type on occasion for other programs. As such, a brief discussion of this type of sample is included herein. Sediment samples may be collected from below the water line, from a dry stream bed, or from any other source in which sediments or soils may collect. The collection technique will depend on conditions. For example, a push tube for collection of sediments below the water surface is generally needed. However, if sediments are being collected from a dry area, then they may be collected using a trowel, hand auger, or push tube of some type. As with all sediment/soil-related samples, VOC samples must be collected in a manner that will minimize the loss of in situ volatiles. As

such, sediment samples for VOC analysis will not be composited or homogenized in the field. Samples for VOC analysis are to be collected first.

In the event that the discreet-interval sediment sampler is used for collection of sediments, the procedure for device operation is as follows:

1. Insert the lower-half of the lead internal rod using a $\frac{3}{8}$ -inch coupler (first stage) into the internal drive tip. Pull down on the brass ring, push the grooved end of the lead internal rod into the recess, and gently release the brass ring.
2. Insert the internal drive tip and lead internal-rod assembly into the external drive tip.
3. Connect the upper lead internal rod using the $\frac{3}{8}$ -inch coupler (second stage) to the lower lead internal rod (first stage).
4. Insert a four-ft liner, with the hole in the liner oriented to the top, into the sample tube (the sample tube has a two-inch outside diameter and consists of two parts, a double female lead section and a male \times female extension). If the EAA staff chooses to use a two-ft liner instead of a four-ft liner, the process is the same, except that the male \times female upper extension is not used.
5. Insert a plastic core catcher (white) in the bottom of the sample tube, with the dome pointing toward the top.
6. Insert the internal drive tip/external drive tip assembly into the sampler tube.
7. Insert the metal core catcher into the top of the main sampler tube, with the dome pointing upward.
8. Install the internal tip chamber to the top of the main sampler tube.
9. Install the top drive head adapter to the top of the internal tip chamber.
10. Install the thread protector cap or internal rod with external drive extensions (if using $1\frac{1}{8} \times 3$ ft external extensions with $\frac{3}{8}$ -inch internal rods, place a $\frac{3}{8}$ -inch coupler on the top of the internal rods prior to installing the top drive head adapter). Install the thread protector cap at the top of the internal rod prior to connecting the vented drive head (install the correct number of internal/external extensions necessary to lower the sampler to the surface and arrive at the desired sampling point).
11. Install the vented hammer adapter, already attached to the slide hammer.

The field notebook will note details related to the sediment samples; for example, was the sediment dry or below water, how was it collected, was it discolored, at what depth (from the surface) was the sample collected? If sediments are field screened with a photoionization detector (PID), readings from the various intervals will be recorded. Other details will be recorded as deemed appropriate by the sampler.

Also, if a hand trowel is used, it must be constructed of stainless steel, and it must be decontaminated prior to each use. For sites at which multiple samples will be collected, multiple hand trowels may be used, or a single trowel may be used if it is decontaminated in the field (Alconox wash, double rinse in potable water, followed by a DI water rinse).

5.5.5 Stormwater Samples

Stormwater samples are scheduled for collection under the EAHCP program at each spring group, twice annually. Stormwater sample collection offers additional challenges and safety issues, as compared with that of other samples collected under EAA programs. This section provides a general summary of stormwater sampling, additional detail regarding this sample type being provided in Appendix F.

Stormwater samples are scheduled for collection across three points on the storm hydrograph. One sample collected from the initial rise on the hydrograph, a second sample from the peak area of the hydrograph, and a final sample along the recession limb of the graph. In addition, water quality parameters obtained from EAA-installed real-time water quality monitors, flow data from the U.S.G.S. springs gauges, and local weather radar maps will be used to define the behavior of the systems and help guide sample-collection timing. The real-time monitors collect data at 15-minute intervals for conductivity, DO, pH, temperature, and turbidity.

A stormwater event will be dictated by a rainfall event sufficient to cause a significant rise in springflow at either Comal or San Marcos springs. The significant rise in springflow is to be further defined in conjunction with real-time data systems. See Appendix F for details on stormwater sampling procedures.

SECTION 6

ANNUAL REVIEW OF PLAN

6.1 ANNUAL REVIEW OF GROUNDWATER QUALITY PLAN

Data collection described in this plan will be reviewed by May 31 each year. The review will be directed at ensuring that all data collection herein is necessary, properly performed, and properly staffed. Furthermore, the review will ascertain whether the methodologies in use remain appropriate for their intended purpose. The review process will include all sample types and programs, as well as methods used to collect and analyze these samples.

Postreview, modifications will be made, if needed, to accommodate changes to EAA sampling. Changes will be initiated by the management and staff of the EAA Aquifer Science Team.

SECTION 7

CONTINUING EDUCATION CREDITS FOR SAMPLE-COLLECTION PERSONNEL

7.1 CONTINUING EDUCATION

Staff members assigned to sample-collection teams must attain a minimum of 12 hours of continuing education each year. Opportunities for continuing education will be provided either in-house by the EAA, or, in some cases, staff may be sent to an offsite facility to attend a class. One hour of credit is considered to be one classroom or contact hour. Staff may also carry credits over into the following year if more than 12 hours of credit are obtained in a calendar year. It is the responsibility of each staff member to document his/her credit hours annually and submit them to the hydrogeology supervisor by December 1 of each year.

SECTION 8

REFERENCES CITED

- AFCEE, 2001, Quality Assurance Project Plan, Version 3.1: Air Force Center for Environmental Excellence, Brooks AFB, Texas.
- EAA, 2012, Water Quality Monitoring Program Strategy for Comal Springs and San Marcos Springs in Support of the Edwards Aquifer Habitat Conservation Plan.
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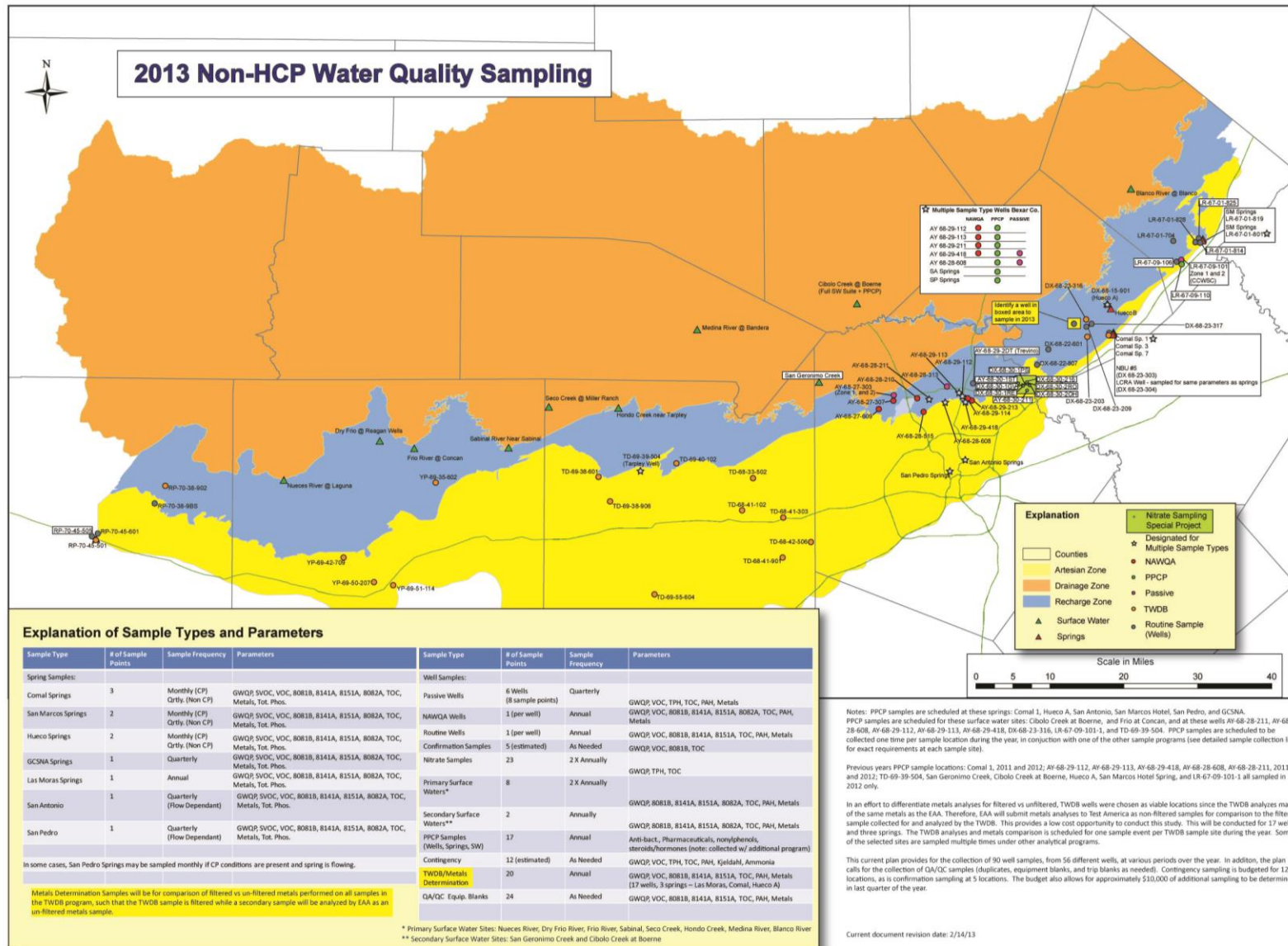
SECTION 9

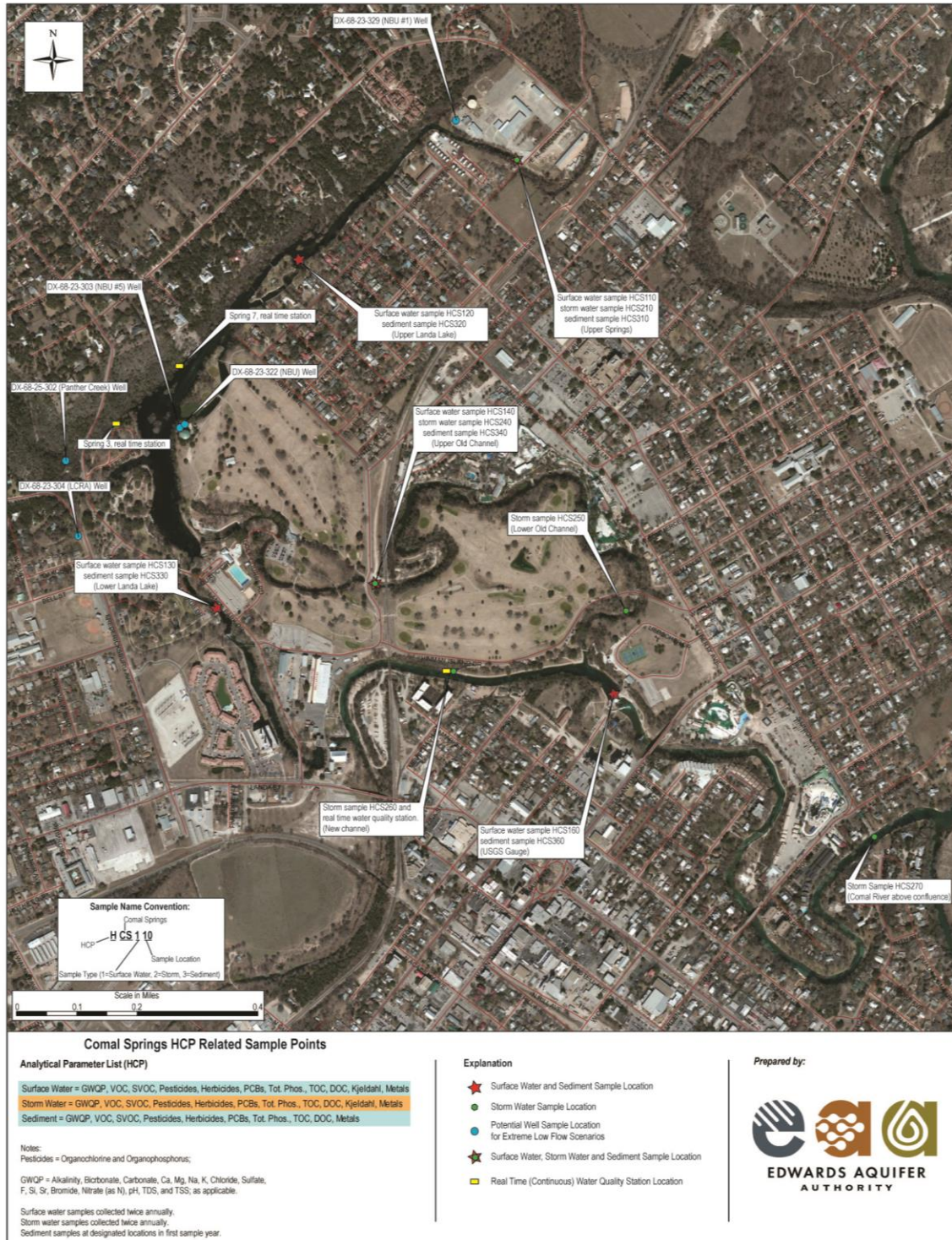
REFERENCES NOT CITED

- Driscoll, F.G., 1995, Groundwater and Wells: St. Paul, Johnson Screens, 1089 p.
- Nielsen, D.M., 2006, Environmental Site Characterization and Ground-Water Monitoring: New York, Taylor and Francis, 1318 p.

APPENDIX A—Sample Locations (2013)

Groundwater Quality Monitoring Plan Edwards Aquifer Authority







APPENDIX B—Glossary of Terms

| | |
|---------------|--|
| Ambient blank | Sample known not to contain target analytes, which are used to assess airborne contaminants at the site. The ambient blank [AB] is |
|---------------|--|

| | |
|-----------------|--|
| | opened at the site and exposed to site (ambient) conditions and subsequently treated as an environmental sample thereafter. AB samples are applicable to VOC analysis only. |
| Anion | Negatively charged ion. |
| Aquifer | Underground geological formation or group of formations containing water; source of groundwater for wells and springs. |
| Cation | Positively charged ion. |
| DOC | Abbreviation for dissolved organic carbon, a broad classification of organic molecules of varied origin and composition within aquatic systems. Organic carbon compounds are a result of decomposition processes from dead organic matter, such as plants. |
| DQO | Abbreviation for data quality objectives, a process used to develop performance and acceptance criteria or data quality objectives that clarify study objectives, define the appropriate type of data, and specify tolerable levels of data needed to support decisions. |
| Equipment blank | Sample used to assess the effectiveness of the decontamination process on sampling equipment. The equipment blank is prepared by pouring reagent-grade water over/through sampling equipment and analyzing for parameters of concern (to match the sampling routine applicable to the site). |
| Field duplicate | Second sample collected simultaneously from the same source as the parent sample, but which is submitted and analyzed as a separate sample. This sample should generally be identified such that the laboratory is unaware that it is a field duplicate. |
| Field replicate | Sometimes referred to as a <i>split sample</i> , a single sample divided into two (or more) samples. |
| Groundwater | Water found beneath Earth's surface that fills pores between materials, such as sand, soil, or gravel. |
| Initial rise | Initial surface runoff of a rainstorm. During this phase, water pollution entering storm drains in areas with high proportions of impervious surfaces is typically more concentrated during first flush than it is during the remainder of the storm. |

| | |
|----------------|---|
| Matrix spike | Sample used to determine the effect of the matrix on a method's recovery efficiency. A known amount of the target analyte is added to a specified amount of matrix sample for which an independent estimate of the target analyte concentration is available. Duplicate samples must be available as well (matrix spike duplicate, or MSD). |
| MDL | Abbreviation for method detection limit, minimum concentration of a substance that can be measured and reported with 99% confidence that the analyte concentration is greater than zero, as determined from analysis of a sample containing the analyte in a given matrix. |
| Peak | Maximum instantaneous flow at a specific location resulting from a given storm condition. |
| PQL | Abbreviation for practical quantitation limit, which is the smallest concentration of the analyte that can be reported with a specific degree of confidence. |
| Precision | State or quality of being precise; exactness. The ability of a measurement to be consistently reproduced. |
| Purge | To remove standing water in a well. |
| Recession | End of runoff event, which is defined as the point in time when the recession limb of the hydrograph is <two% of the peak or is within ten % of the prestorm base flow, whichever is greater. |
| Recharge zone | Where an aquifer is replenished with water by the downward percolation of precipitation through soil and rock. |
| Representative | Said of samples collected that are similar to those of groundwater in its in situ condition. |
| RL | Abbreviation for reporting limit [RL], the smallest concentration of an analyte reported by the laboratory to a customer. The RL is never less than the PQL and is generally twice the MDL. |
| Spike sample | One of any known concentrations of specific analytes that have been added to minimize change in the matrix of the original |

| | |
|---------------|---|
| | sample. Every spike sample analyzed should have an associated reference to the spike solution and the volume added. |
| Spring | Water coming naturally out of the ground. |
| Surface water | That which forms and remains above ground, such as lakes, ponds, rivers, streams, bays, and oceans. |
| SVOC | Abbreviation for semivolatile organic compounds, which is a group of chemicals composed primarily of carbon and hydrogen that have a tendency to evaporate (volatilize) into the air from water or soil. Some of the compounds that make up asphalt are examples of SVOCs. |
| TDS | Abbreviation for total dissolved solids, or the total amount of all inorganic and organic substances, including minerals, salts, metal, cations, or anions that are dispersed within a volume of water. |
| Temporal | Over a period of time. |
| TKN | Abbreviation for total kjeldahl nitrogen, which is the total concentration of organic and ammonia nitrogen in wastewater. |
| TOC | Abbreviation for total organic carbon, which is the gross amount of organic matter found in natural water. Suspended-particulate, colloidal, and dissolved organic matter are part of the TOC measurement. Settable solids consisting of inorganic sediments and some organic particulate are not transferred from the sample by the lab analyst and are not part of the TOC measurement. |
| Trip blank | Sample known to be free of contamination (for target analytes) that is prepared in the laboratory and treated as an environmental sample after receipt by the sampler. Trip blank [TB] samples are applicable to VOC analysis only. |
| TSS | Abbreviation for total suspended solids, which are the nonfilterable residue retained on a glass-fiber disk filter mesh measuring 1.2 micrometers after filtration of a sample of water or wastewater. |
| VOC | Abbreviation for volatile organic compounds, which are often used as solvents in industrial processes and are either known or suspected carcinogens or mutagens. The five most toxic are vinyl |

chloride, tetrachloroethylene, trichloroethylene, 1,2-dichloroethane, and carbon tetrachloride.

Well Bored, drilled, or driven shaft whose purpose is to reach underground water supplies.

APPENDIX C—Equipment Use and Calibration

DOCUMENTATION PROCEDURES

All equipment maintenance and calibration must be documented in the laboratory notebook kept at the EAA Camden Building. This documentation is an important part of ensuring that data-collection results are “defensible.” Calibration details, equipment type, date, calibration statement, and sampler’s signature must appear in the book for each day that the equipment is used.

EAA currently uses the YSI 556 MPS field instrument to collect pH, DO, conductivity, and temperature at each sample point. Calibration procedures for this instrument are detailed next.

CALIBRATION PROCEDURES

Calibration Procedures for YSI 556 MPS

Accessing the Calibrate Screen

1. Press the **On/Off** key to display the run screen.
2. Press the **Escape** key to display the main menu screen.
3. Use the arrow keys to highlight the **Calibrate** selection

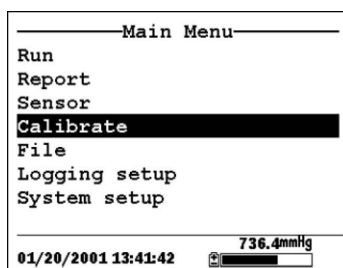


Figure 6.1 Main Menu

4. Press the **Enter** key. The Calibrate screen will be displayed.

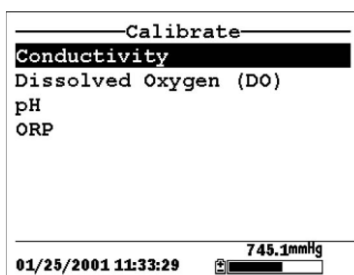


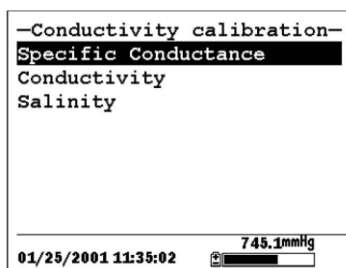
Figure 6.2 Calibrate Screen

Conductivity Calibration

This procedure calibrates specific conductance (recommended), conductivity, and salinity.

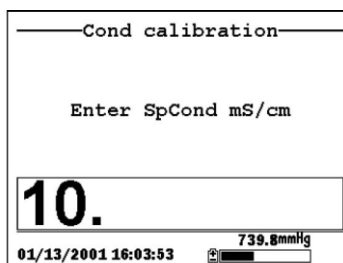
Calibrating any one option automatically calibrates the other two.

1. Go to the Calibrate screen
2. Use the arrow keys to highlight the **Conductivity** selection.
3. Press **Enter**. The Conductivity Calibration Screen is displayed.



Conductivity Calibration Selection Screen

4. Use the arrow keys to highlight the Specific Conductance selection.
5. Press **Enter**. The Conductivity Calibration Entry Screen is displayed.



Conductivity Calibration Selection Screen

6. Place the correct amount of conductivity standard into a clean, dry or pre-rinsed transport/calibration cup.

WARNING: Calibration reagents may be hazardous to health. See information on label.

NOTE: For maximum accuracy, the conductivity standard you choose should be within the same conductivity range as the samples you are preparing to measure. However, we do not recommend using standards less than one mS/cm. For example:

- For freshwater use a one-mS/cm conductivity standard.
- For brackish water use a ten-mS/cm conductivity standard.
- For seawater use a 50-mS/cm conductivity standard.

NOTE: Before proceeding, ensure that the sensor is as dry as possible. Ideally, rinse the conductivity sensor with a small amount of standard that can be discarded. Be certain that cross-contamination of solutions be avoided. Make certain that no salt deposits are around the oxygen or pH/ORP sensors, particularly if standards of low conductivity are being employed.

7. Carefully immerse the sensor end of the probe module into the solution.
8. Gently rotate and/or move the probe module up and down to remove any bubbles from the conductivity cell.

NOTE: The sensor must be completely immersed past its vent hole. Using the recommended volumes and ensure that the vent hole is covered.

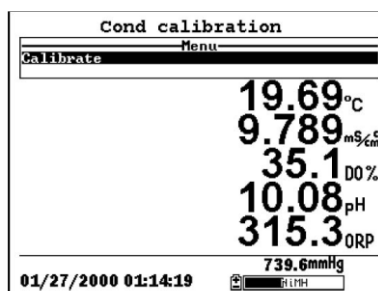
9. Screw the transport/calibration cup onto the threaded end of the probe module and securely tighten.

NOTE: Do not over tighten because doing so could damage the threaded parts.

10. Use the keypad to enter the calibration value of the standard being used.

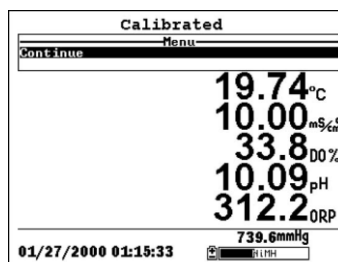
NOTE: Be sure to enter the value in **mS/cm at 25°C**.

11. Press **Enter**. The Conductivity Calibration Screen is displayed.



Conductivity Calibration Screen

12. Allow at least one minute for temperature equilibration before proceeding. The current values of all enabled sensors will appear on the screen and will change with time as they stabilize.
13. Observe the reading under Specific Conductance. When the reading shows no significant change for approximately 30 seconds, press **Enter**. The screen will indicate that the calibration has been accepted and prompt pressing of **Enter** again to Continue.



Calibrated

14. Press **Enter** to return to the Conductivity Calibrate Selection Screen
15. Press **Escape** to return to the Calibrate menu. See Figure 6.2 Calibrate Screen.
16. Rinse the probe module and sensors in tap or purified water and dry.

Dissolved Oxygen Calibration

This procedure calibrates dissolved oxygen. Calibrating any one option (% or mg/L) automatically calibrates the other.

1. Go to the calibrate screen as described in Section 6.2.1 *Accessing the Calibrate Screen*.

NOTE: The instrument must be on for at least 10 to 15 minutes to polarize the DO sensor before calibrating.

2. Use the arrow keys to highlight the **Dissolved Oxygen** selection. See Figure 6.2 Calibrate Screen.
3. Press **Enter**. The Dissolved Oxygen Calibration Screen is displayed.

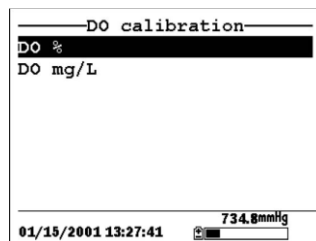


Figure 6.7 DO Calibration Screen

DO Calibration in Percent Saturation

1. Use the arrow keys to highlight the DO% selection.
2. Press **Enter**. The DO Barometric Pressure Entry Screen is displayed.

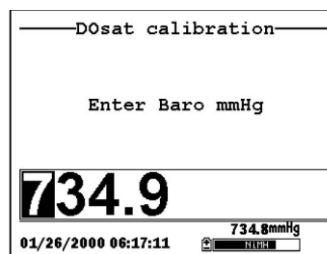


Figure 6.8 DO Barometric Pressure Entry Screen

3. Place approximately 3 mm ($\frac{1}{8}$ inch) of water in the bottom of the transport/calibration cup.
4. Place the probe module into the transport/calibration cup.
NOTE: Ensure that the DO and temperature sensors are **not** immersed in the water.
5. Engage only one or two threads of the transport/calibration cup to ensure that the DO sensor is vented to the atmosphere.
6. Use the keypad to enter the current local barometric pressure.

NOTE: If the unit has the optional barometer, no entry is required.

NOTE: Barometer readings that appear in meteorological reports are generally corrected to sea level and must be uncorrected before use

7. Press **Enter**. The DO% Saturation Calibration screen is displayed.

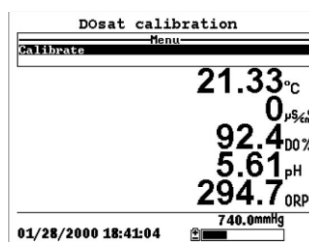


Figure 6.9 DO Sat Calibration Screen

8. Allow approximately ten minutes for the air in the transport/calibration cup to become water saturated and for the temperature to equilibrate before proceeding.
9. Observe the reading under DO %. When the reading shows no significant change for approximately 30 seconds, press **Enter**. The screen will indicate that the calibration has been accepted and prompt pressing of **Enter** again to Continue. See Figure 6.6 Calibrated.
10. Press **Enter** to return to the DO Calibration Screen, See Figure 6.7 DO Calibration Screen.
11. Press **Escape** to return to the calibrate menu. See Figure 6.2 Calibrate Screen.
12. Rinse the probe module and sensors in tap or purified water and dry.

pH Calibration

1. Go to the Calibrate Screen as described in *Section 6.2.1 Accessing the Calibrate Screen*.
2. Use the arrow keys to highlight the **pH** selection. See Figure 6.2 Calibrate Screen.
3. Press **Enter**. The pH Calibration screen is displayed.

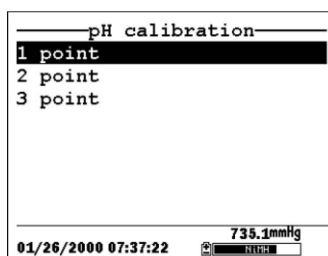


Figure 6.12 pH Calibration Screen

4. Select the **one-point** option only if adjusting a previous calibration. If a two-point or three-point calibration has been performed previously, the calibration can be adjusted by carrying out a one-point calibration. The procedure for this calibration is the same as for a two-point calibration, but the software will prompt a selection of only one pH buffer.
5. Select the **two-point** option to calibrate the pH sensor using only two calibration standards. Use this option if the media being monitored is known to be either basic or acidic. For example, if the pH of a pond is known to vary between 5.5 and seven, a two-point calibration with pH seven and pH four buffers is sufficient. A three-point calibration with an additional pH ten buffer will not increase the accuracy of this measurement because the pH is not within this higher range.
6. Select the **three-point** option to calibrate the pH sensor using three calibration solutions. In this procedure, the pH sensor is calibrated with a pH seven buffer and two additional buffers. The three-point calibration method assures maximum accuracy when the pH of the media to be monitored cannot be anticipated. The procedure for this calibration is the same as for a two-point calibration, but the software will prompt a selection of a third pH buffer.
7. Use the arrow keys to highlight the **two-point** selection.
8. Press **Enter**. The pH Entry Screen is displayed.

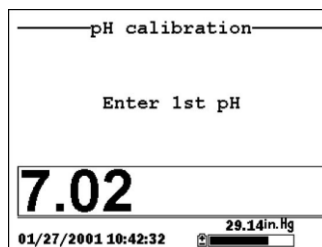


Figure 6.13 pH Entry Screen

9. Place the correct amount (see Table 6.1 Calibration Volumes) of pH buffer into a clean, dry, or prerinsed transport/calibration cup.

NOTE: Always calibrate with buffer seven first, regardless of whether performing a one-, two-, or three-point calibration.

WARNING: Calibration reagents may be hazardous to health. See reagent label for more information.

NOTE: For maximum accuracy, the pH buffers chosen should be within the same pH range as the water being prepared for sampling.

NOTE: Before proceeding, ensure that the sensor is as dry as possible. Ideally, rinse the pH sensor with a small amount of buffer that can be discarded. Be certain to avoid cross-contamination of buffers with other solutions.

10. Carefully immerse the sensor end of the probe module into the solution.
11. Gently rotate and/or move the probe module up and down to remove any bubbles from the pH sensor.

NOTE: The sensor must be completely immersed. Using the recommended volumes from Table 6.1 Calibration Volumes should ensure that the sensor is covered.
12. Screw the transport/calibration cup onto the threaded end of the probe module and securely tighten.

NOTE: Do not overtighten because doing so could damage the threaded parts.

13. Use the keypad to enter the calibration value of the buffer being used **at the current temperature**.

NOTE: pH vs. temperature values are printed on the labels of all YSI pH buffers.

14. Press **Enter**. The pH Calibration Screen is displayed.

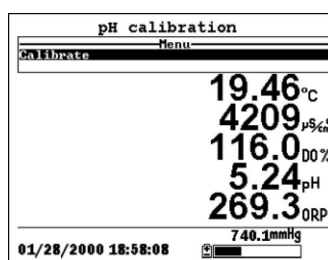


Figure 6.14 pH Calibration Screen

15. Allow at least one minute for temperature equilibration before proceeding. The current values of all enabled sensors will appear on the screen and will change with time as they stabilize.

16. Observe the reading under pH. When the reading shows no significant change for approximately 30 seconds, press **Enter**. The screen will indicate that the calibration has been accepted and prompt **Enter** to be pressed again to continue.
17. Press **Enter** to return to the specified pH Calibration Screen, See Figure 6.13 pH Entry Screen.
18. Rinse the probe module, transport/calibration cup, and sensors in tap or purified water and dry.
19. Repeat steps 6 through 13 using a second pH buffer.
20. Press **Enter** to return to the pH Calibration Screen. See Figure 6.12 pH Calibration Screen.
21. Press **Escape** to return to the Calibrate menu. See Figure 6.2 Calibrate Screen.
22. Rinse the probe module and sensors in tap or purified water and dry.

Return to Factory Settings.

1. Go to the Calibrate screen as described in Section 6.2.1 *Accessing the Calibrate Screen*.
2. Use the arrow keys to highlight the **Conductivity** selection. See Figure 6.2 Calibrate Screen.

NOTE: We will use the Conductivity sensor as an example; however, this process will work for any sensor.

3. Press **Enter**. The Conductivity Calibration Selection Screen is displayed. See Figure 6.3 Conductivity Calibration Selection Screen.
4. Use the arrow keys to highlight the **Specific Conductance** selection.
5. Press **Enter**. The Conductivity Calibration Entry Screen is displayed. See Figure 6.4 Conductivity Calibration Entry Screen.
6. Press and hold the **Enter** key down, and press the **Escape** key.

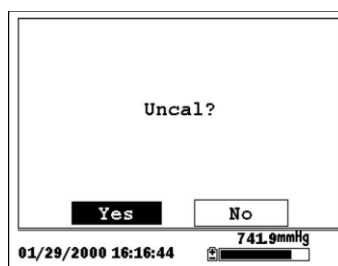


Figure 6.17 ORP Calibration Screen

7. Use the arrow keys to highlight the **YES** selection.

CAUTION: Pressing **YES** returns a sensor to the factory settings. For example, in the selection to return specific conductance to the factory setting, salinity and conductivity will automatically return to their factory settings.

8. Press **Enter** to return you to the Conductivity Calibrate Selection Screen. See Figure 6.3 Conductivity Calibration Selection Screen. .
9. Press **Escape** to return to the Calibrate menu. See Figure 6.2 Calibrate Screen.

HACH Digital Titrator (for Alkalinity) Primary Method

Titration is performed using the HACH digital titrator. This instrument provides precise results when properly operated.

Basic Operation

1. Select a sample volume and titration cartridge corresponding to the expected sample concentration.
2. Insert the cartridge into the titrator slide and lock it into place with the plunger. Remove the polyethylene cap from the cartridge and insert a clean delivery tube into the end of the cartridge. (Note: use a straight tube with a hook on the end for hand-held titrations and a 90° tube with a hook at the end for stationary setups.)
3. To start the titrant flow, hold the tip of the cartridge upward while turning the delivery knob until the air is expelled and several drops of solution flow from the tip of the delivery tube.
4. Use the counter reset knob (the smaller of the two knobs) to set the digital counter back to zero, then blot any titrant from the delivery tube.
5. Proceed with titration by submerging the tip of the delivery tube into the sample and turning the delivery knob to dispense the titrant. (Note: during the titration process, samples must be continuously stirred either manually or with the magnetic stirrer)

Calculations

HACH titration cartridge solutions are designed to give those numbers used in the titrations (reading from the digital meter) to be actual sample concentration in mg/L, or they are marked with conversion factors. If in the process of sample preparation, the amount of SAMPLE becomes less than 100 mL, the titration number must be multiplied by the divisional factor. For example, if the intended 100-mL sample is reduced to 25 mL ($\frac{1}{4}$ of 100 mL) during the sample-preparation process, then the final result must be multiplied by 4 ($25 \text{ mL} \times 4 = 100 \text{ mL}$) to obtain the result.

General Maintenance

1. For long-term storage the delivery tube should be removed, the polyethylene cap reattached, and the cartridge removed from the titrator body. DO NOT attempt to remove the cartridge from the titrator without recapping.
2. After use and removal from the cartridges, rinse the delivery tubes with deionized water to prevent clogging.

The titration process should be checked monthly by titration of a standard solution and recorded in the laboratory notebook. Acceptable results are obtained if the titration is within $\pm 3\%$ of the standard solution.

Alkalinity Determination Using the HACH Digital Titrator

Alkalinity of water is defined by its acid-neutralizing capacity. Once a sample has been collected, geochemical changes can alter the sample's alkalinity. Therefore, alkalinity samples are to be analyzed in the field or immediately upon returning to the EAA laboratory.

Procedure

Sample alkalinity is determined by titration with sulfuric acid to a pH of 4.5 and includes all carbonate, bicarbonate, and hydroxide present within the sample. Values are recorded as mg/L calcium carbonate.

1. Follow the steps outlined in HACH digital titrator usage, with the sulfuric acid cartridge as the active titrant and the 90° delivery tube as a stationary setup.
2. Set up the HACH titrator unit and attach the digital titrator to the rotational holder and clamp securely.
3. The pH and temperature probes should also be connected to the titrastir at the end of the rotational holder. For best results, attempt to have the ends of the delivery tube, pH probe, and temperature probe at the same level.
4. Rinse a 25-mL pipette three times with deionized water and then three times with the sample water to be tested. Pipette 25 mL of this sample into a clean 50-mL beaker. Record this amount on the corresponding field sheet.
5. Place the beaker on the stir plate, put a stir bar in the beaker, and turn on the stirring function.
6. Rotate the titrastir arm toward the sample beaker, submerging the probes and delivery tube. Note: ensure that the titrator counter is reset to zero and the outside of the delivery tube is free of sulfuric acid before submerging.
7. Turn on the pH meter and record the stabilized pH reading of the sample. Record this value on the corresponding field data sheet.
8. Titrate by turning the delivery knob until the pH is reduced to 4.5, which is the endpoint, and the amount of titrant used should be recorded.
9. Calculate the alkalinity by multiplying the amount of titrant used by the dilution factor, and record on the appropriate field data sheet.

Collect a second alkalinity sample every ten samples as a field duplicate, and analyze as outlined above. The field duplicate percent difference should not exceed $\pm 5\%$, where %D is defined as

$$[(X1 - X2) / X1] \times 100 = \%D \text{ (X1 = original sample, X2 = duplicate sample)}$$

(see next page for additional alkalinity procedures)

Additional Procedures for Alkalinity Analyses, University of Minnesota Methodology (to be incorporated into the EAA methodology)

Good Titration Practices

Aliquot Measurement

- ◆ Sample aliquots should be measured with the most accurate method available.
- ◆ Rinse the volumetric flask with sample water.
- ◆ Never rinse the titration flask with sample water.
- ◆ Rinse the titration flask with De-Ionized water between samples and air dry (glass) or shake dry (PMP plastic).
- ◆ An electronic balance is preferred over a volumetric flask is preferred over a graduated cylinder.
- ◆ A 0.1g scale is comparable to a volumetric flask.
- ◆ An electronic balance allows the size of sample aliquots to be varied.
- ◆ An electronic balance allows aliquot size to be reduced in high alkalinity samples which reduces titration time.

Titration Equipment

- ◆ Digital titrator should be periodically lubricated.
- ◆ Titrant cartridges must be kept tightly capped to prevent evaporation.
- ◆ Old, partially used titrant cartridges should be replaced.
- ◆ Don't try to use every drop of acid in the titrant cartridge - when it gets low start a new cartridge.
- ◆ Delivery tubes should be flushed with fresh titrant before use and rinsed after use.
- ◆ A magnetic stirrer (battery powered for field use) helps ensure thorough mixing.

Titration Procedures

- ◆ All chemical analyses should be replicated.
- ◆ Titrations are done in triplicate to allow comparison of results ensuring that reproducible results are obtained.
- ◆ Replicates that vary by more than two percent indicate interference or analytic error.
- ◆ Real time analysis of the results allows additional titrations and/or a change in procedure to identify the sources of the interference or error.
- ◆ Work consistently and quickly to limit degassing and precipitation in your sample bottle.
- ◆ Add acid uniformly to each aliquot as if performing the first titration.

Colorimetric

- ◆ Bromocresol Green / Methyl Red indicator dyes.
- ◆ pH 4.8 to 4.5 buffer solutions.
- ◆ Adding acid too quickly and incomplete equilibration will produce irregular results.
- ◆ Use buffered indicator solutions to define endpoint.
- ◆ Relies on human color interpretation.

Potentiometric

pH Endpoint

- ◆ Meter calibration is critical.
- ◆ Adding acid too quickly and incomplete equilibration will produce irregular results.
- ◆ Must allow for solution equilibration and meter stabilization.
- ◆ Uses one data point to determine endpoint.

Δ pH/ Δ v acid

- ◆ Must be done in uniform steps through the endpoint.
- ◆ Adding acid too quickly and incomplete equilibration will produce irregular results.
- ◆ Organics may shift endpoint.
- ◆ Uses two data points to determine endpoint.

Figure 1 shows a typical “S” shaped titration curve. The inflection point represents the true alkalinity of the sample and may not occur at exactly pH 4.5.

Gran Titration

- ◆ Uses many data points.
- ◆ Must be carried well past the endpoint.
- ◆ Requires graphical interpretation or linear regression.
- ◆ Adding acid too quickly and incomplete equilibration will produce non-linear trend.
- ◆ Presence of organics will produce non-linear trends.
- ◆ Least susceptible to operator error or chemical interference but should still be backed up by replicate measurements - replicate may be by colorimetric or potentiometric methods.

To calculate the alkalinity, use the formula $(V_{\text{aliquot}} + V_{\text{titrant}}) \times 10^{(4.65-\text{pH})}$ to plot an ascending line after the endpoint with apparent alkalinity on the x-axis; V_{aliquot} in ml, $V_{\text{titrant}} = \text{titrator digits}/800$ and 4.65 is the assumed endpoint. A linear regression can then be used to calculate an x-intercept. Use only the points well after the endpoint to get the best regression as shown in Figure 2.

Common Interferences

◆ **Highly colored waters**

Organic-rich waters with humic and fulvic acids.
Often have low pH and correspondingly low alkalinity.

Solutions

Perform Gran Titration - by extrapolating from points below pH 4.5 a fairly precise determination of alkalinity can be made.

Add a second packet of indicator dye to intensify green and red colors.

◆ **Chlorinated waters**

Color change at endpoint goes from green to yellow.

Solution: Add 5 drops and 2N Sodium Thiosulfate to scavenge any free chlorine before titrating.

◆ **Clay-rich waters**

Colors of indicators are "off" often tending towards an orange endpoint.
Commonly associated with poorly developed monitoring wells.

Solution: Filter the sample before titrating.

◆ **Muddy waters**

Suspended sediment may contain carbonates or clays that could react with the acid titrant.
Thick sediment may mask the color changes.

Solution: Filter the sample, preferably after allowing sediment to settle.

References

Determination of the Equivalent Point in Potentiometric Titrations, 1950, Gunnar Gran, Acta Chemica Scandinavica, pp 559-577.

Determination of the Equivalence Point in Potentionmetric Titrations - Part II, 1952, Gunnar Gran, The Analyst, International Congress on Analytical Chemistry, V. 77, pp 661-671.

Field Guide for Collecting and Processing Stream-Water Samples for the National Water-Quality Assessment Program, Larry R. Shelton, 1994, U.S. Geological Survey Open-File Report 94-455, 42 pp.

Field and Laboratory Methods, 1998, Scott C. Alexander and E.C. Alexander Jr., Hydrogeochemistry Lab, Dept. of Geology & Geophysics, Univ. of Minnesota, 21 pp.

EPA Method 310.1: Alkalinity determination to a colorimetric end-point.

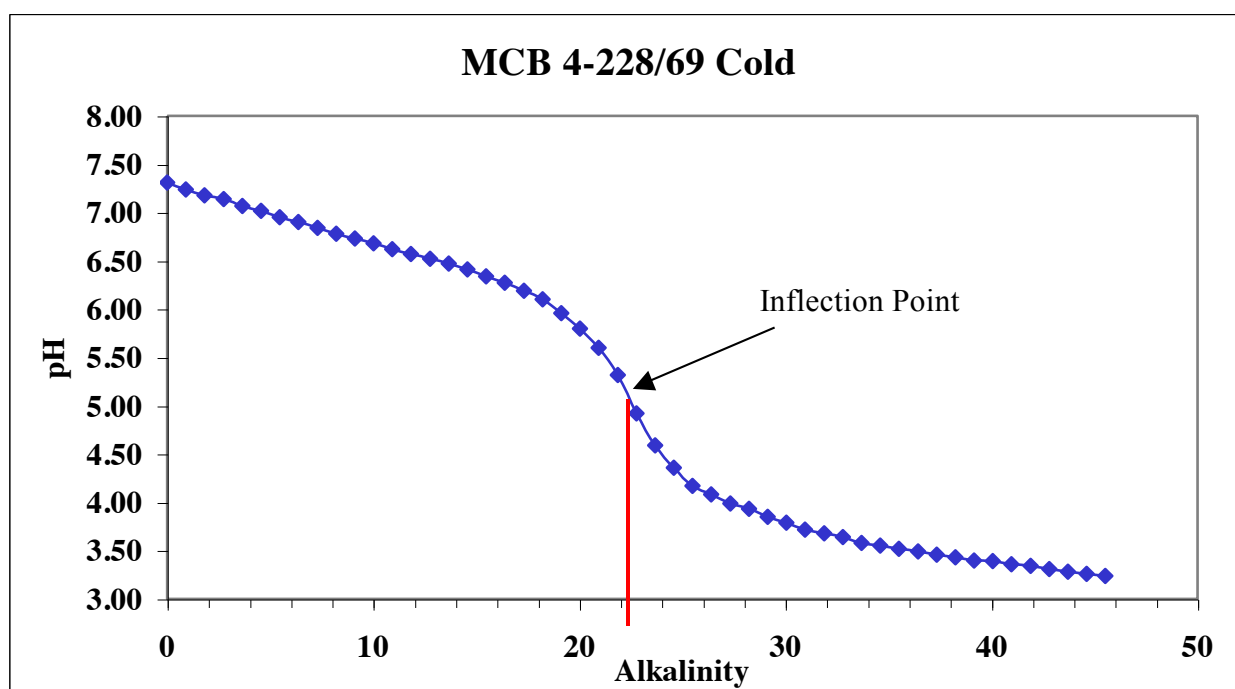


Figure 1. Example of pH titration

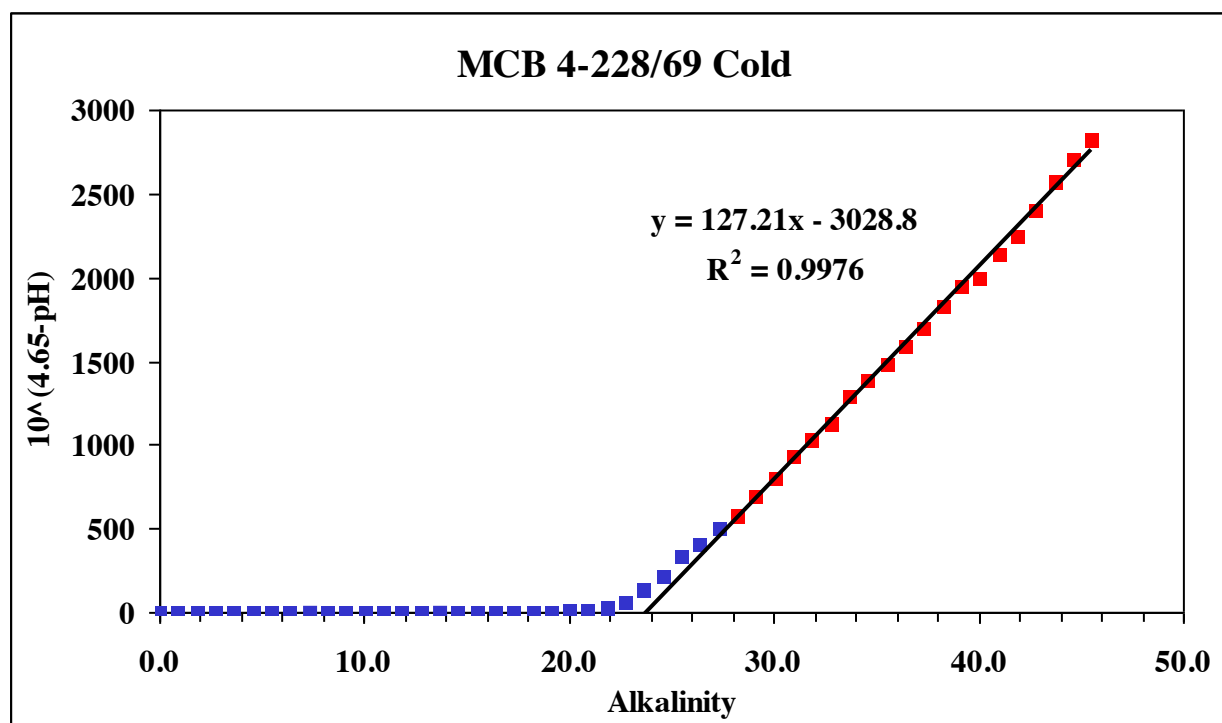


Figure 2. Example of Gran Titration

Use of the DR2800 Portable Spectrophotometer for Alkalinity Measurements (Secondary Method for Alkalinity Determination)

EAA currently uses the DR2800 Portable Spectrophotometer for measuring alkalinity values of samples in the event the Hach Digital Titrator is not available. Measurements are made at the EAA Camden building following the field sample-collection event. All measurements are to be recorded in the alkalinity notebook and on the field sheet. Operation procedures for this instrument are detailed next.

Alkalinity, Total

DOC316.53.01257

Colorimetric Method

Method 10239

25 to 400 mg/L CaCO₃

TNTplus™ 870

Scope and Application: For drinking water, wastewater and boiler water.

Test preparation

Before the test:

DR 2800 only: Install the light shield in Cell Compartment #2 before performing this test.

Read the safety advice and expiration date on the package.

The recommended sample and reagent temperature is 15–25 °C (59–77 °F).

The recommended reagent storage temperature is 15–25 °C (59–77 °F).

TNTplus™ methods are activated from the Main Menu when the sample vial is inserted into the sample cell holder.

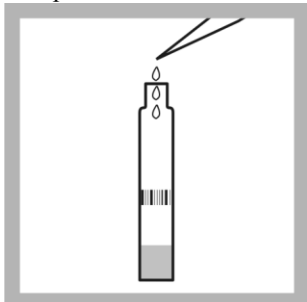
Collect the following items:

| Description | Quantity |
|-------------------------------|----------|
| Total alkalinity TNT870 vials | variable |
| Light shield (DR 2800 only) | 1 |
| Pipette for 2.0-mL sample | 1 |
| Pipette for 0.5-mL sample | 1 |
| Pipette tips | variable |

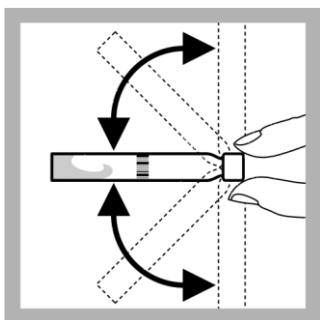
TNTplus™ method



1. Pipette 2.0 mL of **Solution A** into test vial.



2. Pipette 0.5 mL of sample into vial.



3. Cap and invert vial until contents are well mixed.



4. Wait 5 minutes.



5. After timer expires, wipe vial and insert it into cell holder. Instrument reads barcode, selects method, and make measurement. No instrument zero required. Results are in mg/L CaCO_3 .

Interferences

If samples contain particles, remove the particles by filtration through a 0.45- μm filter.

Sample collection, preservation, and storage

- Collect samples in clean plastic or glass bottles. Fill completely and cap tightly.
- Prevent excessive agitation or prolonged exposure to air. Complete the test procedure as soon as possible after collection for best accuracy.
- The sample can be stored for 24 h if cooled to 4 °C (39 °F) or below. Warm to room temperature before the test begins.

Accuracy check

Standard solution method required for accuracy check:

- Alkalinity Voluette® Ampule Standard Solution, 25,000 mg/L CaCO_3 (0.500 N)
- Ampule breaker
- Variable-volume pipette
- Pipette tips
- 100-mL volumetric flask, Class A
- Deionized water

1. Prepare a 250-mg/L CaCO_3 standard solution as follows:

- a.** Pipette 1.0 mL of alkalinity standard solution, 25,000 mg/L as CaCO_3 , into a clean 1.0-mL volumetric flask.
 - b.** Dilute to the mark with deionized water. Mix well. Prepare this solution daily.
- 2.** Use this solution in place of the sample. Follow the TNTplus™ method test procedure. The result should be within 10% of the expected value.

Summary of method

Carbonates and other buffers react with the reagent in the vial to change the pH. The pH affects the color of the indicator, which is measured photometrically at 615 nm.

Calibration Procedures for Backup Instruments

The following pages contain a discussion of proper use of “backup” instrumentation owned by the EAA, but not in regular use. These instruments may be utilized during a contingency sampling event, or in the case where newer instrumentation is not available due to damage or other issues.

Calibration Procedures for Luminescent Dissolved Oxygen Probe Model LDO10101 with HQ30d Meter

Before calibration:

The probe must have the correct service-life time stamp. Set the date and time in the meter before the probe is attached.

It is not necessary to recalibrate when moving a calibrated probe from one HQd meter to another if the additional meter is configured for the same calibration options.

To view the current calibration, push Select View Probe Data, then select View Current Calibration.

If any two probes are connected, push the **UP** or **DOWN** arrow to change to the single display mode in order to show the Calibrate option.

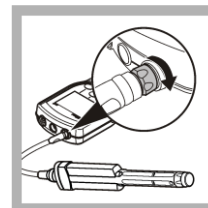
Calibration notes:

- % saturation or mg/L calibration methods are available in the Modify Current Settings menu.
- Slope value is the comparison between the latest calibration and the factory calibration shown as a percentage.
- Calibration is recorded in the probe and the data log. Calibration is also sent to a PC, printer, or flash memory stick if connected.

- Air bubbles under the sensor tip when submerged can cause slow response or error in measurement. If bubbles are present, gently shake the probe until bubbles are removed.

Water-saturated air (100%) calibration procedure:

1. Connect the probe to the meter. Ensure that the cable locking nut is securely connected to the meter. Turn on the meter.



2. Push **Calibrate**.



3. Push **Methods**. Select User Cal-100%. Push **OK**.



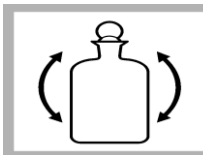
4. Rinse the probe cap with deionized water. Blot dry with a lint-free cloth.



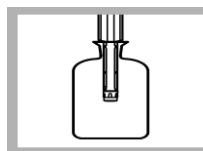
5. Add approximately ¼ inch (6.4 mm) of reagent water to a narrow-neck bottle, such as a BOD bottle.



6. Put a stopper in the bottle and shake the bottle vigorously for approximately 30 seconds to saturate the entrapped air with water. Allow up to 30 minutes for contents to equilibrate to room temperature.



7. Remove the stopper. Carefully dry the probe cap using a nonabrasive cloth. Put the probe in the bottle.



8. Push **Read**. The display shows “Stabilizing” as the probe stabilizes. The display shows the standard value when the reading is stable.

9. Push **Done** to view the calibration summary.

10. Push **Store** to accept the calibration and return the measurement mode. If a rugged probe, install the shroud on the probe.



Calibration Procedures for Conductivity Model 5197500 or 5197503 with sensION™ 5 (Backup Instrumentation)

Before the test:

Collect samples in clean plastic or glass bottles.

Analyze samples as soon as possible after collection. However, samples may be stored at least 24 h by cooling to 4 °C (39 °F) or below (all storage temperatures have changed to 0 to 6 °C as per the EPA MUR, March 2007). When solutions are measured that are not at reference temperature, the meter automatically adjusts the conductivity value to reference temperature from 20 or 25 °C.

Water samples containing oils, grease, or fats will coat the electrode and affect the accuracy of the readings. If this coating occurs, clean the probe with a strong detergent solution, then thoroughly rinse with deionized water.

Mineral buildup on the probe can be removed with a diluted 1:1 hydrochloric acid solution. Refer to the meter user's manual.

Calibration instructions are given in the operation section of the meter manual. For most accurate results, calibrate before use or check the accuracy of the meter using a known conductivity standard.

Calibrating with a Known Standard

1. Place the probe in a conductivity standard that is in the expected range of the samples. On the meter, choose one of four ranges that corresponds to the sample range. Agitate the probe to dislodge bubbles in the cell. Avoid resting the probe on the bottom or sides of the container.
2. Press **CAL**. Functional keys will appear in the lower-left part of the display. **CAL?** and **1.000 1/cm** will appear in the upper display. If the meter has been calibrated, the last calibration value will appear. The numeric keypad will become active.
3. Press the arrow keys to scroll to the factory-calibration options (1000 µS/cm or 18 mS/cm). To calibrate using one of

these standards, press **ENTER**.

4. If using a standard with a different value, use the number keys to enter the standard conductivity at 25 °C, then press **ENTER**. The meter will automatically correct the calibration measurement to 25 °C using the NaCl-based, non-linear temperature coefficient.

If the standard has a value of 25 °C in the $\mu\text{S}/\text{cm}$ range, enter the value when 1000 $\mu\text{S}/\text{cm}$ is displayed. If the standard has a value of 25 °C in the mS/cm range, enter the value when 18 mS/cm is displayed. All four places have a number entered in them. If a number entry error occurs, start over by pressing **SETUP/CE**.

5. When the reading is stable, the calibration is automatically stored, and the instrument returns to reading mode.

Calibration Procedures for Turbidimeter

Note: *for best accuracy, use the same sample cell of four matched sample cells for all measurements during calibration. Always insert the cell so that the orientation mark placed on the cell during the matching procedure is correctly aligned.*

Calibration

1. Rinse a clean sample cell with dilution water several times. Then fill the cell to the line (~15 mL) with dilution water or use StablCal <0.1 NTU standard. *Note: the same dilution water used for preparing the standards must be used in this step.*
2. Insert the sample cell in the cell compartment by aligning the orientation mark on the cell with the mark on the front of the cell compartment. Close the lid, and press **I/O**. *Note: choose signal average mode option (on or off) before pressing CAL— the SIGNAL AVERAGE key in calibration mode.*
3. Press **CAL**. The CAL and S0 icons will be displayed and will flash. The four-digit display will show the value of the S0 standard for the previous calibration. If the blank value were forced to 0.0, the display would be blank (as shown.) Press **→** for a numerical display.

Hach Company recommends the use of StableCal® Stabilized Formazin or formazin standards only for the calibration of Hach turbidimeters. Hach Company cannot guarantee the performance of the turbidimeter if calibrated with co-polymer styrene divinylbenzene beads of other suspension. DO NOT calibrate with Gelex® Secondary Standards.

4. Press **READ**. The instrument will count from 60 to 0 (67 to 0 if signal average is on), read the blank, and use it to calculate a correction factor for the 20 NTU standard measurement. If the dilution water is less than or equal to 0.5 NTU, E 1 will appear when the calibration is calculated. The display will automatically increment to the next standard. Remove the sample cell from the cell compartment. *Note: turbidity of the dilution water can be “forced” to zero by pressing → rather than reading the dilution water. The display will show S0 NTU, and the up arrow key must be pressed to continue with the next standard.*
5. The display will show the S1 (with the 1 flashing) and 20 NTU, or the value of the S1 standard for the previous calibration. If the value is incorrect, edit the value by pressing the → key until the number that needs editing flashes. Use the up arrow key to scroll to the correct number. After editing, fill a clean sample cell to the line with well-mixed 20 NTU StablCal Standard of 20 NTU formazin standard. Insert the sample cell compartment by aligning the orientation mark on the cell with the mark on the front of the cell compartment. Close the lid.
6. Press **READ**. The instrument will count from 60 to 0 (67 to 0 if signal average is on), measure the turbidity, and store the value. The display will automatically increment to the next standard. Remove the sample cell from the cell compartment. *Note: for potable water applications with low turbidity values, instrument calibration may be stopped after the 20 NTU StablCal Standard has been read. Press **CAL** after reading the 20-NTU standard. Instrument calibration is now complete for the range of 0–20 NTU only. The instrument will continue to read turbidity values above 20 NTU. These values were not updated during the 0–20 NTU calibration.*
7. The display will show the S2 (with the 2 flashing) and 100 NTU of the value of the S2 standard for the previous calibration. If the value is incorrect, edit the value by pressing the → key until the number that needs editing flashes. Use the up arrow key to scroll to the correct number. After

editing, fill a clean sample cell to the line with well-mixed 100 NTU StableCal Standard or 100 NTU formazin standard. Insert the sample cell into the cell compartment by aligning the orientation mark on the cell with the mark on the front of the cell compartment. Close the lid.

8. Press **READ**. The instrument will count from 60 to 0 (67 to 0 if signal average is on), measure the turbidity and store the value. Then the display will automatically increment to the next standard. Remove the sample cell from the cell compartment.
9. The display will show the S3 (with 3 flashing) and 800 NTU, or the value of the S3 standard for the previous calibration. If the value is incorrect, edit the value by pressing the → key until the number that needs editing flashes. Use the up arrow key to scroll to the correct number. After editing, fill a clean sample cell to the line with well-mixed 800 NTU formazin standard. Insert the sample cell into the cell compartment by aligning the orientation mark on the cell with the mark on the front of the cell compartment. Close the lid.
10. Press **READ**. The instrument will count from 60 to 0 (67 to 0 if signal average is on), measure the turbidity, and store the value. Then the display will increment back to the S0 display. Remove the sample cell from the cell compartment.
11. Press **CAL** to accept the calibration. The instrument will return to measurement mode automatically. *Note: pressing CAL completes the calculation of the calibration coefficients. If calibration errors occurred during calibration, error messages will appear after CAL is pressed. If E1 or E2 appear, check the standard preparation and review the calibration; repeat the calibration if necessary. If CAL? appears, an error may have occurred during calibration. If CAL? is flashing, the instrument is using the default calibration.*

Notes

- If the **I/O** key is pressed during calibration, the new calibration data are lost, and the old calibration will be used for measurements. Once in calibration mode, only the **READ**, **I/O**, ↑, and → keys function. Signal averaging and range mode must be selected before the

calibration mode can be entered.

- If **E 1** or **E 2** is displayed, an error occurred during calibration. Check the standard preparation and review the calibration; repeat the calibration if necessary. Press **DIAG** to cancel the error message (**E 1** or **E 2**). To continue without repeating the calibration, press **I/O** twice to restore the previous calibration. If **CAL?** is displayed, an error may have occurred during calibration. The previous calibration may not be restored. Either recalibrate or use the calibration as is.
- To review a calibration, press **CAL** and then \uparrow to view the calibration standard values. As long as **READ** is never pressed and **CAL** is not flashing, the calibration will not be updated. Press **CAL** again to return to the measurement mode.

pH Meter Calibration

The pH meter must be calibrated before daily use. The calibration may be accomplished in the laboratory or in the field. In addition to a “preuse” calibration, it is strongly recommended that the meter be checked with a standard buffer solution at least once during the day in order to observe any instrument drift that may have occurred.

Manual Calibration (with two reference solutions)

1. Attach or verify that the pH-indicating electrode and the automatic temperature compensator (ATC) are on the display unit.
2. Remove the rubber filling solution plug (if so equipped) to allow equilibration of the internal solution to the ambient air. Allow approximately five minutes for the equilibration process, and replace the plug.
3. Turn on the unit and select the calibration mode.
4. Rinse both electrodes with deionized water and dry (carefully) any excess water.
5. Rinse the pH electrode in the first pH buffer (reference) solution. After rinsing, immerse the electrode in a container of the first reference solution, and stir to remove bubbles on the electrode.
6. Allow the display to read **READY** and begin flashing. If the pH reading is within the manufacturer’s specifications (see equipment manual), press **YES**. If not, press **NO** and repeat the procedure. The first standard will subsequently be locked into the unit’s memory.
7. To calibrate the meter to the second pH reference solution, repeat steps 4, 5, and 6 USING the second solution.
8. Remove and rinse probes IN deionized water, and begin sample analysis. Otherwise the meter may be turned off; it will keep calibrating as long as the power source remains intact.

Electrode Care and Maintenance for pH Meters

The pH electrodes discussed above are of the temperature-compensating triode design. These probes are delicate and require careful handling. The probes should not be allowed to freeze and MUST be stored in a vial of the storage solution.

1. Inspect the probe for damage before each use. Verify that probes contain the appropriate levels of filling solution.
2. If filling-solution levels are low, more solution should be added. Use the Hach solution for Hach probes and the Orion solution for Orion probes.
3. If the probe appears sluggish when readings are taken, the filling solution should be drained and refilled with fresh solution.
4. During normal operations, the probe will become fouled with scale deposits and oils. Clean with laboratory-grade soap by soaking the probe in the soap solution and rinsing in deionized water. If fouling is not removed by this procedure, then a 0.1-N solution of HCL or HNO₃ can be used as a soaking media.
5. Probes must be stored in the electrode storage solution or in a 4.0-pH buffer solution. If probes are allowed to dry out, irreversible damage to the probe may occur.

Conductivity Probes

Orion Conductivity/Temperature Meters, Models 122, 126, 128, and 1230

Conductance, refers to the ability of a substance to carry an electrical current. These probes are used to define the physical parameters of conductivity. Conductivity is the algebraic reciprocal of electrical resistance and is expressed in SI units of microSeimens per centimeter. Specific conductance is electrical conductance measured across a one-cm cube of liquid (sample) between opposing faces of two platinum electrodes at 25°C. Conductivity is the same parameter measured at ambient temperature that has not been temperature compensated to 25°C.

Calibration

The conductivity meter must be calibrated in the laboratory or in the field daily. Conductance standards should be chosen to closely reflect the values expected in the sample groups. For example, if historical conductivity values for an area to be sampled range below 1000 $\mu\text{S}/\text{cm}$, the 500- $\mu\text{S}/\text{cm}$ solution should be chosen. The meters are designed to provide a nonlinear-function temperature coefficient to correct calculations; however, best results may be obtained when samples are 25°C.

Calibration Steps

- 1 Select conductivity measurement by turning the meter's conductivity/temperature selector knob from **OFF** to **CONDUCTIVITY** (labeled Δ).
- 2 Submerge the probe into THE selected conductivity standard (past the open area within the probe), and stir briefly to eliminate any air bubbles.
- 3 Maintain the probe in solution, wait for the reading to stabilize, and record the final value.
- 4 No manual adjustment for the meter exists; therefore, the process described herein provides a reference check. If the conductivity reading obtained from steps 1 through 3 is within $\pm 3\%$ of the given standard value, the meter is deemed to be within tolerance limits. If repeated attempts fail to obtain readings within the acceptable range, the meter will require factory service.

Maintenance

- 1 The meter electrode must be clean for readings to be accurate. Laboratory-grade soap may be used to clean dirt and oil deposits from the meter. For mineral deposits, a 1-M-HCl solution may be used in ten parts deionized water, and ten parts isopropyl alcohol as a soaking agent for their removal.
- 2 The conductivity probe may be stored dry. After each use, however, the probe should be rinsed in deionized water and blotted dry.
- 3 The unit will indicate a low battery by flashing **LOBAT** in the upper-left-hand corner of the LCD display. The nine-volt disposable battery should be changed out with the unit **OFF**, to prevent damage.

APPENDIX D—Forms



**EDWARDS AQUIFER
AUTHORITY**

Water Quality Field Data Sheet

HCP SEDIMENT

Site Information

| | |
|--|---------------|
| Station Name: | |
| Location: | |
| Owner/Contact: Edwards Aquifer Authority | |
| Address: 900 East Quincy | |
| County: | |
| Point of Collection: | |
| Date: / / 201 | Time: |
| Ambient Temp. | Collector(s): |
| Weather: | |

Equal-Width-Increment Method

| |
|----------------------|
| Transect Width: |
| Number of Verticals: |
| Flow/Appearance: |

| | | | | | |
|---|--------------|-------|------|------|------|
| Type of Analysis: (circle all that apply) | | | | | |
| GWQP | Select. Met. | 8081 | 8082 | 8141 | 8151 |
| T. | | | | | |
| TOC | Phosphorous | SVOCs | TB | DOC | VOC |

Notes

| |
|--|
| |
| |
| |
| |
| |
| |
| |

| | |
|--------------|------------|
| Latitude: | Longitude: |
| Datum: _____ | |

updated 06/06/13



**EDWARDS AQUIFER
AUTHORITY**

Water Quality Field Data Sheet

HCP STORM WATER

Site Information

| |
|--|
| Station Name: |
| Location: |
| Owner/Contact: Edwards Aquifer Authority |
| Address: 900 East Quincy |
| County: |
| Point of Collection: |
| Date: / / 201 Time: |
| Ambient Temp. Collector(s): |
| Weather: |

Field Readings

| |
|-------------------|
| Time Sampled: |
| pH: |
| Temperature: |
| Conductivity: |
| Dissolved Oxygen: |
| Turbidity: |

Equal-Width-Increment Method

| |
|------------------|
| Transect Width: |
| Flow/Appearance: |

Instrument Calibration

| | |
|----------------------|---------------|
| Conductivity Meter # | |
| Standard | Meter Reading |
| 500 | |
| 1000 | |
| 10000 | |
| pH Meter # | |
| Standard | Meter Reading |
| Buffer 4.0 | |
| Buffer 7.0 | |
| Buffer 10.0 | |
| pH Meter # | |
| Standard | Meter Reading |
| DI water in bottle | |

Alkalinity

| | |
|-------------------|----------------|
| Alkalinity Meter: | DR 2800 TNT870 |
| Total Alkalinity | |

| | | | | | |
|---|----------------|------|------------|------|------|
| Type of Analysis: (circle all that apply) | | | | | |
| GWQP | Selet. Met. | 8081 | 8082 | 8141 | 8151 |
| TOC | T. Phosphorous | TKN | E-Coli MPN | DOC | VOC |
| TB | SVOCs | | | | |

| | |
|-----------|------------|
| Latitude: | Longitude: |
|-----------|------------|

Datum: _____

Sampling Conditions

| | | |
|------------------|----------------------|-------|
| Gage Readings | Time | Level |
| Before Sampling | | |
| After Sampling | | |
| Hydrologic Event | Hydrologic Condition | |
| Storm | Stable, Low | |
| Drought | Falling | |
| Spill | Stable, High | |
| Regulated Flow | Rising | |
| Routine Sample | Stable, Normal | |

updated 12/21/12



**EDWARDS AQUIFER
AUTHORITY**

Water Quality Field Data Sheet

HCP SURFACE WATER

Site Information

| |
|--|
| Station Name: |
| Location: |
| Owner/Contact: Edwards Aquifer Authority |
| Address: 900 East Quincy |
| County: |
| Point of Collection: |
| Date: / / 201 Time: |
| Ambient Temp. Collector(s): |
| Weather: |

Field Readings

| |
|-------------------|
| Time Sampled: |
| pH: |
| Temperature: |
| Conductivity: |
| Dissolved Oxygen: |
| Turbidity: |

Equal-Width-Increment Method

| |
|------------------|
| Transect Width: |
| Flow/Appearance: |

Instrument Calibration

| Conductivity Meter # | |
|----------------------|---------------|
| Standard | Meter Reading |
| 500 | |
| 1000 | |
| 10000 | |
| pH Meter # | |
| Standard | Meter Reading |
| Buffer 4.0 | |
| Buffer 7.0 | |
| Buffer 10.0 | |
| pH Meter # | |
| Standard | Meter Reading |
| DI water in bottle | |

Alkalinity

| | mL of Sample | mL of Acid | Total Alk | |
|--------|--------------|------------|-----------|-----------------|
| Rep.1 | | | | Ave. Total Alk. |
| Rep. 2 | | | | |
| Rep3. | | | | |

| | | | | | |
|---|----------------|------|------------|------|------|
| Type of Analysis: (circle all that apply) | | | | | |
| GWQP | Selct. Met. | 8081 | 8082 | 8141 | 8151 |
| TOC | T. Phosphorous | TKN | E-Coli MPN | DOC | VOC |
| TB | SVOCs | | | | |

| | |
|-----------|------------|
| Latitude: | Longitude: |
|-----------|------------|

Datum: _____

Sampling Conditions

| Gage Readings | Time | Level |
|------------------|----------------------|-------|
| Before Sampling | | |
| After Sampling | | |
| Hydrologic Event | Hydrologic Condition | |
| Storm | Stable, Low | |
| Drought | Falling | |
| Spill | Stable, High | |
| Regulated Flow | Rising | |
| Routine Sample | Stable, Normal | |

updated 12/21/12



EDWARDS AQUIFER
AUTHORITY

Water Quality Field Data Sheet

SPRINGS

Spring Information

| |
|---|
| State well ID #: NA |
| Owner/Contact: New Braunfels Parks & Rec. |
| Address: |
| Phone Number: |
| County: Comal |
| Spring Name / #: Comal Springs 7 |
| Point of Collection: Springs Oriface |
| Spring Use: Springs |
| Date: Time: |
| Weather: Collector(s): |

Field Readings

| | | | | |
|---------------|------|-------|----|----|
| Time Sampled: | | | | |
| Turbidity: | | | | |
| Time | Temp | Cond. | pH | DO |
| | | | | |

Alkalinity

| | mL of Sample | mL of Acid | Total Alk | |
|--------|--------------|------------|-----------|-----------------|
| Rep. 1 | | | | Ave. Total Alk. |
| Rep. 2 | | | | |
| Rep. 3 | | | | |

Instrument Calibration

| Conductivity Meter | |
|--------------------|---------------|
| Standard | Meter Reading |
| 500 | |
| 1000 | |

| pH Meter | |
|-------------|---------------|
| Standard | Meter Reading |
| Buffer 4.0 | |
| Buffer 7.0 | |
| Buffer 10.0 | |

| DO Meter | |
|--------------------|---------------|
| Standard | Meter Reading |
| DI water in bottle | |

Type of Analysis: (circle all that apply)

| | | | | | |
|------------|--------------|--------|------|----------------|----------------------|
| GWQP | Select. Met. | 8081 | 8082 | 8141 | 8151 |
| 8260 | Trip Blank | SVOC'S | TOC | T. Phosphorous | Ortho-phosphate as P |
| E-Coli MPN | TPH | PAH | PPCP | | |

| | |
|--------------|------------|
| Latitude: | Longitude: |
| Datum: _____ | |

Sampling Conditions

| | | | |
|-----------------|-------|--------|-------|
| Spring Flow | Low | Medium | High |
| Flow Appearance | Clear | Cloudy | Murky |

updated 12/21/12



EDWARDS AQUIFER
AUTHORITY

Water Quality Field Data Sheet

SURFACE WATER

Site Information

| |
|-----------------------------|
| Station Name: |
| Location: |
| Owner/Contact: |
| Address: |
| County: |
| Point of Collection: |
| Date: Time: |
| Ambient Temp. Collector(s): |
| Weather: |

Field Readings

| |
|-------------------|
| Time Sampled: |
| pH: |
| Temperature: |
| Conductivity: |
| Dissolved Oxygen: |
| Turbidity: |

Equal-Width-Increment Method

| |
|----------------------|
| Transect Width: |
| Number of Verticals: |
| Flow/Appearance: |

Instrument Calibration

| | |
|----------------------|---------------|
| Conductivity Meter # | |
| Standard | Meter Reading |
| 500 | |
| 1000 | |
| 10000 | |
| pH Meter # | |
| Standard | Meter Reading |
| Buffer 4.0 | |
| Buffer 7.0 | |
| Buffer 10.0 | |
| pH Meter # | |
| Standard | Meter Reading |
| DI water in bottle | |

Alkalinity

| | mL of Sample | mL of Acid | Total Alk | |
|--------|--------------|------------|-----------|-----------------|
| Rep.1 | | | | Ave. Total Alk. |
| Rep. 2 | | | | |
| Rep3. | | | | |

Type of Analysis: (circle all that apply)

| | | | | | |
|------|----------------|----------------------|------------|------|------|
| GWQP | Select. Met. | 8081 | 8082 | 8141 | 8151 |
| TOC | T. Phosphorous | Ortho-phosphate as P | E-Coli MPN | TPH | PAH |
| PPCP | | | | | |

| | |
|-----------|------------|
| Latitude: | Longitude: |
|-----------|------------|

Datum: _____

Sampling Conditions

| Gage Readings | Time | Level |
|------------------|----------------------|-------|
| Before Sampling | | |
| After Sampling | | |
| Hydrologic Event | Hydrologic Condition | |
| Storm | Stable, Low | |
| Drought | Falling | |
| Spill | Stable, High | |
| Regulated Flow | Rising | |
| Routine Sample | Stable, Normal | |

updated 12/21/12



Water Quality Field Data Sheet
GROUNDWATER

Well Information

Field Readings

| State well ID #: Owner/Contact: Address: Phone Number: County: Well Name / #: Point of Collection: Well Use: Weather: Date: Time: Flow Rate: gpm Collector(s): Water Level: Well Depth: Casing Dia.: 3 x well volume= | Started Pumping: Time Sampled: Turbidity: <table border="1" style="width: 100%; border-collapse: collapse; text-align: center;"> <tr> <th>Time</th> <th>Temp</th> <th>Cond.</th> <th>pH</th> <th>Do</th> </tr> <tr><td> </td><td> </td><td> </td><td> </td><td> </td></tr> <tr><td> </td><td> </td><td> </td><td> </td><td> </td></tr> <tr><td> </td><td> </td><td> </td><td> </td><td> </td></tr> <tr><td> </td><td> </td><td> </td><td> </td><td> </td></tr> <tr><td> </td><td> </td><td> </td><td> </td><td> </td></tr> <tr><td> </td><td> </td><td> </td><td> </td><td> </td></tr> </table> | Time | Temp | Cond. | pH | Do | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|---|---|-------|------|-------|----|----|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|
| Time | Temp | Cond. | pH | Do | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |

Instrument Calibration

| Conductivity Meter # | |
|----------------------|---------------|
| Standard | Meter Reading |
| 500 | |
| 1000 | |
| 10000 | |
| pH Meter # | |
| Standard | Meter Reading |
| Buffer 4.0 | |
| Buffer 7.0 | |
| Buffer 10.0 | |
| DO Meter | |
| Standard | Meter Reading |
| DI water in bottle | |

Alkalinity

| | mL of Sample | mL of Acid | Total Alk | |
|--------|--------------|------------|-----------|--|
| Rep.1 | | | | Ave. Total Alk. <div style="border: 1px solid black; height: 20px; width: 100%;"></div> |
| Rep. 2 | | | | |
| Rep3. | | | | |

| | |
|------------------|-------------------|
| Latitude: | Longitude: |
|------------------|-------------------|

Datum:

| Type of Analysis: (circle all that apply) | | | | | |
|---|-------------|--------|------|----------------|----------------------|
| GWQP | Selct. Met. | 8081 | 8082 | 8141 | 8151 |
| 8260 | Trip Blank | SVOC'S | TOC | T. Phosphorous | Ortho-phosphate as P |
| E-Coli MPN | TPH | PAH | PPCP | | |

Updated on 12/21/12

TWDB Water Quality Field Data Sheet

FY06

Newly Inventoried Well

SWN: _____
 County: _____
 County Code: _____
 Aquifer Code: _____
 Aquifer Id: _____
 Name: _____
 Address: _____
 Phone Number: _____
 Attention: _____
 Well Name or #: _____

| Calibration Verification Readings | |
|-----------------------------------|------------------------------|
| pH | 7 = _____ 4 or 10 = _____ |
| SLP = | 7.38 = _____ |
| Conductivity | 500 = _____ |
| | 1000 = _____ |
| | 2000 = _____ |
| | 5000 = _____ |

| CIRCLE EACH SAMPLE FRACTION COLLECTED: | | | | | | | | | |
|--|---------------------------|--------------------------------------|---|---|---|---|---|---|----|
| 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 |
| 500ml filtered Anions/T. Alk. | 250 ml filtered Cation | 500ml filtered Nitrate | | | | | | | |
| Ice | (HNO ₃) | Ice + H ₂ SO ₄ | | | | | | | |

All acidified samples pH <2.0. (*) If natural pH <7, then add NaOH until pH is >7. If natural pH is >7, no NaOH required.

Time In: _____

Water Level: _____

Pumping time: _____

Well Use: _____

Lift: _____

Power: _____

Casing Type: _____

Time Out: _____

W.L. remark: _____

M.P. = _____

Sampling Point: _____

FIELD G.P.S. readings

Latitude: _____

Longitude: _____

Datum: _____

Casing Size: _____

Filter pressure: hand pump / line / spring

Sample Time: _____

| Field Alkalinity Titration: | | |
|--|----------|--------|
| | Start pH | End pH |
| 50.0 mL Sample Size | | |
| mL Acid added for Phenol (> 8.3) | | |
| mL Acid added for Total (to pH 4.5) | | |
| Items below calculated from: mL acid added x 20 = Alkalinity | | |
| Phenol Alkalinity (8224): | | mg/L |
| Total Alkalinity (3808): | | mg/L |

Items Below Calculated Later From Results:

| | |
|-----------------------------------|-------|
| Dissolved Solids (mg/L): | _____ |
| Hardness (as CaCO ₃): | _____ |
| Balanced: | _____ |

| Water Quality Stabilization Parameters Table (At least 3 readings @ 5 min. intervals) | | | | | Notes: |
|---|--|--|--|--|--------|
| Time | | | | | |
| pH | | | | | |
| Celsius Temp. | | | | | |
| Conductivity | | | | | |

Page 1 of 1

**APPENDIX E—Drinking-Water Standards and Chemical Health Effects,
from 30 TAC 290, RG-346, and U.S. EPA, July 2002**

(Note, regulatory limits change frequently for certain compounds, the data herein are for general comparisons. The reader should utilized the most recent data available online from TCEQ and EPA if sample results exceed regulatory limits)

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water[^] | Sources of Contaminant in Drinking Water[^] |
|--|--|---|---|
| Temperature (°C) EPA 170.1 | NE | NA | NA |
| pH measured at 25°C EPA 150.1 | >7.0* | NA | NA |
| Turbidity (NTU) | NE | NA | NA |
| Dissolved oxygen (DO) (mg/L) | NE | NA | NA |
| Alkalinity total as CaCO ₃ SM 2320 B (mg/L) | NE | NA | NA |
| Specific conductance µS/cm | NE | NA | NA |
| | | | |
| Laboratory | | NA | NA |
| Alkalinity total as CaCO ₃ SM 2320 B | NE | NA | NA |
| Bicarbonate (HCO ₃) SM 2320 B | NE | NA | NA |
| Fecal coliform (CFU/100 mL) | 0 MCLG ¹ | NA | NA |
| Fecal strep (CFU/100 mL) | 0 MCLG ¹ | NA | NA |
| E. coli (CFU/100 mL) | 0 MCLG ¹ | NA | NA |
| pH measured at 25°C EPA 150.1 | >7.0* | NA | NA |
| Specific conductance µS/cm | NE | NA | NA |
| | | | |
| Nutrients (mg/L) | | | |
| Nitrate-nitrite as N EPA354.1/300.0 | 10 | Infants below the age of six months who drink water containing nitrate in excess of the MCL could become seriously ill and, if untreated, may die. Symptoms include shortness of breath and blue-baby syndrome. | Runoff from fertilizer use; leaching from septic tanks, sewage; erosion of natural deposits |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water [^] | Sources of Contaminant in Drinking Water [^] |
|---|---|---|--|
| Nitrate as N E300 | 10 | Infants below the age of six months who drink water containing nitrate in excess of the MCL could become seriously ill and, if untreated, may die. Symptoms include shortness of breath and blue-baby syndrome. | Runoff from fertilizer use; leaching from septic tanks, sewage; erosion of natural deposits |
| Orthophosphate EPA 365.3 | NE | NA | NA |
| Ammonia as N SM 4500 | NE | NA | NA |
| Phosphorus | NE | NA | NA |
| | | | |
| Major Ions (mg/L) | | NA | NA |
| Sulfate (SO ₄) EPA 300.0 | 300* | NA | NA |
| Solids total dissolved (TDS) EPA 160.1 | 1,000* | NA | NA |
| Solids total suspended (TSS) EPA 160.2 | NE | NA | NA |
| Bromide (Br) EPA 300.0 | NE | NA | NA |
| Chloride (Cl) EPA 300.0 | 300* | NA | NA |
| Fluoride (F) EPA 340.2 | 2.0* | Bone disease (pain and tenderness of the bones); children may get mottled teeth | Water additive that promotes strong teeth, erosion of natural deposits, discharge from fertilizer and aluminum factories |
| | | | |
| Metals by EPA 200.7 and 200.8 (µg/L) | | NA | NA |
| Aluminum | 24,000** | NA | NA |
| | | NA | NA |
| Antimony | 6 | Increase in blood cholesterol; decrease in blood sugar | Discharge from petroleum refineries, fire retardants, ceramics, electronics, solder |
| Arsenic | 5 | Skin damage or problems with circulatory systems and increased risk of cancer | Erosion of natural deposits; runoff from orchards and glass and electronics production wastes |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water[^] | Sources of Contaminant in Drinking Water[^] |
|-------------------------------------|--|--|--|
| Barium | 2,000 | Increase in blood pressure | Discharge of drilling wastes, discharge from metal refineries, erosion of natural deposits |
| Beryllium | 4 | Intestinal lesions | Discharge from metal refineries and coal-burning factories. erosion of natural deposits |
| Boron | 4,900** | | |
| Cadmium | 5 | Kidney damage | Corrosion of galvanized pipe, erosion of natural deposits, discharge from metal refineries, runoff from waste batteries and paints |
| Chromium | 100 | Allergic dermatitis | Discharge from steel and pulp mills, erosion of natural deposits |
| Cobalt | 1,500** | NA | NA |
| Copper | 1,300* | Short-term exposure, gastrointestinal distress; long-term exposure, liver or kidney damage. People with Wilson's disease should consult their personal doctor if the amount of copper in their water exceeds the action level. | Corrosion of household plumbing systems, erosion of natural deposits |
| Iron | 300* | NA | NA |
| Lead | 15 | Infants and children: delays in physical or mental development; children could show slight deficits in attention span and learning abilities. Adults: Kidney problems, high blood pressure | Corrosion of household plumbing systems, erosion of natural deposits |
| Lithium | 490** | NA | NA |
| Manganese | 1,100* | NA | NA |
| Molybdenum | 120** | NA | NA |
| Nickel | 490** | NA | NA |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water [^] | Sources of Contaminant in Drinking Water [^] |
|--|---|--|---|
| Selenium | 50 | Hair or fingernail loss, numbness in fingers or toes, circulatory problems | Discharge from petroleum refineries, erosion of natural deposits, discharge from mines |
| Silver | 120* | NA | NA |
| Strontium | 15,000** | NA | NA |
| Thallium | 2 | Hair loss; changes in blood; kidney, intestine, or liver problems | Leaching from ore processing sites; discharge from electronics, glass, and drug factories |
| Uranium | 30 | NA | NA |
| Vanadium | 1.7** | NA | NA |
| Zinc | 7,300* | NA | NA |
| | | NA | NA |
| Metals by E200.8 (mg/L) | | | |
| Calcium | NE | NA | NA |
| Magnesium | NE | NA | NA |
| Potassium | NE | NA | NA |
| Sodium | NE | NA | NA |
| | | | |
| Metals by SW-7470A (mg/L) | | | |
| Mercury | 0.002 | Kidney damage | Erosion of natural deposits, discharge from refineries and factories, runoff from landfills and croplands |
| | | | |
| Total Organic Carbon by E415.1 (mg/L) | | | |
| TOC | NE | NA | NA |
| | | | |
| Herbicides by SW-8141 (µg/L) | | | |
| Azinphosmethyl | 37** | NA | NA |
| Bolstar (Sulprofos) | 73** | NA | NA |
| Chlorpyrifos | 73** | NA | NA |
| Coumaphos | 170** | NA | NA |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water[^] | Sources of Contaminant in Drinking Water[^] |
|--|--|---|---|
| Demeton-O | 1.0** | NA | NA |
| Demeton-S | 0.98** | NA | NA |
| Diazinon | 22** | NA | NA |
| Dichlorvos | 3.1** | NA | NA |
| Dimethoate | 4.9** | NA | NA |
| Disulfoton | 0.98** | NA | NA |
| EPN | 0.24** | NA | NA |
| Ethoprop | 2.4** | NA | NA |
| Famphur | 0.73** | NA | NA |
| Fensulfothion | 24** | NA | NA |
| Fenthion | 1.7** | NA | NA |
| Malathion | 490** | NA | NA |
| Merphos | 7.3** | NA | NA |
| Methyl parathion | 6.1** | NA | NA |
| Mevinphos (Phosdrin) | 0.61** | NA | NA |
| Mononcrotophos | 15** | NA | NA |
| Naled | 49** | NA | NA |
| Parathion | 150** | NA | NA |
| Phorate | 4.9** | NA | NA |
| Ronnel | 1,200** | NA | NA |
| Stirophos (Tetrachlorvinphos) | 1,000** | NA | NA |
| Sulfotepp (Tetraethyl dithiopyrophosphate) | 12** | NA | NA |
| Tokuthion (Prothiofos) | 2.4** | NA | NA |
| Trichloronate | 73** | NA | NA |
| Thionazin | 1.7** | NA | NA |
| | | | |
| Herbicides by SW-8151 (µg/L) | | | |
| 2,4,5-T | 240 | NA | NA |
| 2,4,5-TP (Silvex) | 50 | Liver problems | Residue of banned herbicide |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water[^] | Sources of Contaminant in Drinking Water[^] |
|---|--|---|--|
| 2,4- D | 70 | Kidney, liver, or adrenal gland problems | Runoff from herbicide used on row crops |
| 2,4-DB | 200 | NA | NA |
| Dalapon | 200 | Minor kidney changes | Runoff from herbicide used on rights of way |
| Dicamba | 730 | NA | NA |
| Dichoroprop | 240 | NA | NA |
| Dinoseb | 7 | Reproductive difficulties | Runoff from herbicide used on soybeans and vegetables |
| MCPA | 12 | NA | NA |
| MCP (mecoprop) | 24 | NA | NA |
| Pentachlorophenol | 1 | Liver or kidney problems, increased cancer risk | Discharge from wood-preserving factories |
| | | | |
| Pesticides by SW-8081 (µg/L) | | | |
| 4, 4'-DDD | 3.8** | NA | NA |
| 4, 4'-DDE | 2.7** | NA | NA |
| 4, 4'-DDT | 2.7** | NA | NA |
| Aldrin | 0.05** | NA | NA |
| Alpha-bhc (Alpha-hexachlorocyclohexane) | 0.1** | NA | NA |
| Alpha-chlordane | 2.6** | NA | NA |
| Beta-bhc (Beta-hexachlorocyclohexane) | 0.5** | NA | NA |
| Chlordane | 2.0** | Liver or nervous system problems, increased risk of cancer | Residue of banned termiticide |
| Delta-bhc (Delta-hexachlorocyclohexane) | 0.5** | NA | NA |
| Diieldrin | 0.57** | NA | NA |
| Endosulfan I | 49** | NA | NA |
| Endosulfan II | 150** | NA | NA |
| Endosulfan sulfate | 150** | NA | NA |
| Endrin | 2.0** | Liver problems | Residue of banned insecticide |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water[^] | Sources of Contaminant in Drinking Water[^] |
|-------------------------------------|--|---|---|
| Endrin aldehyde | 7.3** | NA | NA |
| Endrin ketone | == | | |
| Gamma-bhc (Lindane) | 7.3** | NA | NA |
| Gamma-chlordane | 0.2 | Liver or kidney problems | Runoff/leaching from insecticide used on cattle, lumber, gardens |
| | 2.6** | NA | NA |
| Heptachlor epoxide | 0.4 | Liver damage, increased risk of cancer | Residue of banned termiticide |
| Methoxychlor | 0.2 | Liver damage, increased risk of cancer | Breakdown of heptachlor |
| Toxaphene | 40 | Reproductive difficulties | Runoff/leaching from insecticide used on fruits, vegetables, alfalfa, livestock |
| PCBs by SW-8082 (µg/L) | 3 | Kidney, liver, or thyroid problems; increased risk of cancer | Runoff/leaching from insecticide used on cotton and cattle |
| Aroclor 1016 | | | |
| Aroclor 1221 | 0.5 | Skin changes, thymus gland problems, immune deficiencies, reproductive or nervous system difficulties, increased risk of cancer | Runoff from landfills, discharge of waste chemicals |
| Aroclor 1232 | 0.5 | Skin changes, thymus gland problems, immune deficiencies, reproductive or nervous system difficulties, increased risk of cancer | Runoff from landfills, discharge of waste chemicals |
| Aroclor 1242 | 0.5 | Skin changes, thymus gland problems, immune deficiencies, reproductive or nervous system difficulties, increased risk of cancer | Runoff from landfills, discharge of waste chemicals |
| Aroclor 1248 | 0.5 | Skin changes, thymus gland problems, immune deficiencies, reproductive or nervous system difficulties, increased risk of cancer | Runoff from landfills, discharge of waste chemicals |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water [^] | Sources of Contaminant in Drinking Water [^] |
|---------------------------------|---|---|---|
| Aroclor 1254 | 0.5 | Skin changes, thymus gland problems, immune deficiencies, reproductive or nervous system difficulties, increased risk of cancer | Runoff from landfills, discharge of waste chemicals |
| Aroclor 1260 | 0.5 | Skin changes, thymus gland problems, immune deficiencies, reproductive or nervous system difficulties, increased risk of cancer | Runoff from landfills, discharge of waste chemicals |
| Aroclor 1262 | 0.5 | Skin changes, thymus gland problems, immune deficiencies, reproductive or nervous system difficulties, increased risk of cancer | Runoff from landfills, discharge of waste chemicals |
| Aroclor 1268 | 0.5 | Skin changes, thymus gland problems, immune deficiencies, reproductive or nervous system difficulties, increased risk of cancer | Runoff from landfills, discharge of waste chemicals |
| | 0.5 | Skin changes, thymus gland problems, immune deficiencies, reproductive or nervous system difficulties, increased risk of cancer | Runoff from landfills, discharge of waste chemicals |
| SVOCs by SW-8270C (µg/L) | | | |
| 1,2- dichlorobenzene | 600** | NA | NA |
| 1,2,4- trichlorobenzene | | | |
| 2, 4, 5-trichlorophenol | 70** | Changes in adrenal glands | Discharge from textile finishing factories |
| 2, 4, 6-trichlorophenol | 2,400** | NA | NA |
| 2, 4-dichlorophenol | 24** | NA | NA |
| 2, 4-dimethylphenol | 73** | NA | NA |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water[^] | Sources of Contaminant in Drinking Water[^] |
|---|--|---|---|
| 2, 4-dinitrophenol | 490** | NA | NA |
| 2-chlorophenol | 49** | NA | NA |
| 2-methylnaphthalene | 120** | NA | NA |
| 2-methylphenol (o-cresol) | 98** | NA | NA |
| 2-nitroaniline | 1,200** | NA | NA |
| 2-nitrophenol | 7.3** | NA | NA |
| 3 & 4 methylphenol (m&p cresol) | 49** | NA | NA |
| 3-nitroaniline | 1,200** | NA | NA |
| 4, 6-dinitro-2-methylphenol | 7.3** | NA | NA |
| 4-chloro-3-methylphenol | 2.4** | NA | NA |
| 4- chloroaniline | 120** | NA | NA |
| 4-nitroaniline | 4.6** | NA | NA |
| 4-nitrophenol | 46** | NA | NA |
| Naphthalene | 49** | NA | NA |
| Nitrobenzene | 490** | NA | NA |
| Pentachlorophenol | 49** | NA | NA |
| Phenanthrene | 1 | NA | NA |
| Phenol | 730** | NA | NA |
| Pyrene | 7,300** | NA | NA |
| N-nitrosodi-n-propylamine | 730** | NA | NA |
| N-nitrosodiphenylamine | 0.13** | NA | NA |
| Acenaphthene | 190** | NA | NA |
| Acenaphthylene | 1,500** | NA | NA |
| Anthracene | 1,500** | NA | NA |
| Benzo(a)anthracene (1 2-benzanthracene) | 7,300** | NA | NA |
| Benzo(b)fluoranthene | 1.3** | NA | NA |
| Benzo(k)fluoranthene | 1.3** | NA | NA |
| Benzo(ghi)perylene | 13** | NA | NA |
| Benzo(a)pyrene | 730** | NA | NA |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water[^] | Sources of Contaminant in Drinking Water[^] |
|-------------------------------------|--|--|--|
| Benzyl Alcohol | 0.2 | Reproductive difficulties, increased risk of cancer | Leaching from linings of water storage tanks and distribution lines |
| Butyl benzyl phthalate | 2,400** | NA | NA |
| Bis(2-chloroethoxy)methane | 480** | NA | NA |
| Bis(2-chloroethyl)ether | 0.83** | NA | NA |
| Bis(2-ethylhexyl)phthalate | 0.83** | NA | NA |
| 4-bromophenyl phenyl ether | 6 | NA | NA |
| 4-chloroaniline | 0.061** | NA | NA |
| 2-chloronaphthalene | 4.6** | NA | NA |
| 4-chlorophenyl phenyl ether | 2,000** | NA | NA |
| Chrysene | 0.061** | NA | NA |
| Dibenz(a,h)anthracene | 130** | NA | NA |
| Dibenzofuran | 0.2** | NA | NA |
| 3,3-dichlorobenzidine | 98** | NA | NA |
| Diethyl phthalate | 2** | NA | NA |
| Dimethyl phthalate | 20,000** | NA | NA |
| Di-n-butyl phthalate | 20,000** | NA | NA |
| Di-n-octyl phthalate | 2,400** | NA | NA |
| 2,4-dinitrotoluene | 980** | NA | NA |
| 2,6-dinitrotoluene | 1.3** | NA | NA |
| Fluoranthene | 1.3** | NA | NA |
| Fluorene | 980** | NA | NA |
| Hexachlorobenzene | 980** | NA | NA |
| Hexachlorobutadiene | 1** | Liver or kidney problems, reproductive difficulties, increased risk of cancer | Discharge from metal refineries and agricultural chemical factories |
| Hexachlorocyclopentadiene | 12** | NA | NA |
| Hexachloroethane | 50 | Kidney or stomach problems | Discharge from chemical factories |
| Indeno(1,2,3-cd)pyrene | 24** | NA | NA |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water[^] | Sources of Contaminant in Drinking Water[^] |
|---|--|---|--|
| Isophorone | 1.3** | NA | NA |
| VOCs SW-8260b (µg/L) | 960** | NA | NA |
| 1, 1, 1, 2-tetrachloroethane | | | |
| 1, 1, 1-trichloroethane | 35.0** | NA | NA |
| 1, 1, 2, 2-tetrachloroethane | 200 | Liver, nervous system, or circulatory problems | Discharge from metal degreasing sites and other factories |
| 1, 1, 2-trichloroethane | 4.6** | | |
| 1, 1-dichloroethane | 5 | Liver, kidney, or immune system problems | Discharge from industrial chemical factories |
| 1, 1-dichloropropene | 4,900** | NA | NA |
| 1, 1-dichloroethene (Vinylidene chloride) | 9.1** | NA | NA |
| 1- chlorohexane | 7 | NA | NA |
| 1-octene | 980** | NA | NA |
| 1, 2, 3-trichlorobenzene | NE | NA | NA |
| 1, 2, 3-trichloropropane | 73** | NA | NA |
| 1, 2, 4-trichlorobenzene | 0.03** | NA | NA |
| 1, 2, 4-trimethylbenzene | 72** | NA | NA |
| 1, 2-dibromo-3-chloropropane | 1,200** | NA | NA |
| 1, 2-dibromoethane (EDB) | 0.2 | Reproductive difficulties, increased risk of cancer | Runoff/leaching from soil fumigant used on soybeans, cotton, pineapples, and orchards |
| 1, 2-dichlorobenzene | NE | NA | NA |
| 1, 2-dichloroethane (EDC) | 600** | NA | NA |
| 1, 2-dichloropropane | 5 | Increased risk of cancer | Discharge from industrial chemical factories |
| 1, 3, 5-trimethylbenzene | 5 | Increased risk of cancer | Discharge from industrial chemical factories |
| 1,3- butadiene | 1,200** | NA | NA |
| 1, 3-dichlorobenzene | NE | NA | NA |
| 1, 3-dichloropropane | 730** | NA | NA |
| 1, 4-dichlorobenzene | 9.1** | NA | NA |
| 1, 4-dioxane | 75** | NA | NA |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water[^] | Sources of Contaminant in Drinking Water[^] |
|---|--|--|--|
| 2, 2-dichloropropane | 9.1** | NA | NA |
| 2- chloro-1,3- butadiene | 13 | NA | NA |
| 2-chlorotoluene | NE | NA | NA |
| 2-hexanone | 490** | NA | NA |
| 2-nitropropane | 120** | NA | NA |
| 1,3,5- trichlorobenzene | 3.4** | NA | NA |
| 3- chloro-1- propene | 73** | NA | NA |
| 4-chlorotoluene | NE | NA | NA |
| 4-isopropyltoluene | 490** | NA | NA |
| 4-methyl-2-pentanone (MIBK) | 2,400** | NA | NA |
| Acetone | 1,950** | NA | NA |
| Acetonitrile | 22,000** | NA | NA |
| Benzene | 780** | NA | NA |
| Benzyl chloride | 5 | Anemia, decrease in blood platelets, increased risk of cancer | Discharge from factories, leaching from gas storage tanks and landfills |
| Bromobenzene | 5.4** | NA | NA |
| Bromochloromethane (chlorobromomethane) | 200** | NA | NA |
| Bromodichloromethane | 980** | NA | NA |
| Bromoform (Tribromomethane) | 15** | NA | NA |
| Bromomethane (methyl bromide) | 120** | NA | NA |
| Carbon disulfide | 34** | NA | NA |
| Carbon tetrachloride | 2,400** | NA | NA |
| Chlorobenzene | 5 | Liver problems, increased risk of cancer | Discharge from chemical plants and other industrial activities |
| Chloroethane (ethyl chloride) | 100 | Liver or kidney problems | Discharge from chemical and agricultural chemical factories |
| Chloroform | 9,800** | NA | NA |
| Chloromethane (methyl chloride) | 240** | NA | NA |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water[^] | Sources of Contaminant in Drinking Water[^] |
|--------------------------------------|--|---|---|
| Cis-1, 2-dichloroethene | 70** | NA | NA |
| Cis-1, 3-dichloropropene | 70 | NA | NA |
| Cis-1,4- dichloro-2- butene | 2.0** | NA | NA |
| Cyclohexane | NE | NA | NA |
| Cyclohexanone | 120,000** | NA | NA |
| Dibromochloromethane | 120,000** | NA | NA |
| Dibromomethane | 11** | NA | NA |
| Dichlorodifluoromethane | NE | NA | NA |
| Ethylbenzene | 4,900** | NA | NA |
| Ethyl acetate | 700** | Liver or kidney problems | Discharge from petroleum refineries |
| Ethyl ether | 22,000** | NA | NA |
| Ethylene oxide | 4900** | NA | NA |
| Ethyl methacrylate | 0.89** | NA | NA |
| Hexane | 2,200** | NA | NA |
| Hexachlorobutadiene | 1,500** | NA | NA |
| Iodomethane | 12** | NA | NA |
| Isobutyl alcohol | 34** | NA | NA |
| Isooctane | 7,300** | NA | NA |
| Isopropylbenzene (cumene) | NE | NA | NA |
| Methacrylonitrile | 700 / 2,400** | NA | NA |
| Methyl ethyl ketone (2-butanone) | 2.4** | NA | NA |
| Methyl methacrylate | 15,000** | NA | NA |
| Methylene chloride (dichloromethane) | 34,000** | NA | NA |
| Naphthalene | 5** | NA | NA |
| n-Butylbenzene | 490** | NA | NA |
| n-Heptane | 1,200** | NA | NA |
| n-Propylbenzene | 1,500** | NA | NA |
| Pentachloroethane | 980** | NA | NA |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water[^] | Sources of Contaminant in Drinking Water[^] |
|---|--|---|---|
| Propionitrile | 10** | NA | NA |
| sec-Butylbenzene | 9.8** | NA | NA |
| Styrene | 980** | NA | NA |
| tert-Butylbenzene | 100 | NA | NA |
| Tert-butyl methyl ether (mtbe) | 980** | NA | NA |
| Tetrachloroethene | 240** | NA | NA |
| Toluene | 5 | NA | NA |
| Trans-1, 2-dichloroethene | 1,000 | Nervous system, kidney, or liver problems | Discharge from petroleum factories |
| Trans-1, 3-dichloropropene | 100 | NA | NA |
| Trans-1,4- dicloro-2-butene | 9.1** | NA | NA |
| Trichloroethene | NE | NA | NA |
| Trichlorofluoromethane | 5 | NA | NA |
| Vinyl Acetate | 7,300** | NA | NA |
| Vinyl chloride (chloroethene) | 24,000** | NA | NA |
| m-p-xylene | 2 | Increased risk of cancer | Leaching from PVC pipes, discharge from plastic factories |
| o-xylene | 10,000** | NA | NA |
| Xylenes, Total | 10,000** | NA | NA |
| | 10,000** | Nervous system damage | Discharge from petroleum factories, discharge from chemical factories |
| Total coliforms (including E. Coli MPN) | | | |
| 1694 Pharmaceuticals (LCMS/MS) | 0 | Not a health threat in itself; it is used to indicate whether other potentially harmful bacteria may be present. | Coli forms are naturally present in the environment, as well as feces; fecal coli forms and E. coli only come from human and animal fecal waste. |
| 1694 Pharmaceuticals (LCMS/MS) | NA | NA | NA |
| 1694 Pharmaceuticals (LCMS/MS) | NA | NA | NA |
| 1694 Pharmaceuticals (LCMS/MS) | NA | NA | NA |

| Parameter, Method, and Units | Maximum Contaminant Levels or Secondary Standards | Potential Health Effects from Ingestion of Water[^] | Sources of Contaminant in Drinking Water[^] |
|-------------------------------------|--|---|---|
| Turbidity | NA | NA | NA |
| | NA | Turbidity is a measure of the cloudiness of water. It is used to indicate water quality and filtration effectiveness (e.g., whether disease-causing organisms are present). Higher turbidity levels are often associated with higher levels of disease-causing microorganisms such as viruses, parasites, and some bacteria. These organisms can cause symptoms such as nausea, cramps, diarrhea, and associated headaches. | Soil runoff |

Maximum contaminant level and secondary standards from 30 TAC 290 Subchapter F.

** Numerical value for risk reduction not an MCL, but provides a measure of desirable concentrations, from RG-346 (www.sos.state.tx.us).

[^]From EPA 816-F-02-013 July 2002.

APPENDIX F—Stormwater-Sampling Detail

Stormwater-Sampling Program for Comal and San Marcos Springs in Support of the Edwards Aquifer Habitat Conservation Plan

PURPOSE

The purpose of this technical procedure is to describe the methodology for collecting grab samples from stormwater runoff in surface waters at Comal and San Marcos springs. Sample frequency is twice annually, with samples collected across three points on the hydrograph. The EAA samples storm waters at Comal Springs at the following five locations (see Appendix A for map):

1. Upper Springs (near Blieders Creek),
2. New Channel—(below confluence with Dry Comal Creek),
3. Upper Old Channel—(at Elizabeth Street),
4. Lower Old Channel—(above Hinman Island), and
5. Comal River—(above confluence with Guadalupe River).

The EAA samples stormwaters at San Marcos Springs at the following seven locations (see Appendix A for map):

1. Sink Creek, upstream of Spring Lake,
2. Sessoms Creek,
3. Dog Beach Outflow,
4. Hopkins Street Outflow,
5. Purgatory Creek (above San Marcos River),
6. I-35 Reach, and
7. Willow Creek (above San Marcos River).

SCOPE

This procedure applies to all EAA personnel and subcontractors who sample storm water.

DEFINITIONS

1. Stormwater runoff as stated by the US EPA, “is generated when precipitation from rain and snowmelt events flows over land or impervious surfaces and does not percolate into the ground” (US EPA Stormwater Program, epa.gov).
2. Rivers are sources of water that flow on top of the ground in volume.
3. Sample intervals (for the EAHCP stormwater sampling program) are defined as:
 - a. Initial rise, or rising limb of the hydrograph;
 - b. Peak area of hydrograph; and
 - c. Recession limb of the hydrograph.

GENERAL

Weather permitting, EAA will sample two stormwater events per year to evaluate stormwater quality from urban landscapes that discharge to Comal and San Marcos springs.

STORM-EVENT SELECTION CRITERIA

According to the *Water Quality Monitoring Program Strategy for Comal Springs and San Marcos Springs in Support of the Edwards Aquifer Habitat Conservation Plan*, (EAHCP Workplan “a storm water sampling event will be triggered when a local rainfall event causes a significant increase in spring flow at the historic Comal Springs gauging station and the San Marcos Springs gauging station.” Furthermore, data collected from real-time instrumentation for surface water quality will be used to further refine the type of stormwater event(s) to be sampled. Real-time data are collected for the following parameters at 15-minute intervals from the stations shown on Comal and San Marcos springs EAHCP maps (Appendix A):

- Conductivity,
- DO,
- pH,
- Temperature, and
- Turbidity.

EAA field staff will monitor incoming storms by radar to determine whether the storm will produce one-half inch or more of localized precipitation and determine whether the storm is safe for stormwater sampling. Because of the nature of storms, stormwater sampling may be canceled as a result of false starts, safety issues, or if a new storm interrupts the stormwater sampling. Aquifer Science Management will make the final determination regarding go/no go for stormwater sampling.

Minimum Antecedent Dry Period Requirements

The following is a guideline to determine whether watersheds have returned to “normal” flow conditions. Each watershed will be evaluated separately because one watershed may return to “normal” flow conditions faster and technically be ready for another stormwater sampling event before another watershed, as noted below:

- One day wait if the previous rain event was limited to light rain/drizzle, producing only a surface wetting and no runoff
- Three days wait if the previous rain event did not produce enough rainfall to result in a measurable increase in discharge at the sample location(s)

- Minimum of five days wait if preceded by a rainfall of at least one-half inch at a sample location. The antecedent dry period may be longer if the sample location(s) are still being impacted by runoff from a previous rain event (SARA, 2013).

Canceling a Stormwater-Sampling Event

- A stormwater-sampling event may be canceled because of excessive lightning, hail, high winds, or flooding. If a storm does become severe during a stormwater-sampling event, the event will be postponed, cancelled, or suspended under some circumstances.
- A stormwater sampling event may be suspended because of a new rain event. For example, if samples are collected during the 10% of baseline flow conditions and another storm event interrupts this sampling event, then sampling will be suspended. The second storm will represent a new stormwater event.

RESPONSIBILITIES

CTO and Hydrogeologist Supervisor—Aquifer Science

The CTO and hydrogeologist supervisor—aquifer science will determine which parameters need to be sampled and will ensure that the samples obtained represent the environment being investigated. Sampling parameters are listed in the EAHCP workplan.

Hydrologic Data Coordinator

The hydrologic data coordinator will schedule sampling events and ensure that all field crews are provided with the information and equipment necessary to successfully complete scheduled sampling (i.e., location ID and selected analyses). Furthermore, the coordinator will organize and interface with local entities as needed to ensure that all notifications are in place for each river/spring complex as needed.

Environmental Science Technicians

Environmental science technicians will generally be responsible for collection of samples. Other individuals may also be asked to participate in sample-collection activities. However, each sample team of two people will have a lead sampler who reports back to the hydrogeologist supervisor—aquifer science. Reports will include

problems and issues in the field, inability to sample because of unforeseen or changing circumstances, and any deviations from the sample-collection plan and protocols.

PROCEDURE

Supplies and Equipment

Major Equipment Items

- Sample dipper
- Peristaltic pump with inert sample tubing
- 500- or 1,000-mL Teflon™ beakers affixed to telescoping rods
- Two gallon buckets for field-parameter readings

Equipment Support Items

- Trash bags
- Gloves (nitrile)
- Kim wipes/towels
- Rope
- Garden wagon

Sampling Supplies

- Sample bottles
- COC forms
- Sample labels
- Bailer (for filtration)
- 0.45-micron filter
- Ice chest
- Ice for sample preservation
- Ziplock bags
- Field sheet
- Pen and waterproof permanent marker

Monitoring Equipment

- pH and temperature meter
- Specific conductance meter
- Dissolved-oxygen meter
- Turbidity meter

Health and Safety Items

- First-aid kit and emergency eye-wash kit
- Fire extinguisher
- Mobile phone
- Helmet with head lamp
- Hand sanitizer
- Mud boots
- Raincoat
- Life vests with reflective markings
- Throw rope
- Computer access to real-time flow, water quality, and weather data

Field Equipment Decontamination

Proper decontamination between sites is essential to the avoidance of introducing contaminants from the sampling equipment. Before sampling, all hoses, buckets, water quality probes, and other sampling equipment should be decontaminated at EAA before fieldwork. Procedures specified in the EAA's *Field Sampling Plan* should be followed for decontamination of field equipment.

Instrument Usage and Measurement of Water Quality Parameters

Before going into the field, the environmental science technician should verify that all field instruments are operating properly. Calibration will be done on pH, specific conductivity, dissolved oxygen, and turbidity meters, and calibration information should be recorded in the calibration log book.

Purging

No purging is required for stormwater runoff to be sampled in the Comal and San Marcos rivers.

Sample Collection

According to the EAHCP work plan, “three water quality samples will be collected from each surface water sampling location during the sampling event. Sample times will be spaced to reflect changes in the stream hydrograph.” The first sample will be during the initial rise in the hydrograph. The second sample will be collected near the peak of flow. The final sample will be collected along the recession limb of the storm hydrograph. In some circumstances, additional samples may be collected during the storm event such

that sample groups may be subsequently compared to hydrograph data and the most representative samples groups sent for analyses. Following is the sampling procedure:

At EAA Offices

- EAA field staff will monitor local weather forecasts and Doppler radars to determine whether an incoming storm meets the criteria for a stormwater sampling event.
- If the incoming storm DOES NOT meet the criteria, no action will be taken.
- If the incoming storm DOES meet the criteria, EAA field staff will monitor weather conditions, estimate a time of arrival of the incoming storm, and determine whether weather conditions are safe for stormwater sampling (CTO or hydrogeologist supervisor will make the final go/no go decision).
- EAA field staff will notify the contracted laboratories for the possibility of samples.
- Labels for the sample bottles will be filled out.
- Aquifer Science CTO or Hydrogeologist supervisor will make the final determination regarding go/no-go with regard to the storm event.

In the Field

- Field personnel must wear clean (disposable) nitrile gloves during the sample-collection process.
- Sample water will be collected in a two-gallon bucket for parameter readings, and sample water will be collected in a 500- or 1000-mL Teflon™ beaker attached to telescoping rods, or, if needed, a peristaltic pump with inert tubing will be used.
- Meter(s) will be inserted into a two-gallon bucket and measurements recorded on a field sheet, or, if a peristaltic pump is being used, a flow chamber will be used.
- Samples will be collected using beakers or a peristaltic pump.
 - Herbicides and pesticides
 - General water quality parameters
 - Selected metals
 - Turbidity
 - Bacteria (E-coli most probable number)
 - Total phosphorous
 - Total organic carbon
 - Dissolved organic carbon
 - Total kjeldahl nitrogen
 - All containers will be filled almost full, except for alkalinity and VOCs
 - Alkalinity

- A bailer attached with a filter will be used or a filter will be attached onto tubing from the peristaltic pump
 - Alkalinity must have no head space.
- Selected metals
 - A bailer attached with a filter will be used or a filter will be attached onto tubing from the peristaltic pump
- VOC
 - The VOC sample vial will be completely filled so that the water forms a convex meniscus at the top and then capped so that no air space exists in the vial. The vial must be turned over and tapped to check for bubbles in the vial, which indicate trapped air. If bubbles are observed, the vial should be discarded and another sample collected.
- Any required information will be recorded on the field sheet before, during, and after sampling. Parameter readings will be measured in a two-gallon bucket and recorded on field sheets.
- Preservatives (if any) will be placed in the bottles by EAA-contracted laboratories.
- After the samples have been collected, they will be immediately placed in an ice-filled cooler.
- Prior to departure from the field, field documentation, including the COC form, will be completed, and all EAA field employees will clean their hands with hand sanitizer.
- Field notebooks will be used to record basic information for each event, such as magnitude of storm, issues related to sample collection, weather conditions, time of day samples were collected, and other information deemed pertinent by the lead sampler and/or coordinator.

The second sample will be collected near the peak of flow and will follow the same procedure as that of the initial rise on the hydrograph sample. The third sample will be collected along the recession limb of the hydrograph and will follow the same procedure as that of the other two sampling events. Again, the possibility exists that additional sample may be collected during the event with the most representative three sample groups being submitted for analyses (based on comparison with the appropriate stream hydrograph).

Contracted Laboratories

EAA field staff will drop off samples at EAA-contracted laboratories or have samples picked up at the EAA offices. Samples will be analyzed within proper holding times.

Equipment Blanks

Equipment blanks consist of ASTM II, reagent-grade water poured over/through any sampling equipment used for collection of definitive samples. Most sample-collection equipment is disposable; however, in some cases, an equipment blank may be required. Equipment blanks are used to assess the effectiveness of decontamination procedures (for new materials provided to the EAA or from EAA's decontamination processes) and are designated as *EB* on the COC. The frequency of collection of equipment blanks will depend on the sampling routine and sampling equipment in use. Collection of equipment blanks will be designated prior to sample-collection events.

Trip Blanks

Trip blanks are used to assess potential volatile organic contamination during sample custody in the field and shipment to the receiving laboratory. Trip blanks are submitted with characteristic samples to the laboratory to verify that volatile organic contamination has not occurred from outside influences during sample handling to transport (such as absorption through the septa.)

Trip blanks consist of two 40-mL vials filled with ASTM Type II reagent-grade water prepared by the contracted laboratory. Trip blanks will remain unopened until they are received at the contracted laboratory.

Sample Identification, Handling, and Documentation

Samples will be identified, handled, and recorded as described in the preceding sections of this document.

Records

Field sheets and COCs will be kept in a bound field log book. The following will be recorded using waterproof ink on these sheets and in the field notebook:

- Names of sampling personnel
- Weather conditions
- Project name
- Date and time of sampling
- Analyses to be performed by EAA-contracted laboratory
- Equipment-calibration information
- Field-parameter measurements
- Irregularities, problems, or delays

APPENDIX G—Equipment-Decontamination Procedures

Decontamination

Proper decontamination of all equipment used in the sample-collection process is essential to obtaining quality, representative samples. Improperly decontaminated equipment is capable of causing cross-contamination between sample sites, resulting in samples that are not representative of in situ site conditions. The objective of this appendix is to provide a set of decontamination procedures applicable to various EAA equipment and sampling programs.

Whereas many different protocols exist for decontamination, ASTM Standard D 5088 is perhaps the most commonly referenced protocol. The methods outlined here are tailored to EAA sampling environments and programs.

Basic Decontamination Procedure—Groundwater, Surface Water, and Spring Sampling Equipment

When possible, equipment that comes into contact with sample media will be single-use (disposable) equipment or dedicated equipment. Having such equipment helps reduce the possibility of cross-contamination of samples. However, for many sample types, such dedicated equipment may not be possible. As such, a listing of equipment that may be used to collect a water sample (groundwater, surface water, or spring) would include

- Grundfos submersible pump and associated pump tubing
- Peristaltic pump tubing
- Sample dippers
- Surface water churn

Other equipment that *may* come into direct contact with sample media of concern includes

- Water level measurement devices (steel tape and e-lines)
- Field-parameter probes
- Downhole geophysical equipment

Equipment that will have direct contact with any sample media will be decontaminated prior to use for sample collection or prior to introduction into the well, surface water site, or spring vent, as applicable.

Grundfos Submersible Pumps

Decontamination will be accomplished as follows for submersible well pumps. Sampler will wear new, disposable, nitrile (or equivalent) gloves to perform the decontamination.

Materials needed:

- Submersible pump, pump controller, and pump tubing
- 33-gallon trashcan (dedicated for decon use only)
- Alconox® or laboratory-grade soap
- DI water
- Large plastic bags or foil
- Plastic sheeting
- Clean scrub brush(es)

The designated trashcan will be rinsed with fresh, potable water and subsequently filled with potable water and laboratory-grade soap (per soap label directions).

When the container is approximately 80% full, the pump will be lowered, with heat shield attached, into the trashcan. The pump should be suspended at least six inches off the bottom of the trashcan. The pump will then be activated and allowed to discharge outside of the trashcan for at least 30 seconds. After the initial discharge, pump tubing will be directed into the trashcan such that the decontamination mixture is recirculated through the pump and tubing. The pump should run/recirculate a minimum of ten pump-tubing volumes (about 40 gallons) through the system. This process should take about 15 to 20 minutes.

Note: in the event that the pump or tubing has sediment or other foreign matter on it, a step will be added. A clean scrub brush will be used to remove any sediment or other foreign matter from the equipment manually prior to the circulation process.

Next, the decontamination mixture will be allowed to pump out of the trashcan into the sink (the pump should not be allowed to run dry or cavitate). The pump and tubing will be placed on a clean surface (plastic sheet) and the trashcan rinsed in clean water. The pump will be rinsed and placed back into the trashcan. The pump is to be allowed to discharge outside of the trashcan until the soapy water is evacuated from the tubing. The discharge tubing will then be placed back into the trashcan and more clean water added if needed. The freshwater will be recirculated through the pump and into the trashcan for a minimum of ten volumes (about 40 gallons). Once circulation is complete, the pump will be allowed to discharge outside the trashcan until nearly empty (again, the pump should not be allowed to run dry or cavitate). Next, a final rinse of DI water will be provided on the pump and tubing, an adequate volume being used to ensure that the pump and tubing are well rinsed.

Upon completion of the decontamination procedure, the pump will be sealed in a clean plastic bag, and the end of the pump tubing will be sealed in its own clean plastic bag. A

rubber band can be used to affix the bags around the apparatus. Pump and hose assembly are to be stored indoors when not in use, away from any sources of cross-contamination.

Tubing Decontamination for Peristaltic Pumps:

Decontamination should be accomplished as follows for peristaltic pump tubing. Sampler will wear new, disposable, nitrile (or equivalent) gloves to perform the decontamination.

Materials needed:

- Four five- or seven-gallon plastic buckets (for decon use only)
- Alconox® or laboratory-grade soap
- DI water
- Large plastic bags that can be sealed (large zip-top bags)
- Plastic sheeting
- Clean scrub brush(s)

The designated buckets will be rinsed in fresh, potable water. The first bucket will be subsequently filled with potable water and laboratory-grade soap (per soap label directions). The next two buckets will be filled with clean tap water. All three decontamination buckets are to be placed on top of a clean sheet of plastic sufficiently long to provide a clean surface on which all decontamination can take place. Decon buckets are to be placed in order on the sheet, with the soap bucket first, followed by the two rinse buckets. Decontamination should proceed such that each step is always followed in order from most contaminated to least contaminated (i.e., from prewash if needed, to soap–water mixture, to first rinse bucket, to second rinse bucket, to final DI water rinse).

Any excess foreign material will be removed from the tubing, first by wiping or scrubbing with soap and water mixture (if needed). The suction side of the tubing will be lowered into the soap–water bucket. The pump will be activated and allowed to discharge outside of the bucket until the soap–water mixture has initially purged the tubing. After the initial discharge, the pump tubing will be directed into the bucket such that the decontamination mixture is recirculated through the tubing. The pump will be allowed to run a minimum of ten pump-tubing volumes through the system (or about eight to ten gallons).

Next, the suction end of the tubing will be placed into the first rinse bucket and the pump allowed to discharge into the soap bucket until the soapy water is evacuated from the tubing. The discharge side of the tubing will then be placed back into the first rinse bucket. The freshwater will be allowed to recirculate through the pump and into the first rinse bucket for a minimum of ten volumes (or about eight to ten gallons). Once circulation is complete, the process will be repeated using the second rinse bucket. Final rinse is to be accomplished by pumping/recirculating DI water through the tubing for a minimum of ten volumes, using the third rinse bucket filled with DI water. Next, a final

rinse of DI water will be provided on the outside of the tubing using an adequate volume to ensure that the tubing is well rinsed. This final rinse will complete the decontamination process.

Upon completion of the decontamination procedure, the tubing will be allowed to dry and the tubing seal placed in a plastic bag to prevent exposure to cross-contamination. Bagged tubing is to be stored indoors when not in use away from any sources of cross-contamination.

Note: peristaltic tubing for EAHCP samples is dedicated tubing and is to be stored in *labeled* bags. The bag label will have the name of the sample point written on the outside of it. EAHCP-related tubing is not to be used for any other applications.

Decontamination of Other Equipment Used in Collection of Water or Soil Samples

Decontamination will be accomplished as follows for other equipment that will come into direct contact with sample media (dippers, churns, sample probes—if placed into sample media, water level measurement devices, soil sampling devices, or trowels). Sampler will wear new, disposable, nitrile (or equivalent) gloves to perform the decontamination.

Materials needed:

- Sample-collection device (dipper, churn, etc.) or field meter (applies only to that part of the probe exposed to sample media) or water level measurement device
- Three five- or seven-gallon plastic buckets (for decon use only)
- Alconox® or laboratory-grade soap
- DI water
- Large plastic bags or foil
- Plastic sheeting
- Clean scrub brush

Designated buckets will be rinsed in fresh, potable water. The first bucket will be subsequently filled with potable water and laboratory-grade soap (per soap label directions). The remaining two buckets will be filled with clean tap water. All three decontamination buckets are to be placed on top of a clean sheet of plastic sufficiently long to provide a clean surface on which all decontamination will take place. Decon buckets are to be placed in order on the sheet, with the soap bucket first, followed by the two rinse buckets. Decontamination will proceed such that each step is always followed in order from most contaminated to least contaminated (i.e., from prewash if needed, to soap–water mixture, to first rinse bucket, to second rinse bucket, to final DI water rinse).

Any excess sediment or foreign matter will be removed from the device by gentle scrubbing and rinsing with water prior to placement into the soap–water mixture. The

sampling device will be placed into the soap–water mixture and gently scrubbed (all surfaces that will come into contact with sample media must be cleaned).

Note: the surface water churn may not fit in the bucket(s), as such the churn may be cleaned in the 33-gallon trashcan, or it may be cleaned by some of the soap–water mixture being poured into the churn. The churn will be cleaned with the soap–water mixture; double rinsed in clean, potable water; and provided a final rinse in DI water.

Upon completion of the soap–water wash, each device being decontaminated must be double rinsed (i.e., buckets two and three) in clean, potable water, followed by a final rinse in DI water. Upon completion of decontamination, equipment will be allowed to dry and stored such that it is not exposed to potential contaminants. Equipment should be stored in plastic bags or wrapped in foil to further insulate it from potential contamination.

Note: decontamination buckets are to be monitored when used for multiple items to ensure that the soap–water mixture does not become spent or ineffective. They are to be replaced as needed. Also, rinse water should be replaced regularly when it appears to have a significant accumulation of soap.

Special Decontamination Procedures

Downhole or soil-sampling equipment may be decontaminated generally by one of the applicable processes outlined above. However, in rare cases, a tool or device that is not disposable may be exposed to hydrocarbon residue or, in rarer cases, high concentrations of heavy metals may occur. In such a scenario, the tool may (at the discretion of management) require a more elaborate decontamination procedure.

Exposure to Hydrocarbons

In the event that a tool is exposed to free-product hydrocarbons, an additional step in the decontamination process may be required that will involve spraying the tool with pesticide-grade methanol or hexane prior to the final DI water rinse. Use of solvents in this case serves to remove any hydrocarbon residual from the tool.

Exposure to Heavy Metals

In the event that a tool or device is exposed to heavy metals, and the sample media are being analyzed for these same metals, another step in the decontamination process may be required. In this case, the tool may require a spray rinse with dilute (10%) hydrochloric or nitric acid prior to DI water rinse. Use of acid in this situation will act to remove residual metals from the tool.

Note: use of solvents or acids is only to be pursued if directed by management. Use of these products can be hazardous and can also present issues regarding disposal of the waste products themselves. Use of the products may also damage sampling equipment in

some cases. In the vast majority of cases, the standard washing and rinsing procedures described herein are adequate for proper decontamination of sampling equipment. Analysis of equipment blanks will be pursued when needed so that the decontamination process might be assessed. It is the responsibility of the sampler to notify management if a tool is suspected of any unusual exposure

APPENDIX E

PHOTOGRAPHIC LOG

PHOTOGRAPHIC LOG

Comal and San Marcos springs, Texas



PHOTOGRAPH 1.
Surface water grab sample collection, March 16, 2015;
sample location HCS160.



PHOTOGRAPH 2.
Field filtering surface water grab samples, March 24, 2015;
sample location HSM150.

PHOTOGRAPHIC LOG

Comal and San Marcos springs, Texas



PHOTOGRAPH 3.
Stormwater sample collection, January 23, 2015;
sample location HCS240.



PHOTOGRAPH 4.
Alkalinity analysis of stormwater samples, January 23, 2015.

PHOTOGRAPHIC LOG

Comal and San Marcos springs, Texas



PHOTOGRAPH 5.
Sediment sample collected with core sampler, June 4, 2015;
sample location HCS330.



PHOTOGRAPH 6.
Sediment sample collected with trowel, June 5, 2015;
sample location HSM310.

PHOTOGRAPHIC LOG

Comal and San Marcos springs, Texas



PHOTOGRAPH 7.
Sediment sample collected with trowel, June 5, 2015;
sample location HSM320.



PHOTOGRAPH 8.
PDS deployment device in water, April 2015;
sample location HSM440.

PHOTOGRAPHIC LOG

Comal and San Marcos springs, Texas



PHOTOGRAPH 9.

**Removing sampler from PDS deployment device upon retrieval,
April 21, 2015; sample location HCS420.**



PHOTOGRAPH 10.

**Installing PDS deployment device, April 7, 2015;
sample location HSM410.**

APPENDIX F

RECORD OF STORMWATER SAMPLING

January 2015 – Necessary supplies and equipment were restocked and assembled.

January 2015 (and prior to each subsequent event) – Sample kits, labels and Chain-of-Custody (COC) forms were received from the contract laboratory and sampling containers, and coolers were labeled by SWCA staff. Staff went on standby for sampling events.

May 11, 2015 – SWCA contacted the EAA regarding the timeline for stormwater sampling. The overall goal of the program is to sample each system at the beginning and end of the year. EAA requested SWCA hold stormwater sampling until after June 1, 2015.

June 3, 2015 – EAA requested SWCA extend stormwater sampling hold until July 1, 2015.

July 1, 2015 – SWCA sampling teams returned to on-call status and began monitoring forecasts for stormwater sampling.

COMAL SPRINGS COMPLEX

January 21–23, 2015 – Staff mobilized to New Braunfels in response to a storm event forecast in the Comal Springs area. Light intermittent rain fell from approximately 17:00 to 22:30 on January 21, 2015, without causing a significant change to water quality parameters provided by the RTI data. More intense rain was forecasted for approximately 1:00 on January 22, 2015, SWCA staff contacted EAA staff Gizelle Luevano and Marcus Gary to confirm that sampling would still be valid if that rainfall occurred. EAA staff concurred that sampling later in the night would still be valid and SWCA staff continued to monitor water quality parameters on Environet and discharge readings from the USGS Gauge 08170500.

More intense rain did begin late on January 21, 2015, and SWCA staff determined that discharge and water quality parameters were beginning to change and the storm system would produce enough precipitation to trigger a sampling event. The EAA was notified and sampling began at approximately 00:20 on January 22, 2015.

The storm was slow moving and rain continued through January 23, 2015. Conductivity began decreasing at a slower rate between 02:30 and 02:45 and began to increase slightly at 03:00. It appeared the storm had reached its peak but more rain appeared on the radar so peak sampling was not initiated. Conductivity began to decrease again until 08:00, sampling teams mobilized and collected a peak sample. By 09:15 conductivity began decreasing again and discharge increased indicating the storm had not peaked. Sampling teams mobilized to collect peak sample sets again at 10:00 and 11:30 only to have discharge water quality parameters continue to change. The true peak of the storm was reached and sampled at approximately 16:15 on January 22, 2015. All samples collected during earlier peak runs were discarded and not analyzed.

SWCA staff continued to monitor water quality and collected the trail samples for the event beginning at approximately 03:00 on January 23, 2015. At that time discharge had reached 50% recovery and additional rain was beginning to move into the area that could have potentially affected the recovery of

the water quality parameters. The area received 2 to 2.5 inches of rain (NOAA 2015). Discharge increased from approximately 141 to 420 cfs at the peak of the storm (USGS 2015). Staging and alkalinity processing occurred at Heidelberg Lodges in New Braunfels, Texas.

August 20, 2015 – SWCA staff monitored weather forecasts and radar imagery during the morning and early afternoon hours of August 20, 2015. Forecasts varied widely and only briefly showed a slightly greater than 50% chance of 0.5 inch of rain in the area surrounding the Comal Springs system. Chances of precipitation did not exceed 50% in San Marcos. SWCA did not mobilize to either system. Some precipitation did occur, but real-time water quality data did not meet appropriate sampling conditions. RTI data is shown in Figure 1 below.

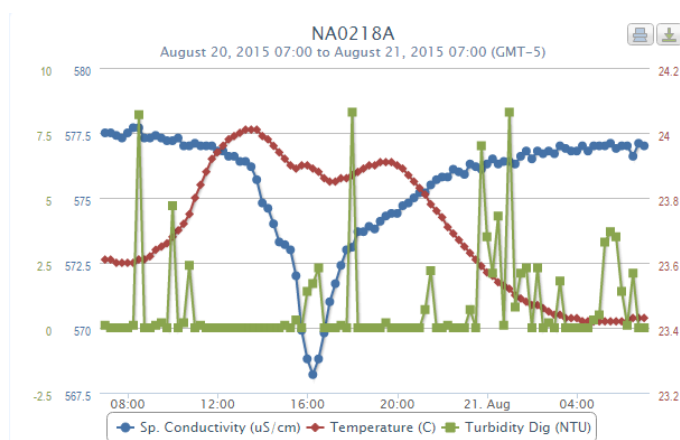


Figure 1. Real-time water quality data Aug 20, 2015.

September 10, 2015 – SWCA staff monitored weather forecasts for both Comal and San Marcos springs as some precipitation was forecasted in both areas. The forecasts were inconsistent with the amount and locations of predicted precipitation but SWCA staff anticipated potential rain events in the afternoon of September 10, 2015. Surface water collection was scheduled for the San Marcos Springs system on September 10, while sampling a short, but intense, rainfall event did occur. The surface water samples were discarded and the team returned to SWCA offices. The forecast in New Braunfels appeared very favorable for the evening hours and SWCA staff mobilized to the area. It rained for a short time and staff collected the lead sample set. Unfortunately, the rain stopped and a full event was not captured. The samples were discarded. A full event did occur in the San Marcos Springs complex but the precipitation occurring earlier in the day violated the antecedent dry period for stormwater sampling.

October 21, 2015 – SWCA staff monitored weather forecasts for both Comal and San Marcos spring complexes. Greater than 50% chance of greater than 0.5 inch of precipitation was forecasted for both spring complexes. One team was deployed to San Marcos. Sample equipment and sample kits were stationed at the San Marcos Nature Center. Insufficient rainfall occurred on October 21, 2015, but more precipitation was forecasted for early the next morning.

Forecasts were checked throughout the night and teams were informed at 02:30 to mobilize to both spring complexes by 06:00 on October 22, 2015. No precipitation occurred during the morning of October 22, 2015, and teams demobilized. More precipitation was forecasted for October 23, 2015.

Weather forecasts were monitored throughout the night and one team mobilized to each spring complex in the early morning hours. Precipitation began in New Braunfels late in the morning. As soon as the RTI data indicated a notable change in water quality from stormwater, sampling of the lead samples began. Monitoring continued and the peak and trail sample sets were collected in the afternoon and early evening of October 23, 2015.

SAN MARCOS SPRINGS COMPLEX

January 29–30, 2015 – Staff were aware of a storm forecast for the San Marcos area January 29–30, 2015. As the date approached, staff prepared for the event and continued to monitor the storm. The storm changed course and only the edge of the storm passed over San Marcos. Sampling teams did not deploy but monitored radar, forecasts and water quality parameters over the weekend. The system did not produce enough precipitation to cause a change in water quality parameters.

March 7–8, 2015 – A storm was forecasted for the San Marcos area in the early morning hours of March 8, 2015. One sample team deployed to the San Marcos Nature Center late on the evening on March 7, 2015, ahead of the anticipated storm. Rain began at approximately 4:00 on March 8, 2015. Additional sampling teams were called to the area. A lead sample collection began at approximately 5:00 when water quality parameters began to change. Unfortunately, the storm dissipated before enough precipitation fell to qualify as a full storm event. More rain was forecasted for the evening of March 8, 2015, but after consultation with EAA staff it was determined that rain later that day would constitute a separate event. The effort was cancelled and samples were disposed of.

March 17, 2015 – SWCA sampling teams monitored weather forecasts and radar and deployed to the Old Fish Hatchery building in San Marcos at approximately 17:00. The majority of the storm did not pass directly over San Marcos. Light rain fell intermittently but did not change water quality parameters. Sampling teams remained in place because radar continued to show rain moving toward the area. The sampling effort was cancelled at approximately 01:00 on March 18, 2015.

March 20, 2015 – A storm was forecasted for the San Marcos area in the afternoon of March 20, 2015. Sampling teams deployed to the area at approximately 14:30. Rain began to fall and water quality parameters began to change. Lead sample collection began at 15:45. The rain stopped but more was forecasted throughout the evening. The team remained in place overnight, ready to continue sampling. Additional rain that fell overnight was light and water quality parameters recovered before reaching levels that would qualify the event for sampling. The effort was cancelled and samples were discarded.

April 16, 2015 – SWCA sampling teams deployed to the San Marcos Nature Center in response to a storm occurring in the San Marcos area in the afternoon of April 16, 2015. Lead sample collection was initiated at approximately 17:00. Sampling continued through the evening until the final peak sample was

collected at approximately 23:45. A full event was sampled however a delay in shipping caused the samples to reach the laboratory well outside of acceptable temperature ranges and some holding times. The samples were not analyzed and were discarded.

May 5–6, 2015 – SWCA sampling teams deployed to the San Marcos Nature Center in the evening on May 5, 2015. Lead sampling was initiated at approximately 21:45 after significant rain began and water quality parameters began to change. The river responded rapidly to the storm event and sampling teams initiated peak sampling immediately following lead sample collection. Peak samples were initiated at approximately 23:15. SWCA staff transported samples back to the SWCA offices in San Antonio for processing and packaging for shipment. One sampling team remained in place in San Marcos to collect the trail samples. Trail samples were initiated at approximately 01:30 on May 6, 2015, after specific conductivity passed 50% recovery.

Filtration for the metals and alkalinity samples was performed at the SWCA San Antonio office utilizing a peristaltic pump. New disposable bailers were used to collect HSM240 Peak and HSM240 Trail samples. *E. coli* samples were delivered to the San Antonio River Authority (SARA) laboratory when it opened the morning of May 6, 2015. Due to the overnight collection of samples, it was not possible to deliver samples to the SARA laboratory within the 6-hour hold time.

October 21, 2015 – SWCA staff monitored weather forecasts for both Comal and San Marcos spring complexes. A greater than 50% chance of greater than 0.5 inch of precipitation was forecasted for both spring complexes. One team was deployed to San Marcos. Sample equipment and sample kits were stationed at the San Marcos Nature Center. Insufficient rainfall occurred on October 21, 2015, but more precipitation was forecasted for early the next morning.

Forecasts were checked throughout the night and teams were informed at 02:30 to mobilize to both spring complexes by 06:00 on October 22, 2015. No precipitation occurred during the morning of October 22, 2015, and teams demobilized. More precipitation was forecasted for October 23, 2015.

Weather forecasts were monitored throughout the night and one team mobilized to each spring complex in the early morning hours. Precipitation began in San Marcos in the early afternoon. As soon as the RTI data indicated a notable change in water quality from stormwater, sampling of the lead samples began. Monitoring continued and the peak and trail sample sets were collected in the afternoon and evening of October 23, 2015.

APPENDIX G

LABORATORY REPORTS

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HCS 110
Sample Number: AA93406
Sample Matrix: Non Potable Water

Collection Date/Time: 3/16/15 12:44
Receipt Date/Time: 3/16/15 15:16

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93406-A | E. coli | √ | SM 9223B-2004 | 4 | MPN/100 mL | | 1 | 42074 | 3/16/15 | 16:20 | RSC |
| AA93406-A | E. Coli Holding Time - IDEXX Colilert | | NA | 3.60 | hours | | 0.00 | 42073 | 3/16/15 | 16:20 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 20, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42074

QC Analyte Name

Initial Blank for E. coli

Log Range for E. coli

Result

Absent

0.0000

Units

Qualifier

Lower

0.0

Acceptance Criteria

Target

Absent

Upper

0.5



Jeanette Hernandez
Water Quality Planner / QAO

3/20/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HCS 120
Sample Number: AA93407
Sample Matrix: Non Potable Water

Collection Date/Time: 3/16/15 13:20
Receipt Date/Time: 3/16/15 15:16

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93407-A | E. coli | √ | SM 9223B-2004 | 47 | MPN/100 mL | | 1 | 42074 | 3/16/15 | 16:20 | RSC |
| AA93407-A | E. Coli Holding Time - IDEXX Colilert | | NA | 3.00 | hours | | 0.00 | 42073 | 3/16/15 | 16:20 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 20, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42074

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

3/20/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 20, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HCS 130
Sample Number: AA93408
Sample Matrix: Non Potable Water

Collection Date/Time: 3/16/15 11:29
Receipt Date/Time: 3/16/15 15:16

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93408-A | E. coli | √ | SM 9223B-2004 | 59 | MPN/100 mL | | 1 | 42074 | 3/16/15 | 16:20 | RSC |
| AA93408-A | E. Coli Holding Time - IDEXX Colilert | | NA | 4.85 | hours | | 0.00 | 42073 | 3/16/15 | 16:20 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 20, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42074

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

3/20/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HCS 140
Sample Number: AA93409
Sample Matrix: Non Potable Water

Collection Date/Time: 3/16/15 13:52
Receipt Date/Time: 3/16/15 15:16

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93409-A | E. coli | √ | SM 9223B-2004 | 80 | MPN/100 mL | | 1 | 42074 | 3/16/15 | 16:20 | RSC |
| AA93409-A | E. Coli Holding Time - IDEXX Colilert | | NA | 2.47 | hours | | 0.00 | 42073 | 3/16/15 | 16:20 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 20, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42074

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

3/20/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 20, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HCS 160
Sample Number: AA93410
Sample Matrix: Non Potable Water

Collection Date/Time: 3/16/15 14:13
Receipt Date/Time: 3/16/15 15:16

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93410-A | E. coli | √ | SM 9223B-2004 | 60 | MPN/100 mL | | 1 | 42074 | 3/16/15 | 16:20 | RSC |
| AA93410-A | E. Coli Holding Time - IDEXX Colilert | | NA | 2.12 | hours | | 0.00 | 42073 | 3/16/15 | 16:20 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 20, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42074

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

3/20/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: FD HCS 120
Sample Number: AA93411
Sample Matrix: Non Potable Water

Collection Date/Time: 3/16/15 13:20
Receipt Date/Time: 3/16/15 15:16

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93411-A | E. coli | √ | SM 9223B-2004 | 51 | MPN/100 mL | | 1 | 42074 | 3/16/15 | 16:20 | RSC |
| AA93411-A | E. Coli Holding Time - IDEXX Colilert | | NA | 3.00 | hours | | 0.00 | 42073 | 3/16/15 | 16:20 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 20, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42074

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

3/20/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable



WORK ORDER NUMBER: 15-03-1281

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: SWCA Environmental Consultants

Client Project Name: EAA 27122

Attention: Philip Pearce
6200 UTSA Blvd.
Suite 102
San Antonio, TX 78249-1618

A handwritten signature in black ink, appearing to read "Don Burley".

Approved for release on 03/31/2015 by:
Don Burley
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

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Client Project Name: EAA 27122
Work Order Number: 15-03-1281

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Work Order Narrative

Work Order: 15-03-1281

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Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 03/17/15. They were assigned to Work Order 15-03-1281.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.



Calscience

Sample Summary

| | | | |
|---------|--------------------------------|-----------------------|----------------|
| Client: | SWCA Environmental Consultants | Work Order: | 15-03-1281 |
| | 6200 UTSA Blvd., Suite 102 | Project Name: | EAA 27122 |
| | San Antonio, TX 78249-1618 | PO Number: | |
| | | Date/Time Received: | 03/17/15 10:00 |
| | | Number of Containers: | 63 |
| Attn: | Philip Pearce | | |

| Sample Identification | Lab Number | Collection Date and Time | Number of Containers | Matrix |
|-----------------------|--------------|--------------------------|----------------------|---------|
| HCS 110 | 15-03-1281-1 | 03/16/15 12:44 | 9 | Aqueous |
| HCS 120 | 15-03-1281-2 | 03/16/15 13:20 | 9 | Aqueous |
| HCS 130 | 15-03-1281-3 | 03/16/15 11:29 | 9 | Aqueous |
| HCS 140 | 15-03-1281-4 | 03/16/15 13:52 | 9 | Aqueous |
| HCS 160 | 15-03-1281-5 | 03/16/15 14:13 | 9 | Aqueous |
| FDHCS 120 | 15-03-1281-6 | 03/16/15 13:20 | 9 | Aqueous |
| TB02 | 15-03-1281-7 | 03/16/15 00:00 | 9 | Aqueous |


Return to Contents



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

Page 1 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 110 | 15-03-1281-1-F | 03/16/15 12:44 | Aqueous | ICP 7300 | 03/30/15 | 03/30/15 14:56 | 150330LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 64.1 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 16.8 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.39 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.678 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 0.655 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS 120 | 15-03-1281-2-F | 03/16/15 13:20 | Aqueous | ICP 7300 | 03/30/15 | 03/30/15 14:58 | 150330LA2F |
|---------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 87.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 16.3 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.65 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.690 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.44 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS 130 | 15-03-1281-3-F | 03/16/15 11:29 | Aqueous | ICP 7300 | 03/30/15 | 03/30/15 15:00 | 150330LA2F |
|---------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 84.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 16.4 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.63 | 0.500 | 0.103 | 1.00 | |
| Sodium | 14.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.696 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.35 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

Page 2 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 140 | 15-03-1281-4-F | 03/16/15 13:52 | Aqueous | ICP 7300 | 03/30/15 | 03/30/15 15:02 | 150330LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 85.1 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 16.6 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.68 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.7 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.708 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.47 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS 160 | 15-03-1281-5-F | 03/16/15 14:13 | Aqueous | ICP 7300 | 03/30/15 | 03/30/15 15:05 | 150330LA2F |
|---------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 86.7 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.0 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.71 | 0.500 | 0.103 | 1.00 | |
| Sodium | 14.5 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.722 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.48 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|-----------|----------------|----------------|---------|----------|----------|----------------|------------|
| FDHCS 120 | 15-03-1281-6-F | 03/16/15 13:20 | Aqueous | ICP 7300 | 03/30/15 | 03/30/15 15:07 | 150330LA2F |
|-----------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 85.6 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 16.2 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.67 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.0 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.668 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.44 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-683-1193 | N/A | Aqueous | ICP 7300 | 03/30/15 | 03/30/15 17:20 | 150330LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | ND | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 0.00760 | 0.100 | 0.00336 | 1.00 | J |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 1 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 110 | 15-03-1281-1-F | 03/16/15 12:44 | Aqueous | ICP/MS 03 | 03/19/15 | 03/23/15 15:42 | 150319LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000118 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0371 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000547 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00170 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000295 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0245 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0520 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00243 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 120 | 15-03-1281-2-F | 03/16/15 13:20 | Aqueous | ICP/MS 03 | 03/19/15 | 03/23/15 15:46 | 150319LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0555 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000538 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00201 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000334 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0239 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0460 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00178 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 130 | 15-03-1281-3-F | 03/16/15 11:29 | Aqueous | ICP/MS 03 | 03/19/15 | 03/23/15 15:49 | 150319LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0574 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000570 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00184 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000347 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00384 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0468 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00459 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 140 | 15-03-1281-4-F | 03/16/15 13:52 | Aqueous | ICP/MS 03 | 03/19/15 | 03/23/15 15:53 | 150319LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|-----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0584 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000907 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | 0.0000905 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00197 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000446 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0118 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0399 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00632 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 160 | 15-03-1281-5-F | 03/16/15 14:13 | Aqueous | ICP/MS 03 | 03/19/15 | 03/23/15 15:57 | 150319LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000107 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0598 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000607 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00216 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000334 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0265 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0508 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00991 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 120 | 15-03-1281-6-F | 03/16/15 13:20 | Aqueous | ICP/MS 03 | 03/19/15 | 03/23/15 16:00 | 150319LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000166 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0562 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000426 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.000783 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | 0.000117 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00225 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000414 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0391 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00459 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0555 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00607 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-769 | N/A | Aqueous | ICP/MS 03 | 03/19/15 | 03/23/15 13:26 | 150319LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 110 | 15-03-1281-1-F | 03/16/15 12:44 | Aqueous | Mercury 04 | 03/18/15 | 03/18/15 19:02 | 150318L07F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 120 | 15-03-1281-2-F | 03/16/15 13:20 | Aqueous | Mercury 04 | 03/18/15 | 03/18/15 19:04 | 150318L07F |
|---------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 130 | 15-03-1281-3-F | 03/16/15 11:29 | Aqueous | Mercury 04 | 03/18/15 | 03/18/15 19:06 | 150318L07F |
|---------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 140 | 15-03-1281-4-F | 03/16/15 13:52 | Aqueous | Mercury 04 | 03/18/15 | 03/18/15 19:08 | 150318L07F |
|---------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 160 | 15-03-1281-5-F | 03/16/15 14:13 | Aqueous | Mercury 04 | 03/18/15 | 03/18/15 19:15 | 150318L07F |
|---------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|-----------|----------------|----------------|---------|------------|----------|----------------|------------|
| FDHCS 120 | 15-03-1281-6-F | 03/16/15 13:20 | Aqueous | Mercury 04 | 03/18/15 | 03/18/15 19:17 | 150318L07F |
|-----------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-763-517 | N/A | Aqueous | Mercury 04 | 03/18/15 | 03/18/15 18:48 | 150318L07F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|-----------------------|-----------------------|----------------|-----------------------|-----------------|-----------------------|-------------------|
| HCS 110 | 15-03-1281-1-I | 03/16/15 12:44 | Aqueous | LC/TQ 2 | 03/19/15 | 03/23/15 15:29 | 150319L20 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 39 | | 9.7 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 45 | | 31-200 | | | |
| HCS 120 | 15-03-1281-2-I | 03/16/15 13:20 | Aqueous | LC/TQ 2 | 03/19/15 | 03/23/15 15:38 | 150319L20 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 9.8 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 44 | | 31-200 | | | |
| HCS 130 | 15-03-1281-3-I | 03/16/15 11:29 | Aqueous | LC/TQ 2 | 03/19/15 | 03/23/15 15:46 | 150319L20 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 9.8 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 43 | | 31-200 | | | |
| HCS 140 | 15-03-1281-4-I | 03/16/15 13:52 | Aqueous | LC/TQ 2 | 03/19/15 | 03/23/15 15:54 | 150319L20 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 9.9 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 43 | | 31-200 | | | |
| HCS 160 | 15-03-1281-5-I | 03/16/15 14:13 | Aqueous | LC/TQ 2 | 03/19/15 | 03/23/15 16:03 | 150319L20 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 9.7 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 43 | | 31-200 | | | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|----------------|-----------------|-----------------------|------------------|
| FDHCS 120 | 15-03-1281-6-I | 03/16/15 13:20 | Aqueous | LC/TQ 2 | 03/19/15 | 03/23/15 16:11 | 150319L20 |

| | | | | |
|------------------|---------------|-----------|-----------|-------------------|
| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | ND | 9.8 | 1.00 | |

| | | | |
|--------------------------|-----------------|-----------------------|-------------------|
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
| Caffeine-C13 (Surrogate) | 42 | 31-200 | |

| | | | | | | | |
|---------------------|----------------------|------------|----------------|----------------|-----------------|-----------------------|------------------|
| Method Blank | 099-16-376-11 | N/A | Aqueous | LC/TQ 2 | 03/19/15 | 03/23/15 15:21 | 150319L20 |
|---------------------|----------------------|------------|----------------|----------------|-----------------|-----------------------|------------------|

| | | | | |
|------------------|---------------|-----------|-----------|-------------------|
| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | ND | 10 | 1.00 | |

| | | | |
|--------------------------|-----------------|-----------------------|-------------------|
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
| Caffeine-C13 (Surrogate) | 46 | 31-200 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

Page 1 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 110 | 15-03-1281-1-I | 03/16/15 12:44 | Aqueous | GC 44 | 03/19/15 | 03/23/15 13:10 | 150319L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 93 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 89 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 120 | 15-03-1281-2-I | 03/16/15 13:20 | Aqueous | GC 44 | 03/19/15 | 03/23/15 13:24 | 150319L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 95 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 130 | 15-03-1281-3-I | 03/16/15 11:29 | Aqueous | GC 44 | 03/19/15 | 03/23/15 13:38 | 150319L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 114 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 98 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 140 | 15-03-1281-4-I | 03/16/15 13:52 | Aqueous | GC 44 | 03/19/15 | 03/23/15 13:53 | 150319L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 118 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 106 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 160 | 15-03-1281-5-I | 03/16/15 14:13 | Aqueous | GC 44 | 03/19/15 | 03/23/15 14:07 | 150319L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 101 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 90 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 120 | 15-03-1281-6-I | 03/16/15 13:20 | Aqueous | GC 44 | 03/19/15 | 03/23/15 14:36 | 150319L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 99 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 90 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-792 | N/A | Aqueous | GC 44 | 03/19/15 | 03/20/15 13:25 | 150319L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 107 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 113 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 110 | 15-03-1281-1-I | 03/16/15 12:44 | Aqueous | GC 58 | 03/19/15 | 03/20/15 19:21 | 150319L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 84 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 73 | 50-135 | |

| HCS 120 | 15-03-1281-2-I | 03/16/15 13:20 | Aqueous | GC 58 | 03/19/15 | 03/20/15 19:39 | 150319L02 |
|---------|----------------|----------------|---------|-------|----------|----------------|-----------|
|---------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 92 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 77 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

Page 2 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 130 | 15-03-1281-3-I | 03/16/15 11:29 | Aqueous | GC 58 | 03/19/15 | 03/20/15 19:57 | 150319L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 103 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 84 | 50-135 | |

| HCS 140 | 15-03-1281-4-I | 03/16/15 13:52 | Aqueous | GC 58 | 03/19/15 | 03/20/15 20:15 | 150319L02 |
|---------|----------------|----------------|---------|-------|----------|----------------|-----------|
|---------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 107 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 88 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 160 | 15-03-1281-5-I | 03/16/15 14:13 | Aqueous | GC 58 | 03/19/15 | 03/20/15 20:33 | 150319L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 93 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 77 | 50-135 | |

| FDHCS 120 | 15-03-1281-6-I | 03/16/15 13:20 | Aqueous | GC 58 | 03/19/15 | 03/20/15 20:51 | 150319L02 |
|-----------|----------------|----------------|---------|-------|----------|----------------|-----------|
|-----------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 92 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 79 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-533-1015 | N/A | Aqueous | GC 58 | 03/19/15 | 03/20/15 19:03 | 150319L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 91 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 1 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 110 | 15-03-1281-1-I | 03/16/15 12:44 | Aqueous | GC 35 | 03/19/15 | 03/20/15 13:01 | 150319L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 103 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 120 | 15-03-1281-2-I | 03/16/15 13:20 | Aqueous | GC 35 | 03/19/15 | 03/20/15 13:47 | 150319L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0051 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0051 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0051 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0051 | 0.0025 | 1.00 | |
| Diazinon | ND | 0.0051 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0051 | 0.0037 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0051 | 0.0026 | 1.00 | |
| Fensulfothion | ND | 0.0051 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0051 | 0.0027 | 1.00 | |
| Merphos | ND | 0.0051 | 0.0027 | 1.00 | |
| Methyl Parathion | ND | 0.0051 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0051 | 0.0028 | 1.00 | |
| Naled | ND | 0.041 | 0.020 | 1.00 | |
| Phorate | ND | 0.0051 | 0.0026 | 1.00 | |
| Ronnel | ND | 0.0051 | 0.0033 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0090 | 1.00 | |
| Tokuthion | ND | 0.0051 | 0.0029 | 1.00 | |
| Trichloronate | ND | 0.0051 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0051 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 98 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 3 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 130 | 15-03-1281-3-I | 03/16/15 11:29 | Aqueous | GC 35 | 03/19/15 | 03/20/15 14:33 | 150319L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 96 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 4 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 140 | 15-03-1281-4-I | 03/16/15 13:52 | Aqueous | GC 35 | 03/19/15 | 03/20/15 15:19 | 150319L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 102 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 160 | 15-03-1281-5-I | 03/16/15 14:13 | Aqueous | GC 35 | 03/19/15 | 03/20/15 16:04 | 150319L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 98 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 120 | 15-03-1281-6-I | 03/16/15 13:20 | Aqueous | GC 35 | 03/19/15 | 03/20/15 16:50 | 150319L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 94 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-90 | N/A | Aqueous | GC 35 | 03/19/15 | 03/20/15 10:44 | 150319L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 95 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

Page 1 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 110 | 15-03-1281-1-I | 03/16/15 12:44 | Aqueous | GC 40 | 03/23/15 | 03/24/15 16:43 | 150323L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPD | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 82 | 0-123 | |

| HCS 120 | 15-03-1281-2-I | 03/16/15 13:20 | Aqueous | GC 40 | 03/23/15 | 03/24/15 17:06 | 150323L02 |
|---------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|---------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPD | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 76 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

Page 2 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 130 | 15-03-1281-3-I | 03/16/15 11:29 | Aqueous | GC 40 | 03/23/15 | 03/24/15 17:29 | 150323L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 77 | 0-123 | |

| HCS 140 | 15-03-1281-4-I | 03/16/15 13:52 | Aqueous | GC 40 | 03/23/15 | 03/24/15 17:52 | 150323L02 |
|---------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|---------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 77 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

Page 3 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 160 | 15-03-1281-5-I | 03/16/15 14:13 | Aqueous | GC 40 | 03/23/15 | 03/24/15 18:15 | 150323L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 88 | 0-123 | |

| FDHCS 120 | 15-03-1281-6-I | 03/16/15 13:20 | Aqueous | GC 40 | 03/23/15 | 03/24/15 18:38 | 150323L02 |
|-----------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|-----------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 81 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-034-639 | N/A | Aqueous | GC 40 | 03/23/15 | 03/24/15 16:20 | 150323L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 118 | 0-123 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 1 of 21

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 110 | 15-03-1281-1-H | 03/16/15 12:44 | Aqueous | GC/MS TT | 03/18/15 | 03/19/15 17:57 | 150318L08 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.8 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 9.8 | 2.8 | 1.00 | |
| Aniline | ND | 9.8 | 1.5 | 1.00 | |
| Anthracene | ND | 9.8 | 3.0 | 1.00 | |
| Azobenzene | ND | 9.8 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.4 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.8 | 4.6 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.8 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.8 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.8 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.8 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.8 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.8 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.8 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.8 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.8 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.8 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.8 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.8 | 2.6 | 1.00 | |
| Chrysene | ND | 9.8 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.8 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.8 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.8 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.8 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.8 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.8 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.8 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.8 | 3.0 | 1.00 | |
| Fluorene | ND | 9.8 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.8 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.8 | 1.00 | |
| Hexachloroethane | ND | 9.8 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.8 | 2.1 | 1.00 | |
| Isophorone | ND | 9.8 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.8 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.8 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.8 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.8 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 9.8 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 9.8 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.8 | 4.6 | 1.00 | |
| Phenanthrene | ND | 9.8 | 2.9 | 1.00 | |
| Phenol | ND | 9.8 | 2.0 | 1.00 | |
| Pyrene | ND | 9.8 | 2.9 | 1.00 | |
| Pyridine | ND | 9.8 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 78 | 33-120 | |
| 2-Fluorophenol | 46 | 24-120 | |
| Nitrobenzene-d5 | 87 | 38-120 | |
| p-Terphenyl-d14 | 84 | 41-137 | |
| Phenol-d6 | 30 | 16-120 | |
| 2,4,6-Tribromophenol | 77 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 4 of 21

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 120 | 15-03-1281-2-H | 03/16/15 13:20 | Aqueous | GC/MS TT | 03/18/15 | 03/19/15 18:16 | 150318L08 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 86 | 33-120 | |
| 2-Fluorophenol | 78 | 24-120 | |
| Nitrobenzene-d5 | 97 | 38-120 | |
| p-Terphenyl-d14 | 92 | 41-137 | |
| Phenol-d6 | 57 | 16-120 | |
| 2,4,6-Tribromophenol | 92 | 27-159 | |



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 7 of 21

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 130 | 15-03-1281-3-H | 03/16/15 11:29 | Aqueous | GC/MS TT | 03/18/15 | 03/19/15 18:35 | 150318L08 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 86 | 33-120 | |
| 2-Fluorophenol | 55 | 24-120 | |
| Nitrobenzene-d5 | 92 | 38-120 | |
| p-Terphenyl-d14 | 88 | 41-137 | |
| Phenol-d6 | 35 | 16-120 | |
| 2,4,6-Tribromophenol | 87 | 27-159 | |



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 140 | 15-03-1281-4-H | 03/16/15 13:52 | Aqueous | GC/MS TT | 03/18/15 | 03/19/15 18:53 | 150318L08 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 90 | 33-120 | |
| 2-Fluorophenol | 56 | 24-120 | |
| Nitrobenzene-d5 | 96 | 38-120 | |
| p-Terphenyl-d14 | 93 | 41-137 | |
| Phenol-d6 | 36 | 16-120 | |
| 2,4,6-Tribromophenol | 90 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 160 | 15-03-1281-5-H | 03/16/15 14:13 | Aqueous | GC/MS TT | 03/18/15 | 03/19/15 19:12 | 150318L08 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 14 of 21

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 85 | 33-120 | |
| 2-Fluorophenol | 71 | 24-120 | |
| Nitrobenzene-d5 | 89 | 38-120 | |
| p-Terphenyl-d14 | 85 | 41-137 | |
| Phenol-d6 | 51 | 16-120 | |
| 2,4,6-Tribromophenol | 84 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 120 | 15-03-1281-6-H | 03/16/15 13:20 | Aqueous | GC/MS TT | 03/18/15 | 03/19/15 19:31 | 150318L08 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 88 | 33-120 | |
| 2-Fluorophenol | 74 | 24-120 | |
| Nitrobenzene-d5 | 95 | 38-120 | |
| p-Terphenyl-d14 | 92 | 41-137 | |
| Phenol-d6 | 54 | 16-120 | |
| 2,4,6-Tribromophenol | 88 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-4010 | N/A | Aqueous | GC/MS TT | 03/18/15 | 03/19/15 16:38 | 150318L08 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 88 | 33-120 | |
| 2-Fluorophenol | 59 | 24-120 | |
| Nitrobenzene-d5 | 90 | 38-120 | |
| p-Terphenyl-d14 | 86 | 41-137 | |
| Phenol-d6 | 38 | 16-120 | |
| 2,4,6-Tribromophenol | 86 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 1 of 16

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 110 | 15-03-1281-1-A | 03/16/15 12:44 | Aqueous | GC/MS BB | 03/20/15 | 03/21/15 05:25 | 150320L025 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 105 | 78-126 | |
| 1,2-Dichloroethane-d4 | 107 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 120 | 15-03-1281-2-A | 03/16/15 13:20 | Aqueous | GC/MS BB | 03/20/15 | 03/21/15 05:52 | 150320L025 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 95 | 80-120 | |
| Dibromofluoromethane | 102 | 78-126 | |
| 1,2-Dichloroethane-d4 | 106 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 130 | 15-03-1281-3-A | 03/16/15 11:29 | Aqueous | GC/MS BB | 03/20/15 | 03/21/15 03:37 | 150320L025 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 92 | 80-120 | |
| Dibromofluoromethane | 108 | 78-126 | |
| 1,2-Dichloroethane-d4 | 110 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 140 | 15-03-1281-4-A | 03/16/15 13:52 | Aqueous | GC/MS BB | 03/20/15 | 03/21/15 06:19 | 150320L025 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 104 | 78-126 | |
| 1,2-Dichloroethane-d4 | 107 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 160 | 15-03-1281-5-A | 03/16/15 14:13 | Aqueous | GC/MS BB | 03/20/15 | 03/21/15 06:46 | 150320L025 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 103 | 78-126 | |
| 1,2-Dichloroethane-d4 | 110 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 120 | 15-03-1281-6-A | 03/16/15 13:20 | Aqueous | GC/MS BB | 03/20/15 | 03/21/15 07:13 | 150320L025 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 105 | 78-126 | |
| 1,2-Dichloroethane-d4 | 112 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| TB02 | 15-03-1281-7-A | 03/16/15 00:00 | Aqueous | GC/MS BB | 03/20/15 | 03/21/15 02:43 | 150320L025 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 105 | 78-126 | |
| 1,2-Dichloroethane-d4 | 108 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 15 of 16

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-16646 | N/A | Aqueous | GC/MS BB | 03/20/15 | 03/21/15 02:16 | 150320L025 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 91 | 80-120 | |
| Dibromofluoromethane | 104 | 78-126 | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

03/17/15

Work Order:

15-03-1281

Page 1 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 110 | 15-03-1281-1 | 03/16/15 12:44 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.25 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Chloride (24) | 17 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Nitrate (as N) (24) | 0.35 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Sulfate (24) | 25 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 03/23/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 189 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 189 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Solids, Total Dissolved (24) | 295 | 1.00 | 0.820 | 1.00 | | mg/L | 03/21/15 | 03/21/15 | SM 2540 C |
| Solids, Total Suspended (24) | 43 | 1.0 | 0.83 | 1.00 | | mg/L | 03/23/15 | 03/23/15 | SM 2540 D |
| pH (24) | 7.78 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/17/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.84 | 0.50 | 0.28 | 1.00 | | mg/L | 03/20/15 | 03/20/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 8.1 | 2.5 | 1.2 | 5.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 10 | 0.50 | 0.24 | 1.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |

| | | | |
|----------------|---------------------|-----------------------|----------------|
| HCS 120 | 15-03-1281-2 | 03/16/15 13:20 | Aqueous |
|----------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.23 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Chloride (24) | 17 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.7 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Sulfate (24) | 29 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 03/23/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 235 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 235 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Solids, Total Dissolved (24) | 350 | 1.00 | 0.820 | 1.00 | | mg/L | 03/21/15 | 03/21/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 03/23/15 | 03/23/15 | SM 2540 D |
| pH (24) | 7.36 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/17/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 03/20/15 | 03/20/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 11 | 0.50 | 0.24 | 1.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.9 | 0.50 | 0.24 | 1.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

03/17/15

Work Order:

15-03-1281

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 130 | 15-03-1281-3 | 03/16/15 11:29 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.23 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Chloride (24) | 19 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Bromide (24) | 0.10 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.8 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Sulfate (24) | 34 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 03/23/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 230 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 230 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Solids, Total Dissolved (24) | 305 | 1.00 | 0.820 | 1.00 | | mg/L | 03/21/15 | 03/21/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 03/23/15 | 03/23/15 | SM 2540 D |
| pH (24) | 7.42 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/17/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 03/20/15 | 03/20/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 3.9 | 0.50 | 0.24 | 1.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.4 | 0.50 | 0.24 | 1.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |

| | | | |
|----------------|---------------------|-----------------------|----------------|
| HCS 140 | 15-03-1281-4 | 03/16/15 13:52 | Aqueous |
|----------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.20 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Chloride (24) | 18 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Bromide (24) | 0.11 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.7 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Sulfate (24) | 29 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 03/23/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 229 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 229 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Solids, Total Dissolved (24) | 330 | 1.00 | 0.820 | 1.00 | | mg/L | 03/21/15 | 03/21/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 03/23/15 | 03/23/15 | SM 2540 D |
| pH (24) | 7.68 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/17/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 03/20/15 | 03/20/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 8.2 | 0.50 | 0.24 | 1.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 5.8 | 2.5 | 1.2 | 5.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

03/17/15

Work Order:

15-03-1281

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 160 | 15-03-1281-5 | 03/16/15 14:13 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.20 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Chloride (24) | 18 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Bromide (24) | 0.14 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.7 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Sulfate (24) | 32 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 03/23/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 229 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 229 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Solids, Total Dissolved (24) | 315 | 1.00 | 0.820 | 1.00 | | mg/L | 03/21/15 | 03/21/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 03/23/15 | 03/23/15 | SM 2540 D |
| pH (24) | 7.81 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/17/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 03/20/15 | 03/20/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 6.8 | 0.50 | 0.24 | 1.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 7.6 | 0.50 | 0.24 | 1.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |

| FDHCS 120 | 15-03-1281-6 | 03/16/15 13:20 | Aqueous |
|-----------|--------------|----------------|---------|
|-----------|--------------|----------------|---------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.22 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Chloride (24) | 17 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.7 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Sulfate (24) | 28 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 03/23/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 235 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 235 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Solids, Total Dissolved (24) | 335 | 1.00 | 0.820 | 1.00 | | mg/L | 03/21/15 | 03/21/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 03/23/15 | 03/23/15 | SM 2540 D |
| pH (24) | 7.41 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/17/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.35 | 0.50 | 0.28 | 1.00 | J | mg/L | 03/20/15 | 03/20/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 10 | 0.50 | 0.24 | 1.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.1 | 0.50 | 0.24 | 1.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

03/17/15

Work Order:

15-03-1281

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|-------------------|---------------------|----------------|
| Method Blank | | N/A | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|-------|---------------|---------------|-----------------|
| Fluoride (24) | ND | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Chloride (24) | ND | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Nitrate (as N) (24) | ND | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Sulfate (24) | ND | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/17/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 03/23/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 03/26/15 | SM 2320B |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.82 | 1.00 | | mg/L | 03/21/15 | 03/21/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 03/23/15 | 03/23/15 | SM 2540 D |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 03/20/15 | 03/20/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 03/26/15 | 03/27/15 | SM 5310 B |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-03-1250-1 | Sample | Aqueous | IC 7 | N/A | 03/17/15 14:47 | 150317S01 |
| 15-03-1250-1 | Matrix Spike | Aqueous | IC 7 | N/A | 03/17/15 15:21 | 150317S01 |
| 15-03-1250-1 | Matrix Spike Duplicate | Aqueous | IC 7 | N/A | 03/17/15 15:39 | 150317S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.4690 | 250.0 | 229.8 | 92 | 239.5 | 96 | 80-120 | 4 | 0-20 | |
| Chloride | 36.76 | 5000 | 4952 | 98 | 4956 | 98 | 80-120 | 0 | 0-20 | |
| Bromide | 0.3610 | 500.0 | 510.3 | 102 | 509.4 | 102 | 80-120 | 0 | 0-20 | |
| Nitrate (as N) | 5.892 | 500.0 | 514.7 | 102 | 505.8 | 100 | 80-120 | 2 | 0-20 | |
| Sulfate | 294.4 | 5000 | 5132 | 97 | 5153 | 97 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HCS 110 | Sample | Aqueous | ACA 1 | N/A | 03/23/15 17:40 | 150323LO1 |
| HCS 110 | Matrix Spike | Aqueous | ACA 1 | N/A | 03/23/15 17:40 | 150323LO1 |
| HCS 110 | Matrix Spike Duplicate | Aqueous | ACA 1 | N/A | 03/23/15 17:40 | 150323LO1 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | ND | 0.2000 | 0.1374 | 69 | 0.1364 | 68 | 90-110 | 1 | 0-25 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|---------------------|-----------------|------------|---------------|-------------------|
| HCS 110 | Sample | Aqueous | TOC 8 | 03/26/15 | 03/27/15 13:32 | F0326TOCS2 | | | | |
| HCS 110 | Matrix Spike | Aqueous | TOC 8 | 03/26/15 | 03/27/15 13:32 | F0326TOCS2 | | | | |
| HCS 110 | Matrix Spike Duplicate | Aqueous | TOC 8 | 03/26/15 | 03/27/15 13:32 | F0326TOCS2 | | | | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Carbon, Total Organic | 8.100 | 50.00 | 72.00 | 128 | 67.50 | 119 | 31-145 | 6 | 0-23 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| FDHCS 120 | Sample | Aqueous | TOC 8 | 03/26/15 | 03/27/15 01:43 | F0326DOCS1 | | | | |
| FDHCS 120 | Matrix Spike | Aqueous | TOC 8 | 03/26/15 | 03/27/15 01:43 | F0326DOCS1 | | | | |
| FDHCS 120 | Matrix Spike Duplicate | Aqueous | TOC 8 | 03/26/15 | 03/27/15 01:43 | F0326DOCS1 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Dissolved Organic | 2.140 | 10.00 | 14.60 | 125 | 18.40 | 163 | 31-145 | 23 | 0-25 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15

Work Order: 15-03-1281

Preparation: EPA 3010A Total

Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-03-2224-1 | Sample | Aqueous | ICP 7300 | 03/30/15 | 03/30/15 17:51 | 150330SA2 |
| 15-03-2224-1 | Matrix Spike | Aqueous | ICP 7300 | 03/30/15 | 03/30/15 17:29 | 150330SA2 |
| 15-03-2224-1 | Matrix Spike Duplicate | Aqueous | ICP 7300 | 03/30/15 | 03/30/15 17:31 | 150330SA2 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Calcium | 69.32 | 0.5000 | 76.50 | 4X | 74.12 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 25.54 | 0.5000 | 28.50 | 4X | 27.13 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 6.721 | 5.000 | 12.84 | 122 | 12.05 | 107 | 83-131 | 6 | 0-7 | |
| Sodium | 324.9 | 5.000 | 344.3 | 4X | 337.2 | 4X | 73-127 | 4X | 0-9 | Q |
| Strontium | 0.7751 | 0.5000 | 1.358 | 117 | 1.338 | 112 | 81-123 | 2 | 0-6 | |
| Silicon | 4.873 | 0.5000 | 5.801 | 4X | 5.559 | 4X | 24-180 | 4X | 0-15 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|---------------------------|------------------------|-------------|----------|------------|---------------|----------------|---------------------|-----|--------|------------|
| HCS 110 | Sample | | Aqueous | ICP/MS 03 | 03/19/15 | 03/23/15 15:42 | 150319SA1 | | | |
| HCS 110 | Matrix Spike | | Aqueous | ICP/MS 03 | 03/19/15 | 03/23/15 14:35 | 150319SA1 | | | |
| HCS 110 | Matrix Spike Duplicate | | Aqueous | ICP/MS 03 | 03/19/15 | 03/23/15 14:38 | 150319SA1 | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.09644 | 96 | 0.09549 | 95 | 85-133 | 1 | 0-11 | |
| Arsenic | ND | 0.1000 | 0.09379 | 94 | 0.09376 | 94 | 73-127 | 0 | 0-11 | |
| Barium | 0.03710 | 0.1000 | 0.1371 | 100 | 0.1416 | 105 | 74-128 | 3 | 0-10 | |
| Beryllium | ND | 0.1000 | 0.09495 | 95 | 0.09408 | 94 | 56-122 | 1 | 0-11 | |
| Cadmium | ND | 0.1000 | 0.09338 | 93 | 0.09297 | 93 | 84-114 | 0 | 0-8 | |
| Chromium | ND | 0.1000 | 0.1037 | 104 | 0.1040 | 104 | 73-133 | 0 | 0-11 | |
| Copper | ND | 0.1000 | 0.1013 | 101 | 0.1006 | 101 | 72-108 | 1 | 0-10 | |
| Lead | ND | 0.1000 | 0.1011 | 101 | 0.1021 | 102 | 79-121 | 1 | 0-10 | |
| Nickel | 0.001702 | 0.1000 | 0.1003 | 99 | 0.1031 | 101 | 68-122 | 3 | 0-10 | |
| Selenium | ND | 0.1000 | 0.07938 | 79 | 0.07723 | 77 | 59-125 | 3 | 0-12 | |
| Silver | ND | 0.05000 | 0.04975 | 100 | 0.04920 | 98 | 68-128 | 1 | 0-14 | |
| Thallium | ND | 0.1000 | 0.1026 | 103 | 0.1031 | 103 | 73-121 | 0 | 0-11 | |
| Zinc | 0.02449 | 0.1000 | 0.09241 | 68 | 0.09593 | 71 | 43-145 | 4 | 0-39 | |
| Aluminum | ND | 0.1000 | 0.1631 | 163 | 0.2019 | 202 | 47-161 | 21 | 0-24 | 3 |
| Iron | 0.05202 | 5.100 | 5.922 | 115 | 6.415 | 125 | 27-201 | 8 | 0-24 | |
| Manganese | 0.002430 | 0.1000 | 0.1089 | 107 | 0.1133 | 111 | 72-126 | 4 | 0-42 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15

Work Order: 15-03-1281

Preparation: EPA 7470A Total

Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-03-1285-1 | Sample | Aqueous | Mercury 04 | 03/18/15 | 03/18/15 18:53 | 150318S07 |
| 15-03-1285-1 | Matrix Spike | Aqueous | Mercury 04 | 03/18/15 | 03/18/15 18:55 | 150318S07 |
| 15-03-1285-1 | Matrix Spike Duplicate | Aqueous | Mercury 04 | 03/18/15 | 03/18/15 18:57 | 150318S07 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | 0.0007127 | 0.01000 | 0.01092 | 102 | 0.01046 | 97 | 57-141 | 4 | 0-10 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HCS 130 | Sample | Aqueous | LC/TQ 2 | 03/19/15 | 03/23/15 15:46 | 150319S20 | | | | |
| HCS 130 | Matrix Spike | Aqueous | LC/TQ 2 | 03/19/15 | 03/23/15 16:19 | 150319S20 | | | | |
| HCS 130 | Matrix Spike Duplicate | Aqueous | LC/TQ 2 | 03/19/15 | 03/23/15 16:28 | 150319S20 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Caffeine | ND | 100.0 | 107.1 | 107 | 101.5 | 101 | 70-130 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HCS 130 | Sample | Aqueous | GC/MS BB | 03/20/15 | 03/21/15 03:37 | 150320S014 | | | | |
| HCS 130 | Matrix Spike | Aqueous | GC/MS BB | 03/20/15 | 03/21/15 04:04 | 150320S014 | | | | |
| HCS 130 | Matrix Spike Duplicate | Aqueous | GC/MS BB | 03/20/15 | 03/21/15 04:31 | 150320S014 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 47.89 | 96 | 46.22 | 92 | 74-122 | 4 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 44.63 | 89 | 45.67 | 91 | 60-144 | 2 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 50.61 | 101 | 51.11 | 102 | 73-120 | 1 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 51.91 | 104 | 51.07 | 102 | 80-122 | 2 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 48.22 | 96 | 49.71 | 99 | 70-120 | 3 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 53.56 | 107 | 51.63 | 103 | 64-142 | 4 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 47.34 | 95 | 47.20 | 94 | 52-136 | 0 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 49.39 | 99 | 49.25 | 99 | 77-125 | 0 | 0-24 | |
| Toluene | ND | 50.00 | 47.89 | 96 | 47.19 | 94 | 72-126 | 1 | 0-23 | |
| Trichloroethene | ND | 50.00 | 48.68 | 97 | 46.71 | 93 | 74-128 | 4 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 39.28 | 79 | 42.61 | 85 | 67-133 | 8 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 101.7 | 102 | 101.9 | 102 | 63-129 | 0 | 0-25 | |
| o-Xylene | ND | 50.00 | 51.46 | 103 | 51.54 | 103 | 62-128 | 0 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 48.83 | 98 | 47.65 | 95 | 68-134 | 2 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|--------------|-------------|------------|----------------|----------------|-----------------------|
| HCS 110 | Sample | Aqueous | ICP/MS 03 | 03/19/15 00:00 | 03/23/15 15:42 | 150319SA1 |
| HCS 110 | PDS | Aqueous | ICP/MS 03 | 03/19/15 00:00 | 03/23/15 14:42 | 150319SA1 |
| Parameter | Sample Conc. | Spike Added | PDS Conc. | PDS %Rec. | %Rec. CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.1019 | 102 | 75-125 | |
| Arsenic | ND | 0.1000 | 0.09725 | 97 | 75-125 | |
| Barium | 0.03710 | 0.1000 | 0.1415 | 104 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.09996 | 100 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.09881 | 99 | 75-125 | |
| Chromium | ND | 0.1000 | 0.1069 | 107 | 75-125 | |
| Copper | ND | 0.1000 | 0.1015 | 101 | 75-125 | |
| Lead | ND | 0.1000 | 0.1061 | 106 | 75-125 | |
| Nickel | 0.001702 | 0.1000 | 0.1042 | 103 | 75-125 | |
| Selenium | ND | 0.1000 | 0.08272 | 83 | 75-125 | |
| Silver | ND | 0.05000 | 0.05097 | 102 | 75-125 | |
| Thallium | ND | 0.1000 | 0.1074 | 107 | 75-125 | |
| Zinc | 0.02449 | 0.1000 | 0.1144 | 90 | 75-125 | |
| Aluminum | ND | 0.1000 | 0.1335 | 133 | 75-125 | 5 |
| Iron | 0.05202 | 5.100 | 5.439 | 106 | 75-125 | |
| Manganese | 0.002430 | 0.1000 | 0.1135 | 111 | 75-125 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-03-1608-1 | Sample | Aqueous | PH1/BUR03 | N/A | 03/26/15 17:35 | F0326ALKD1 |
| 15-03-1608-1 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 03/26/15 17:35 | F0326ALKD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|---|---------------------|------------------|------------|---------------|-------------------|
| Alkalinity, Total (as CaCO ₃) | 268.0 | 265.0 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-03-1608-1 | Sample | Aqueous | PH1/BUR03 | N/A | 03/26/15 17:35 | F0326HCOD1 |
| 15-03-1608-1 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 03/26/15 17:35 | F0326HCOD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Bicarbonate (as CaCO ₃) | 268.0 | 265.0 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-03-1608-1 | Sample | Aqueous | PH1/BUR03 | N/A | 03/26/15 17:35 | F0326C03D1 |
| 15-03-1608-1 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 03/26/15 17:35 | F0326C03D1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------------------|------------------|----------------|----------------|------------------------|
| HCS 110 | Sample | Aqueous | SC 5 | 03/21/15 00:00 | 03/21/15 15:00 | F0321TDSD1 |
| HCS 110 | Sample Duplicate | Aqueous | SC 5 | 03/21/15 00:00 | 03/21/15 15:00 | F0321TDSD1 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Solids, Total Dissolved | | 295.0 | 310.0 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-03-1366-1 | Sample | Aqueous | N/A | 03/23/15 00:00 | 03/23/15 17:00 | F0323TSSD2 |
| 15-03-1366-1 | Sample Duplicate | Aqueous | N/A | 03/23/15 00:00 | 03/23/15 17:00 | F0323TSSD2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Suspended | 170.0 | 167.0 | 2 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------------------|------------------|---------------|----------------|------------------------|
| HCS 110 | Sample | Aqueous | PH 1 | N/A | 03/17/15 11:56 | F0317PHD1 |
| HCS 110 | Sample Duplicate | Aqueous | PH 1 | N/A | 03/17/15 11:56 | F0317PHD1 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| pH | | 7.780 | 7.790 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 4500 N Org B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-03-1267-1 | Sample | Aqueous | BUR05 | 03/20/15 00:00 | 03/20/15 17:46 | F0320TKND1 |
| 15-03-1267-1 | Sample Duplicate | Aqueous | BUR05 | 03/20/15 00:00 | 03/20/15 17:46 | F0320TKND1 |

| Parameter | Sample Conc. | DUP Conc. | RPD | RPD CL | Qualifiers |
|-------------------------|--------------|-----------|-----|--------|------------|
| Total Kjeldahl Nitrogen | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|--------------------|------------------------|------------------|-----------------|-------------------|
| 099-12-906-5530 | LCS | Aqueous | IC 7 | N/A | 03/17/15 11:45 | 150317L01 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Fluoride | | 2.500 | 2.478 | 99 | 90-110 | |
| Chloride | | 50.00 | 49.34 | 99 | 90-110 | |
| Bromide | | 5.000 | 5.168 | 103 | 90-110 | |
| Nitrate (as N) | | 5.000 | 5.019 | 100 | 90-110 | |
| Sulfate | | 50.00 | 49.20 | 98 | 90-110 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-739-101 | LCS | Aqueous | ACA 1 | N/A | 03/23/15 17:40 | 150323SO1 | | | |
| 099-12-739-101 | LCSD | Aqueous | ACA 1 | N/A | 03/23/15 17:40 | 150323SO1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Phosphorus, Total | 0.2000 | 0.1914 | 96 | 0.1979 | 99 | 90-110 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
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San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-859-629 | LCS | Aqueous | PH1/BUR03 | N/A | 03/26/15 17:35 | F0326ALKB1 | | | |
| 099-15-859-629 | LCSD | Aqueous | PH1/BUR03 | N/A | 03/26/15 17:35 | F0326ALKB1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 100.0 | 99.00 | 99 | 99.00 | 99 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4482 | LCS | Aqueous | SC 5 | 03/21/15 | 03/21/15 15:00 | F0321TDSL1 | | | |
| 099-12-180-4482 | LCSD | Aqueous | SC 5 | 03/21/15 | 03/21/15 15:00 | F0321TDSL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 105.0 | 105 | 110.0 | 110 | 80-120 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-09-010-7112 | LCS | Aqueous | N/A | 03/23/15 | 03/23/15 17:00 | F0323TSSB2 | | | |
| 099-09-010-7112 | LCSD | Aqueous | N/A | 03/23/15 | 03/23/15 17:00 | F0323TSSB2 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Suspended | 100.0 | 89.00 | 89 | 93.00 | 93 | 80-120 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-097-5569 | LCS | Aqueous | TOC 8 | 03/26/15 | 03/27/15 13:32 | F0326TOCL2 | | | |
| 099-05-097-5569 | LCSD | Aqueous | TOC 8 | 03/26/15 | 03/27/15 13:32 | F0326TOCL2 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 10.00 | 9.970 | 100 | 10.70 | 107 | 80-120 | 7 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-115-1414 | LCS | Aqueous | TOC 8 | 03/26/15 | 03/27/15 01:43 | F0326DOCL1 | | | |
| 099-05-115-1414 | LCSD | Aqueous | TOC 8 | 03/26/15 | 03/27/15 01:43 | F0326DOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Dissolved Organic | 10.00 | 10.30 | 103 | 10.30 | 103 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-15-683-1193 | LCS | Aqueous | ICP 7300 | 03/30/15 | 03/30/15 17:22 | 150330LA2F |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Calcium | | 0.5000 | 0.5074 | 101 | 80-120 | |
| Magnesium | | 0.5000 | 0.5119 | 102 | 80-120 | |
| Potassium | | 5.000 | 5.177 | 104 | 80-120 | |
| Sodium | | 5.000 | 5.109 | 102 | 80-120 | |
| Strontium | | 0.5000 | 0.5166 | 103 | 80-120 | |
| Silicon | | 0.5000 | 0.4310 | 86 | 80-120 | |


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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-769 | LCS | Aqueous | ICP/MS 03 | 03/19/15 | 03/23/15 13:29 | 150319LA1F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.1037 | 104 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.1035 | 103 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.09899 | 99 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.1014 | 101 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.1039 | 104 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.1068 | 107 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.1039 | 104 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.1002 | 100 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.1017 | 102 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.1056 | 106 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.05267 | 105 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.09832 | 98 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.1050 | 105 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.1146 | 115 | 80-120 | 73-127 | |
| Iron | 5.100 | 5.039 | 99 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.1050 | 105 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|------------------|
| 099-15-763-517 | LCS | Aqueous | Mercury 04 | 03/18/15 | 03/18/15 18:51 | 150318L07F |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.01000 | 0.009958 | 100 | 85-121 | |


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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 099-16-376-11 | LCS | Aqueous | LC/TQ 2 | 03/19/15 | 03/23/15 15:13 | 150319L20 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Caffeine | | 100.0 | 106.4 | 106 | 80-120 | |


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Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|--------|-----|--------|------------|
| 099-12-529-792 | LCS | Aqueous | GC 44 | 03/19/15 | 03/20/15 12:16 | 150319L01 | | | | |
| 099-12-529-792 | LCSD | Aqueous | GC 44 | 03/19/15 | 03/20/15 12:31 | 150319L01 | | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Alpha-BHC | 0.5000 | 0.5797 | 116 | 0.5728 | 115 | 50-135 | 36-149 | 1 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.5939 | 119 | 0.5913 | 118 | 50-135 | 36-149 | 0 | 0-25 | |
| Beta-BHC | 0.5000 | 0.5539 | 111 | 0.5643 | 113 | 50-135 | 36-149 | 2 | 0-25 | |
| Heptachlor | 0.5000 | 0.5715 | 114 | 0.5307 | 106 | 50-135 | 36-149 | 7 | 0-25 | |
| Delta-BHC | 0.5000 | 0.5636 | 113 | 0.5327 | 107 | 50-135 | 36-149 | 6 | 0-25 | |
| Aldrin | 0.5000 | 0.5480 | 110 | 0.4992 | 100 | 50-135 | 36-149 | 9 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.5681 | 114 | 0.5506 | 110 | 50-135 | 36-149 | 3 | 0-25 | |
| Endosulfan I | 0.5000 | 0.5836 | 117 | 0.5593 | 112 | 50-135 | 36-149 | 4 | 0-25 | |
| Dieldrin | 0.5000 | 0.5837 | 117 | 0.5690 | 114 | 50-135 | 36-149 | 3 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.5680 | 114 | 0.5658 | 113 | 50-135 | 36-149 | 0 | 0-25 | |
| Endrin | 0.5000 | 0.5576 | 112 | 0.5328 | 107 | 50-135 | 36-149 | 5 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.6081 | 122 | 0.5490 | 110 | 50-135 | 36-149 | 10 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.5604 | 112 | 0.5329 | 107 | 50-135 | 36-149 | 5 | 0-25 | |
| Endosulfan II | 0.5000 | 0.5716 | 114 | 0.5623 | 112 | 50-135 | 36-149 | 2 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.5418 | 108 | 0.5378 | 108 | 50-135 | 36-149 | 1 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.5492 | 110 | 0.5394 | 108 | 50-135 | 36-149 | 2 | 0-25 | |
| Methoxychlor | 0.5000 | 0.5491 | 110 | 0.5496 | 110 | 50-135 | 36-149 | 0 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-533-1015 | LCS | Aqueous | GC 58 | 03/19/15 | 03/20/15 18:27 | 150319L02 | | | |
| 099-12-533-1015 | LCSD | Aqueous | GC 58 | 03/19/15 | 03/20/15 18:45 | 150319L02 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Aroclor-1016 | 2.000 | 1.924 | 96 | 1.971 | 99 | 50-135 | 2 | 0-25 | |
| Aroclor-1260 | 2.000 | 1.880 | 94 | 1.989 | 99 | 50-135 | 6 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-963-90 | LCS | Aqueous | | GC 35 | 03/19/15 | 03/20/15 11:29 | 150319L03 | | | |
| 099-15-963-90 | LCSD | Aqueous | | GC 35 | 03/19/15 | 03/20/15 12:15 | 150319L03 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.04451 | 111 | 0.04563 | 114 | 30-130 | 13-147 | 2 | 0-30 | |
| Bolstar | 0.04000 | 0.04440 | 111 | 0.04192 | 105 | 30-130 | 13-147 | 6 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.04042 | 101 | 0.04003 | 100 | 30-130 | 13-147 | 1 | 0-30 | |
| Coumaphos | 0.04000 | 0.04421 | 111 | 0.04463 | 112 | 30-130 | 13-147 | 1 | 0-30 | |
| Diazinon | 0.04000 | 0.04504 | 113 | 0.04489 | 112 | 30-130 | 13-147 | 0 | 0-30 | |
| Disulfoton | 0.04000 | 0.04279 | 107 | 0.04050 | 101 | 30-130 | 13-147 | 5 | 0-30 | |
| Ethoprop | 0.04000 | 0.04074 | 102 | 0.03955 | 99 | 30-130 | 13-147 | 3 | 0-30 | |
| Fensulfothion | 0.04000 | 0.04534 | 113 | 0.04549 | 114 | 30-130 | 13-147 | 0 | 0-30 | |
| Fenthion | 0.04000 | 0.04212 | 105 | 0.04209 | 105 | 30-130 | 13-147 | 0 | 0-30 | |
| Merphos | 0.04000 | 0.03744 | 94 | 0.03799 | 95 | 30-130 | 13-147 | 1 | 0-30 | |
| Methyl Parathion | 0.04000 | 0.04153 | 104 | 0.04150 | 104 | 30-130 | 13-147 | 0 | 0-30 | |
| Phorate | 0.04000 | 0.05059 | 126 | 0.04860 | 122 | 30-130 | 13-147 | 4 | 0-30 | |
| Ronnel | 0.04000 | 0.03901 | 98 | 0.03838 | 96 | 30-130 | 13-147 | 2 | 0-30 | |
| Stirophos | 0.04000 | 0.02787 | 70 | 0.02752 | 69 | 30-130 | 13-147 | 1 | 0-30 | |
| Tokuthion | 0.04000 | 0.04108 | 103 | 0.04055 | 101 | 30-130 | 13-147 | 1 | 0-30 | |
| Trichloronate | 0.04000 | 0.04114 | 103 | 0.04071 | 102 | 30-130 | 13-147 | 1 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-034-639 | LCS | Aqueous | GC 40 | 03/23/15 | 03/24/15 15:34 | 150323L02 | | | |
| 095-01-034-639 | LCSD | Aqueous | GC 40 | 03/23/15 | 03/24/15 15:57 | 150323L02 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| 2,4-D | 20.00 | 18.92 | 95 | 22.02 | 110 | 30-130 | 15 | 0-30 | |
| 2,4,5-T | 2.000 | 2.065 | 103 | 2.435 | 122 | 30-130 | 16 | 0-30 | |
| 2,4-DB | 20.00 | 19.48 | 97 | 22.44 | 112 | 30-130 | 14 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-4010 | LCS | Aqueous | | GC/MS TT | 03/18/15 | 03/19/15 16:57 | 150318L08 | | | |
| 095-01-003-4010 | LCSD | Aqueous | | GC/MS TT | 03/18/15 | 03/19/15 17:16 | 150318L08 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 200.0 | 217.1 | 109 | 214.1 | 107 | 61-120 | 51-130 | 1 | 0-20 | |
| Acenaphthylene | 200.0 | 208.2 | 104 | 206.6 | 103 | 55-120 | 44-131 | 1 | 0-20 | |
| Butyl Benzyl Phthalate | 200.0 | 213.4 | 107 | 207.1 | 104 | 56-122 | 45-133 | 3 | 0-20 | |
| 4-Chloro-3-Methylphenol | 200.0 | 185.4 | 93 | 180.2 | 90 | 52-120 | 41-131 | 3 | 0-20 | |
| 2-Chlorophenol | 200.0 | 181.9 | 91 | 178.7 | 89 | 47-120 | 35-132 | 2 | 0-20 | |
| 1,4-Dichlorobenzene | 200.0 | 170.3 | 85 | 166.3 | 83 | 36-120 | 22-134 | 2 | 0-20 | |
| Dimethyl Phthalate | 200.0 | 208.7 | 104 | 206.1 | 103 | 60-120 | 50-130 | 1 | 0-20 | |
| 2,4-Dinitrotoluene | 200.0 | 206.1 | 103 | 206.3 | 103 | 61-121 | 51-131 | 0 | 0-20 | |
| Fluorene | 200.0 | 218.0 | 109 | 214.5 | 107 | 67-120 | 58-129 | 2 | 0-20 | |
| N-Nitroso-di-n-propylamine | 200.0 | 189.5 | 95 | 187.8 | 94 | 39-123 | 25-137 | 1 | 0-20 | |
| Naphthalene | 200.0 | 183.5 | 92 | 176.9 | 88 | 54-120 | 43-131 | 4 | 0-20 | |
| 4-Nitrophenol | 200.0 | 170.6 | 85 | 173.5 | 87 | 14-120 | 0-138 | 2 | 0-20 | |
| Pentachlorophenol | 200.0 | 172.5 | 86 | 173.4 | 87 | 31-127 | 15-143 | 1 | 0-20 | |
| Phenol | 200.0 | 134.4 | 67 | 134.5 | 67 | 17-120 | 0-137 | 0 | 0-20 | |
| Pyrene | 200.0 | 194.1 | 97 | 187.5 | 94 | 58-124 | 47-135 | 3 | 0-20 | |
| 1,2,4-Trichlorobenzene | 200.0 | 174.0 | 87 | 168.3 | 84 | 49-120 | 37-132 | 3 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/17/15
Work Order: 15-03-1281
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-16646 | LCS | Aqueous | GC/MS BB | 03/20/15 | 03/21/15 01:22 | 150320L025 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 46.43 | 93 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 44.63 | 89 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 50.78 | 102 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 50.82 | 102 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 50.17 | 100 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 50.79 | 102 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 46.81 | 94 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 49.15 | 98 | 80-123 | 73-130 | |
| Toluene | 50.00 | 46.65 | 93 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 48.43 | 97 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 42.24 | 84 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 101.3 | 101 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 51.39 | 103 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 47.13 | 94 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Sample Analysis Summary Report

Work Order: 15-03-1281

Page 1 of 1

| <u>Method</u> | <u>Extraction</u> | <u>Chemist ID</u> | <u>Instrument</u> | <u>Analytical Location</u> |
|-----------------------|-------------------|-------------------|-------------------|----------------------------|
| EPA 1694 (M) Caffeine | EPA 1694 | 262 | LC/TQ 2 | 1 |
| EPA 300.0 | N/A | 650 | IC 7 | 1 |
| EPA 365.1 | N/A | 735 | ACA 1 | 1 |
| EPA 6010B | EPA 3005A Filt. | 935 | ICP 7300 | 1 |
| EPA 6020 | EPA 3005A Filt. | 598 | ICP/MS 03 | 1 |
| EPA 7470A | EPA 7470A Filt. | 915 | Mercury 04 | 1 |
| EPA 8081A | EPA 3510C | 960 | GC 44 | 1 |
| EPA 8082 | EPA 3510C | 944 | GC 58 | 1 |
| EPA 8141A | EPA 3510C | 949 | GC 35 | 1 |
| EPA 8151A | EPA 8151A | 944 | GC 40 | 1 |
| EPA 8260B | EPA 5030C | 959 | GC/MS BB | 2 |
| EPA 8270C | EPA 3510C | 923 | GC/MS TT | 1 |
| SM 2320B | N/A | 688 | PH1/BUR03 | 1 |
| SM 2540 C | N/A | 689 | SC 5 | 1 |
| SM 2540 D | N/A | 689 | N/A | 1 |
| SM 4500 H+ B | N/A | 688 | PH 1 | 1 |
| SM 4500 N Org B | N/A | 685 | BUR05 | 1 |
| SM 5310 B | N/A | 735 | TOC 8 | 1 |

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Location 1: 7440 Lincoln Way, Garden Grove, CA 92841

Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841

Glossary of Terms and Qualifiers

Work Order: 15-03-1281

Page 1 of 1

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|--|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDS or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.

ORIGIN ID: SARA (210) 877-2847
SWCA INC

6200 UTSA BLVD STE 102

SAN ANTONIO, TX 782491618
UNITED STATES US

SHIP DATE: 16MAR15
ACTWGT: 62.0 LB
CAD: 7OFFC1522
DIMS: 24x14x13 IN

BILL SENDER

ORIGIN ID: SARA (210) 877-2847
SWCA INC

6200 UTSA BLVD STE 102

SAN ANTONIO, TX 782491618
UNITED STATES US

WGT: 62.0 LB
OFFC1522
Page 1 of 126
L SENDER

CALSCIENCE ENVIRONMENTAL LAB
7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895-5494

REF:

DEPT:

FedEx
Express



1 of 6

TRK# 8062 7398 5690

0215
MASTER

A7 APVA

TUE - 17 MAR 3:00P
STANDARD OVERNIGHT

92841

CA-US SNA

CALSCIENCE ENVIRONMENTAL LAB
7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895-5494

REF:

DEPT:

FedEx
Express



2 of 6

MPS# 7803 5520 1130

0681
ORIGIN ID: SARA (210) 877-2847
SWCA INC

6200 UTSA BLVD STE 102

SAN ANTONIO, TX 782491618
UNITED STATES US

TUE - 17 MAR 3:00P
STANDARD OVERNIGHT

SHIP DATE: 16MAR15
ACTWGT: 60.0 LB
CAD: 7OFFC1522
DIMS: 24x14x13 IN

BILL SENDER

CALSCIENCE ENVIRONMENTAL LAB
7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895-5494

REF:

DEPT:

FedEx
Express



3 of 6

MPS# 7803 5520 1140

0681
Mstr# 8062 7398 5690

A7 APVA

TUE - 17 MAR 3:00P
STANDARD OVERNIGHT

92841

CA-US SNA

MPS# 7803 5520 1150

0681
Mstr# 8062 7398 5690

A7 APVA

TUE - 17 MAR 3:00P
STANDARD OVERNIGHT

92841

CA-US SNA

TUE - 17 MAR 3:00P
STANDARD OVERNIGHT

MPS# 7803 5520 1161

0681
Mstr# 8062 7398 5690

A7 APVA

92841

CA-US SNA

MPS# 7803 5520 1172

0681
Mstr# 8062 7398 5690

A7 APVA

TUE - 17 MAR 3:00P
STANDARD OVERNIGHT

92841

CA-US SNA

SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 6

CLIENT:

SWCA

DATE: 03/17/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 2.6 °C (w/ CF): 2.8 °C ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Not Present ☐ Not Intact ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☒ Not Present ☐ Not Intact ☐ N/A

Checked by: 826

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|--------------------------|--------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date/timeSampler's name indicated on COC..... ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC..... ☐ Yes ☒ No ☐ N/ASample container(s) intact and in good condition..... ☒ Yes ☐ No ☐ N/AProper containers for analyses requested..... ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested..... ☒ Yes ☐ No ☐ N/ASamples received within holding time..... ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen..... ☐ Yes ☒ No ☒ N/AProper preservation chemical(s) noted on COC and/or sample container..... ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace..... ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation..... ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: 3

(Trip Blank Lot Number: 15115A)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna2 ☐ 100PJ ☒ 100PJna2 ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznn ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBnf ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna2 ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 GAL ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® ☐ TerraCores® ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF Other Matrix (____): ☐ _____ ☐ _____

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO3, na=NaOH, na2=Na2S2O3, p=H3PO4,

s = H2SO4, u = ultra-pure, znn = Zn(CH3CO2)2 + NaOH

Labeled/Checked by: 826

Reviewed by: 300

WORK ORDER NUMBER: 15-03-1281

SAMPLE RECEIPT CHECKLIST

COOLER 2 OF 6

CLIENT:

SWCA

DATE: 03/17/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 1.9 °C (w/ CF): 2.1 °C ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Not Present ☐ Not Intact ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☒ Not Present ☐ Not Intact ☐ N/A

Checked by: 836

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete..... ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date/timeSampler's name indicated on COC..... ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC..... ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition..... ☒ Yes ☐ No ☐ N/AProper containers for analyses requested..... ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested..... ☒ Yes ☐ No ☐ N/ASamples received within holding time..... ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☒ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace..... ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation..... ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na2} ☐ 100PJ ☐ 100PJ_{na2} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☒ 250PB_h ☐ 500AGB ☐ 500AG_J ☐ 500AG_J_s☐ 500PB ☒ 1AGB ☐ 1AGB_{na2} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5 GAL ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® ☐ TerraCores® ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF Other Matrix (____): ☐ _____ ☐ _____

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Labeled/Checked by: 836

Reviewed by: 762

SAMPLE RECEIPT CHECKLIST

COOLER 3 OF 6

CLIENT:

SWCA

DATE: 03/17/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 3.1 °C (w/ CF): 3.3 °C ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Not Present ☐ Not Intact ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☒ Not Present ☐ Not Intact ☐ N/A

Checked by: 876

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|--------------------------|--------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date/timeSampler's name indicated on COC..... ☒ ☐ ☐Sample container label(s) consistent with COC..... ☒ ☐ ☐Sample container(s) intact and in good condition..... ☒ ☐ ☐Proper containers for analyses requested..... ☒ ☐ ☐Sufficient volume/mass for analyses requested..... ☒ ☐ ☐Samples received within holding time..... ☒ ☐ ☐

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen..... ☐ ☒ ☒Proper preservation chemical(s) noted on COC and/or sample container..... ☒ ☐ ☐

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace..... ☒ ☐ ☐☐ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation..... ☐ ☐ ☒

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 GA ☐ ☐Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® ☐ TerraCores® ☐Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF Other Matrix (____): ☐ ☐

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Labeled/Checked by: 876

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 4 OF 6

CLIENT:

SWCA

DATE: 03/17/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 2.7 °C (w/ CF): 2.9 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Not Present ☐ Not Intact ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☒ Not Present ☐ Not Intact ☐ N/A

Checked by: 836

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete..... ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date/time

Sampler's name indicated on COC..... ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC..... ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition..... ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested..... ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested..... ☒ Yes ☐ No ☐ N/A

Samples received within holding time..... ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☒ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☒ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace..... ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation..... ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na2} ☐ 100PJ ☐ 100PJ_{na2} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☒ 250PB_{ng} ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s
☐ 500PB ☒ 1AGB ☐ 1AGB_{na2} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5 G_{AM} ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® ☐ TerraCores® ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF Other Matrix (____): ☐ _____ ☐ _____

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Labeled/Checked by: 836

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 5 OF 6

CLIENT: SWCA

DATE: 03/17/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 2.8 °C (w/ CF): 3.0 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Not Present ☐ Not Intact ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☒ Not Present ☐ Not Intact ☐ N/A

Checked by: 836

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|---|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date/time | | | |
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> 3/17/15 |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na2} ☐ 100PJ ☐ 100PJ_{na2} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☒ 250PB_{na} ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s

☐ 500PB ☒ 1AGB ☐ 1AGB_{na2} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5 GAL ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® ☐ TerraCores® ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF Other Matrix (_____) ☐ _____ ☐ _____

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Labeled/Checked by: 836

Reviewed by: 780

SAMPLE RECEIPT CHECKLIST

COOLER 6 OF 6

CLIENT: SWCA

DATE: 03/17/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 2.4 °C (w/ CF): 2.6 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Not Present ☐ Not Intact ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☒ Not Present ☐ Not Intact ☐ N/A

Checked by: 876

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|--------------------------|--------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date/time

Sampler's name indicated on COC..... ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC..... ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition..... ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested..... ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested..... ☒ Yes ☐ No ☐ N/A

Samples received within holding time..... ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☒ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☒ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace..... ☒ Yes ☐ No ☐ N/A

☐ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation..... ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na2} ☐ 100PJ ☐ 100PJ_{na2} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☒ 250PB_h ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s
☐ 500PB ☒ 1AGB ☐ 1AGB_{na2} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5 GAL ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® ☐ TerraCores® ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF Other Matrix (_____) ☐ _____ ☐ _____

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Labeled/Checked by: 876

Reviewed by: 700

SAMPLE ANOMALY REPORT

DATE: 03 / 17 / 2015

SAMPLES, CONTAINERS, AND LABELS:

- ☐ Sample(s) NOT RECEIVED but listed on COC
- ☐ Sample(s) received but NOT LISTED on COC
- ☐ Holding time expired (list client or ECI sample ID and analysis)
- ☐ Insufficient sample amount for requested analysis (list analysis)
- ☐ Improper container(s) used (list analysis)
- ☐ Improper preservative used (list analysis)
- ☐ No preservative noted on COC or label (list analysis and notify lab)
- ☐ Sample container(s) not labeled
- ☐ Client sample label(s) illegible (list container type and analysis)
- ☒ Client sample label(s) do not match COC (comment)
 - ☐ Project information
 - ☐ Client sample ID
 - ☐ Sampling date and/or time
 - ☒ Number of container(s)
 - ☐ Requested analysis
- ☐ Sample container(s) compromised (comment)
 - ☐ Broken
 - ☐ Water present in sample container
- ☐ Air sample container(s) compromised (comment)
 - ☐ Flat
 - ☐ Very low in volume
 - ☐ Leaking (not transferred; duplicate bag submitted)
 - ☐ Leaking (transferred into ECI Tedlar™ bags*)
 - ☐ Leaking (transferred into client's Tedlar™ bags*)

* Transferred at client's request.

MISCELLANEOUS: (Describe)

HEADSPACE:

(Containers with bubble > 6 mm or ¼ inch for volatile organic or dissolved gas analysis)

| ECI Sample ID | ECI Container ID | Total Number** | ECI Sample ID | ECI Container ID | Total Number** |
|---------------|------------------|----------------|---------------|------------------|----------------|
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

Comments: _____

** Record the total number of containers (i.e., vials or bottles) for the affected sample.

Comments

(-7) Received only 1-vial
(TB) .

Comments

(Containers with bubble for other analysis)

| ECI Sample ID | ECI Container ID | Total Number** | Requested Analysis |
|------------------|---------------------|-------------------|--------------------|
| | | | |
| | | | |
| | | | |
| | | | |

Reported by: 826

Reviewed by: 3/00

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HSM 110
Sample Number: AA93572
Sample Matrix: Non Potable Water

Collection Date/Time: 3/25/15 10:28
Receipt Date/Time: 3/25/15 14:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93572-A | E. coli | √ | SM 9223B-2004 | <1 | MPN/100 mL | | 1 | 42188 | 3/25/15 | 15:26 | RSC |
| AA93572-A | E. Coli Holding Time - IDEXX Colilert | | NA | 4.97 | hours | | 0.00 | 42186 | 3/25/15 | 15:26 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 30, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42188

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

3/30/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 120
Sample Number: AA93573
Sample Matrix: Non Potable Water

Collection Date/Time: 3/25/15 10:51
Receipt Date/Time: 3/25/15 14:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93573-A | E. coli | √ | SM 9223B-2004 | 13 | MPN/100 mL | | 1 | 42188 | 3/25/15 | 15:26 | RSC |
| AA93573-A | E. Coli Holding Time - IDEXX Colilert | | NA | 4.58 | hours | | 0.00 | 42186 | 3/25/15 | 15:26 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 30, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42188

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

3/30/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 130
Sample Number: AA93574
Sample Matrix: Non Potable Water

Collection Date/Time: 3/25/15 11:17
Receipt Date/Time: 3/25/15 14:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93574-A | E. coli | √ | SM 9223B-2004 | 12 | MPN/100 mL | | 1 | 42188 | 3/25/15 | 15:26 | RSC |
| AA93574-A | E. Coli Holding Time - IDEXX Colilert | | NA | 4.15 | hours | | 0.00 | 42186 | 3/25/15 | 15:26 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 30, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42188

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
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3/30/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 140
Sample Number: AA93575
Sample Matrix: Non Potable Water

Collection Date/Time: 3/25/15 11:30
Receipt Date/Time: 3/25/15 14:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: slight air space.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93575-A | E. coli | √ | SM 9223B-2004 | 200 | MPN/100 mL | | 1 | 42189 | 3/25/15 | 15:26 | RSC |
| AA93575-A | E. Coli Holding Time - IDEXX Colilert | | NA | 3.93 | hours | | 0.00 | 42187 | 3/25/15 | 15:26 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 30, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42189

QC Analyte Name

Initial Blank for E. coli

Log Range for E. coli

Result

Absent

0.1621

Units

Qualifier

Lower

0.0

Acceptance Criteria

Target

Absent

Upper

0.5



Jeanette Hernandez
Water Quality Planner / QAO

3/30/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 30, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 150
Sample Number: AA93576
Sample Matrix: Non Potable Water

Collection Date/Time: 3/25/15 12:25
Receipt Date/Time: 3/25/15 14:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93576-A | E. coli | √ | SM 9223B-2004 | 66 | MPN/100 mL | | 1 | 42189 | 3/25/15 | 15:26 | RSC |
| AA93576-A | E. Coli Holding Time - IDEXX Colilert | | NA | 3.02 | hours | | 0.00 | 42187 | 3/25/15 | 15:26 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 30, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42189

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

3/30/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 160
Sample Number: AA93577
Sample Matrix: Non Potable Water

Collection Date/Time: 3/25/15 12:50
Receipt Date/Time: 3/25/15 14:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93577-A | E. coli | √ | SM 9223B-2004 | 40 | MPN/100 mL | | 1 | 42189 | 3/25/15 | 15:26 | RSC |
| AA93577-A | E. Coli Holding Time - IDEXX Colilert | | NA | 2.60 | hours | | 0.00 | 42187 | 3/25/15 | 15:26 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 30, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42189

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

3/30/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 170
Sample Number: AA93578
Sample Matrix: Non Potable Water

Collection Date/Time: 3/25/15 13:18
Receipt Date/Time: 3/25/15 14:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93578-A | E. coli | √ | SM 9223B-2004 | 34 | MPN/100 mL | | 1 | 42189 | 3/25/15 | 15:26 | RSC |
| AA93578-A | E. Coli Holding Time - IDEXX Colilert | | NA | 2.13 | hours | | 0.00 | 42187 | 3/25/15 | 15:26 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 30, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42189

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

3/30/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: FD - HSM 110
Sample Number: AA93579
Sample Matrix: Non Potable Water

Collection Date/Time: 3/25/15 10:28
Receipt Date/Time: 3/25/15 14:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA93579-A | E. coli | √ | SM 9223B-2004 | 9 | MPN/100 mL | | 1 | 42189 | 3/25/15 | 15:26 | RSC |
| AA93579-A | E. Coli Holding Time - IDEXX Colilert | | NA | 4.97 | hours | | 0.00 | 42187 | 3/25/15 | 15:26 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



March 30, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42189

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

3/30/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable



Calscience



WORK ORDER NUMBER: 15-03-2020

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: SWCA Environmental Consultants

Client Project Name: EAA 27122

Attention: Philip Pearce
6200 UTSA Blvd.
Suite 102
San Antonio, TX 78249-1618

H. Burley FOR

Approved for release on 04/14/2015 by:
Don Burley
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

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Work Order Number: 15-03-2020

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Work Order Narrative

Work Order: 15-03-2020

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Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 03/26/15. They were assigned to Work Order 15-03-2020.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.



Calscience

Sample Summary

| | |
|--|------------------------------------|
| Client: SWCA Environmental Consultants | Work Order: 15-03-2020 |
| 6200 UTSA Blvd., Suite 102 | Project Name: EAA 27122 |
| San Antonio, TX 78249-1618 | PO Number: |
| | Date/Time Received: 03/26/15 10:40 |
| | Number of Containers: 73 |
| Attn: Philip Pearce | |

| Sample Identification | Lab Number | Collection Date and Time | Number of Containers | Matrix |
|-----------------------|--------------|--------------------------|----------------------|---------|
| HSM110 | 15-03-2020-1 | 03/25/15 10:28 | 9 | Aqueous |
| HSM120 | 15-03-2020-2 | 03/25/15 10:51 | 9 | Aqueous |
| HSM130 | 15-03-2020-3 | 03/25/15 11:17 | 9 | Aqueous |
| HSM140 | 15-03-2020-4 | 03/25/15 11:30 | 9 | Aqueous |
| HSM150 | 15-03-2020-5 | 03/25/15 12:25 | 9 | Aqueous |
| HSM160 | 15-03-2020-6 | 03/25/15 12:50 | 9 | Aqueous |
| HSM170 | 15-03-2020-7 | 03/25/15 13:18 | 9 | Aqueous |
| FDHSM110 | 15-03-2020-8 | 03/25/15 10:28 | 9 | Aqueous |
| TB04 | 15-03-2020-9 | 03/25/15 00:00 | 1 | Aqueous |

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Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-03-2020-1-F | 03/25/15 10:28 | Aqueous | ICP 7300 | 03/27/15 | 04/09/15 22:10 | 150327L4FF |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 93.9 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 18.0 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.31 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.7 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.568 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.63 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|
| HSM120 | 15-03-2020-2-F | 03/25/15 10:51 | Aqueous | ICP 7300 | 03/27/15 | 04/09/15 22:16 | 150327L4FF |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 86.6 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 20.0 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.85 | 0.500 | 0.103 | 1.00 | |
| Sodium | 15.2 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.680 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.60 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|
| HSM130 | 15-03-2020-3-F | 03/25/15 11:17 | Aqueous | ICP 7300 | 03/27/15 | 04/09/15 22:18 | 150327L4FF |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 91.5 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 13.6 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.55 | 0.500 | 0.103 | 1.00 | |
| Sodium | 14.6 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.511 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.15 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-03-2020-4-F | 03/25/15 11:30 | Aqueous | ICP 7300 | 03/27/15 | 04/09/15 22:20 | 150327L4FF |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 93.1 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 18.6 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.48 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.574 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.70 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|
| HSM150 | 15-03-2020-5-F | 03/25/15 12:25 | Aqueous | ICP 7300 | 03/27/15 | 04/09/15 22:22 | 150327L4FF |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 93.5 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 18.2 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.29 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.7 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.577 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.60 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|
| HSM160 | 15-03-2020-6-F | 03/25/15 12:50 | Aqueous | ICP 7300 | 03/27/15 | 04/09/15 22:24 | 150327L4FF |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 92.4 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 18.1 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.49 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.9 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.573 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.59 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-03-2020-7-F | 03/25/15 13:18 | Aqueous | ICP 7300 | 03/27/15 | 04/09/15 22:32 | 150327L4FF |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 91.6 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 18.0 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.43 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.3 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.571 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.62 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|----------|----------------|----------------|---------|----------|----------|----------------|------------|
| FDHSM110 | 15-03-2020-8-F | 03/25/15 10:28 | Aqueous | ICP 7300 | 03/27/15 | 04/09/15 22:34 | 150327L4FF |
|----------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 87.4 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 20.1 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.75 | 0.500 | 0.103 | 1.00 | |
| Sodium | 15.3 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.688 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.59 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|-----------------|-----|---------|----------|----------|----------------|------------|
| Method Blank | 099-15-683-1214 | N/A | Aqueous | ICP 7300 | 03/27/15 | 04/09/15 22:07 | 150327L4FF |
|--------------|-----------------|-----|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | ND | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 0.0449 | 0.100 | 0.00336 | 1.00 | J |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-03-2020-1-F | 03/25/15 10:28 | Aqueous | ICP/MS 03 | 03/27/15 | 03/28/15 05:36 | 150327LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0409 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000788 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00193 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000310 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00287 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0411 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.0222 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM120 | 15-03-2020-2-F | 03/25/15 10:51 | Aqueous | ICP/MS 03 | 03/27/15 | 03/28/15 05:40 | 150327LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0382 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000682 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00185 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000351 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0344 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.000220 | 0.00100 | 0.000139 | 1.00 | J |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-03-2020-3-F | 03/25/15 11:17 | Aqueous | ICP/MS 03 | 03/27/15 | 03/28/15 05:43 | 150327LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0418 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000662 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00194 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000422 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00447 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0295 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00148 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-03-2020-4-F | 03/25/15 11:30 | Aqueous | ICP/MS 03 | 03/27/15 | 03/28/15 05:47 | 150327LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0392 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000951 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00208 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000314 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.000711 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0338 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.000545 | 0.00100 | 0.000139 | 1.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM150 | 15-03-2020-5-F | 03/25/15 12:25 | Aqueous | ICP/MS 03 | 03/27/15 | 03/28/15 05:50 | 150327LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0388 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000728 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00189 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000282 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00135 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0288 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.000616 | 0.00100 | 0.000139 | 1.00 | J |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-03-2020-6-F | 03/25/15 12:50 | Aqueous | ICP/MS 03 | 03/27/15 | 03/28/15 05:54 | 150327LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0379 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000978 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00207 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000314 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00520 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0342 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00128 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-03-2020-7-F | 03/25/15 13:18 | Aqueous | ICP/MS 03 | 03/27/15 | 03/28/15 05:58 | 150327LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0398 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000824 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00180 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000334 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0348 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00167 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM110 | 15-03-2020-8-F | 03/25/15 10:28 | Aqueous | ICP/MS 03 | 03/27/15 | 03/28/15 06:01 | 150327LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0429 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000673 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00203 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0485 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.0220 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-771 | N/A | Aqueous | ICP/MS 03 | 03/27/15 | 03/28/15 04:39 | 150327LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-03-2020-1-F | 03/25/15 10:28 | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 17:16 | 150401L04F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HSM120 | 15-03-2020-2-F | 03/25/15 10:51 | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 17:19 | 150401L04F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HSM130 | 15-03-2020-3-F | 03/25/15 11:17 | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 17:21 | 150401L04F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HSM140 | 15-03-2020-4-F | 03/25/15 11:30 | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 17:23 | 150401L04F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HSM150 | 15-03-2020-5-F | 03/25/15 12:25 | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 17:30 | 150401L04F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HSM160 | 15-03-2020-6-F | 03/25/15 12:50 | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 17:32 | 150401L04F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-03-2020-7-F | 03/25/15 13:18 | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 17:34 | 150401L04F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|----------|----------------|----------------|---------|------------|----------|----------------|------------|
| FDHSM110 | 15-03-2020-8-F | 03/25/15 10:28 | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 17:37 | 150401L04F |
|----------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|-----|---------|------------|----------|----------------|------------|
| Method Blank | 099-15-763-527 | N/A | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 16:36 | 150401L04F |
|--------------|----------------|-----|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|-----------------------|-----------------------|----------------|-----------------------|-----------------|-----------------------|-------------------|
| HSM110 | 15-03-2020-1-I | 03/25/15 10:28 | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 10:15 | 150328L13 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 16 | | 9.8 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 42 | | 31-200 | | | |
| HSM120 | 15-03-2020-2-I | 03/25/15 10:51 | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 10:23 | 150328L13 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 41 | | 31-200 | | | |
| HSM130 | 15-03-2020-3-I | 03/25/15 11:17 | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 10:32 | 150328L13 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 9.9 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 44 | | 31-200 | | | |
| HSM140 | 15-03-2020-4-I | 03/25/15 11:30 | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 10:40 | 150328L13 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 9.8 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 47 | | 31-200 | | | |
| HSM150 | 15-03-2020-5-I | 03/25/15 12:25 | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 10:48 | 150328L13 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 14 | | 9.7 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 45 | | 31-200 | | | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|-----------------------|-----------------------|----------------|-----------------------|-----------------|-----------------------|-------------------|
| HSM160 | 15-03-2020-6-I | 03/25/15 12:50 | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 10:57 | 150328L13 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 42 | | 31-200 | | | |
| HSM170 | 15-03-2020-7-I | 03/25/15 13:18 | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 11:05 | 150328L13 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 9.8 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 47 | | 31-200 | | | |
| FDHSM110 | 15-03-2020-8-I | 03/25/15 10:28 | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 11:13 | 150328L13 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 10 | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 43 | | 31-200 | | | |
| Method Blank | 099-16-376-12 | N/A | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 10:07 | 150328L13 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 47 | | 31-200 | | | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-03-2020-1-I | 03/25/15 10:28 | Aqueous | GC 44 | 03/27/15 | 03/30/15 20:23 | 150327L02A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 102 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 110 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM120 | 15-03-2020-2-I | 03/25/15 10:51 | Aqueous | GC 44 | 03/27/15 | 03/30/15 20:38 | 150327L02A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 109 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 118 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-03-2020-3-I | 03/25/15 11:17 | Aqueous | GC 44 | 03/27/15 | 03/31/15 12:51 | 150327L02A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 111 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 115 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-03-2020-4-I | 03/25/15 11:30 | Aqueous | GC 44 | 03/27/15 | 03/31/15 13:05 | 150327L02A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 98 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 98 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM150 | 15-03-2020-5-I | 03/25/15 12:25 | Aqueous | GC 44 | 03/27/15 | 03/31/15 15:21 | 150327L02A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 93 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 103 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-03-2020-6-I | 03/25/15 12:50 | Aqueous | GC 44 | 03/27/15 | 03/31/15 15:36 | 150327L02A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 103 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 109 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-03-2020-7-I | 03/25/15 13:18 | Aqueous | GC 44 | 03/27/15 | 03/31/15 18:07 | 150327L02A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 93 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 98 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM110 | 15-03-2020-8-I | 03/25/15 10:28 | Aqueous | GC 44 | 03/27/15 | 03/31/15 18:21 | 150327L02A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 99 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 108 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-796 | N/A | Aqueous | GC 44 | 03/27/15 | 03/30/15 13:14 | 150327L02A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 98 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 101 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

Page 1 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-03-2020-1-I | 03/25/15 10:28 | Aqueous | GC 31 | 03/27/15 | 03/30/15 14:29 | 150327L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 99 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

| HSM120 | 15-03-2020-2-I | 03/25/15 10:51 | Aqueous | GC 31 | 03/27/15 | 03/30/15 14:48 | 150327L03 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 106 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 90 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-03-2020-3-I | 03/25/15 11:17 | Aqueous | GC 31 | 03/27/15 | 03/30/15 15:07 | 150327L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 99 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 87 | 50-135 | |

| HSM140 | 15-03-2020-4-I | 03/25/15 11:30 | Aqueous | GC 31 | 03/27/15 | 03/30/15 15:26 | 150327L03 |
|--------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|--------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 87 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 77 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

Page 3 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM150 | 15-03-2020-5-I | 03/25/15 12:25 | Aqueous | GC 31 | 03/27/15 | 03/30/15 15:45 | 150327L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 93 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 85 | 50-135 | |

| HSM160 | 15-03-2020-6-I | 03/25/15 12:50 | Aqueous | GC 31 | 03/27/15 | 03/30/15 16:05 | 150327L03 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 102 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 90 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-03-2020-7-I | 03/25/15 13:18 | Aqueous | GC 31 | 03/27/15 | 03/30/15 16:34 | 150327L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 93 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 85 | 50-135 | |

| FDHSM110 | 15-03-2020-8-I | 03/25/15 10:28 | Aqueous | GC 31 | 03/27/15 | 03/30/15 16:53 | 150327L03 |
|----------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|----------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 95 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-533-1019 | N/A | Aqueous | GC 31 | 03/27/15 | 03/30/15 13:13 | 150327L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 84 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 80 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-03-2020-1-I | 03/25/15 10:28 | Aqueous | GC 35 | 03/27/15 | 03/31/15 15:21 | 150327L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 102 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM120 | 15-03-2020-2-I | 03/25/15 10:51 | Aqueous | GC 35 | 03/27/15 | 03/31/15 16:07 | 150327L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0049 | 0.0028 | 1.00 | |
| Bolstar | ND | 0.0049 | 0.0028 | 1.00 | |
| Chlorpyrifos | ND | 0.0049 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0049 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0049 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0049 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.0098 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0049 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0049 | 0.0028 | 1.00 | |
| Fenthion | ND | 0.0049 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0049 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0049 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0049 | 0.0027 | 1.00 | |
| Naled | ND | 0.039 | 0.019 | 1.00 | |
| Phorate | ND | 0.0049 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0049 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0086 | 1.00 | |
| Tokuthion | ND | 0.0049 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0049 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0049 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 92 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-03-2020-3-I | 03/25/15 11:17 | Aqueous | GC 35 | 03/27/15 | 03/31/15 16:53 | 150327L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0048 | 0.0027 | 1.00 | |
| Bolstar | ND | 0.0048 | 0.0027 | 1.00 | |
| Chlorpyrifos | ND | 0.0048 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0048 | 0.0023 | 1.00 | |
| Diazinon | ND | 0.0048 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0048 | 0.0035 | 1.00 | |
| Disulfoton | ND | 0.0096 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0048 | 0.0024 | 1.00 | |
| Fensulfothion | ND | 0.0048 | 0.0027 | 1.00 | |
| Fenthion | ND | 0.0048 | 0.0025 | 1.00 | |
| Merphos | ND | 0.0048 | 0.0025 | 1.00 | |
| Methyl Parathion | ND | 0.0048 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0048 | 0.0026 | 1.00 | |
| Naled | ND | 0.038 | 0.019 | 1.00 | |
| Phorate | ND | 0.0048 | 0.0024 | 1.00 | |
| Ronnel | ND | 0.0048 | 0.0031 | 1.00 | |
| Stirophos | ND | 0.019 | 0.0085 | 1.00 | |
| Tokuthion | ND | 0.0048 | 0.0027 | 1.00 | |
| Trichloronate | ND | 0.0048 | 0.0020 | 1.00 | |
| Demeton-o/s | ND | 0.0048 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 92 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 4 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-03-2020-4-I | 03/25/15 11:30 | Aqueous | GC 35 | 03/27/15 | 03/31/15 17:40 | 150327L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0028 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0028 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.0099 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0028 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0087 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 94 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM150 | 15-03-2020-5-I | 03/25/15 12:25 | Aqueous | GC 35 | 03/27/15 | 03/31/15 18:26 | 150327L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 86 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-03-2020-6-I | 03/25/15 12:50 | Aqueous | GC 35 | 03/27/15 | 03/31/15 19:12 | 150327L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0048 | 0.0027 | 1.00 | |
| Bolstar | ND | 0.0048 | 0.0027 | 1.00 | |
| Chlorpyrifos | ND | 0.0048 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0048 | 0.0023 | 1.00 | |
| Diazinon | ND | 0.0048 | 0.0027 | 1.00 | |
| Dichlorvos | ND | 0.0048 | 0.0035 | 1.00 | |
| Disulfoton | ND | 0.0095 | 0.0024 | 1.00 | |
| Ethoprop | ND | 0.0048 | 0.0024 | 1.00 | |
| Fensulfothion | ND | 0.0048 | 0.0027 | 1.00 | |
| Fenthion | ND | 0.0048 | 0.0025 | 1.00 | |
| Merphos | ND | 0.0048 | 0.0025 | 1.00 | |
| Methyl Parathion | ND | 0.0048 | 0.0028 | 1.00 | |
| Mevinphos | ND | 0.0048 | 0.0026 | 1.00 | |
| Naled | ND | 0.038 | 0.018 | 1.00 | |
| Phorate | ND | 0.0048 | 0.0024 | 1.00 | |
| Ronnel | ND | 0.0048 | 0.0031 | 1.00 | |
| Stirophos | ND | 0.019 | 0.0084 | 1.00 | |
| Tokuthion | ND | 0.0048 | 0.0027 | 1.00 | |
| Trichloronate | ND | 0.0048 | 0.0020 | 1.00 | |
| Demeton-o/s | ND | 0.0048 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 92 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-03-2020-7-I | 03/25/15 13:18 | Aqueous | GC 35 | 03/27/15 | 03/31/15 19:58 | 150327L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0049 | 0.0028 | 1.00 | |
| Bolstar | ND | 0.0049 | 0.0028 | 1.00 | |
| Chlorpyrifos | ND | 0.0049 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0049 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0049 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0049 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.0098 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0049 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0049 | 0.0028 | 1.00 | |
| Fenthion | ND | 0.0049 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0049 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0049 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0049 | 0.0027 | 1.00 | |
| Naled | ND | 0.039 | 0.019 | 1.00 | |
| Phorate | ND | 0.0049 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0049 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0086 | 1.00 | |
| Tokuthion | ND | 0.0049 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0049 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0049 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 83 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| FDHSM110 | 15-03-2020-8-I | 03/25/15 10:28 | Aqueous | GC 35 | 03/27/15 | 03/31/15 20:44 | 150327L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 81 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 9 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-91 | N/A | Aqueous | GC 35 | 03/27/15 | 03/31/15 12:17 | 150327L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 83 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

Page 1 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-03-2020-1-J | 03/25/15 10:28 | Aqueous | GC 40 | 03/27/15 | 04/06/15 13:42 | 150327L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 80 | 0-123 | |

| HSM120 | 15-03-2020-2-I | 03/25/15 10:51 | Aqueous | GC 40 | 03/27/15 | 04/06/15 14:05 | 150327L05 |
|--------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|--------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 87 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-03-2020-3-I | 03/25/15 11:17 | Aqueous | GC 40 | 03/27/15 | 04/06/15 14:28 | 150327L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 81 | 0-123 | |

| HSM140 | 15-03-2020-4-I | 03/25/15 11:30 | Aqueous | GC 40 | 03/27/15 | 04/06/15 14:51 | 150327L05 |
|--------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|--------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 65 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM150 | 15-03-2020-5-I | 03/25/15 12:25 | Aqueous | GC 40 | 03/27/15 | 04/06/15 15:14 | 150327L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 91 | 0-123 | |

| HSM160 | 15-03-2020-6-I | 03/25/15 12:50 | Aqueous | GC 40 | 03/27/15 | 04/06/15 15:37 | 150327L05 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 85 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-03-2020-7-I | 03/25/15 13:18 | Aqueous | GC 40 | 03/27/15 | 04/06/15 16:00 | 150327L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 80 | 0-123 | |

| FDHSM110 | 15-03-2020-8-I | 03/25/15 10:28 | Aqueous | GC 40 | 03/27/15 | 04/06/15 16:23 | 150327L05 |
|----------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|----------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 95 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-034-640 | N/A | Aqueous | GC 40 | 03/27/15 | 03/31/15 16:04 | 150327L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 94 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-03-2020-1-G | 03/25/15 10:28 | Aqueous | GC/MS CCC | 03/27/15 | 03/30/15 17:02 | 150327L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 87 | 33-120 | |
| 2-Fluorophenol | 86 | 24-120 | |
| Nitrobenzene-d5 | 81 | 38-120 | |
| p-Terphenyl-d14 | 81 | 41-137 | |
| Phenol-d6 | 88 | 16-120 | |
| 2,4,6-Tribromophenol | 85 | 27-159 | |


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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Method: EPA 8270C
Units: ug/L

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM120 | 15-03-2020-2-G | 03/25/15 10:51 | Aqueous | GC/MS CCC | 03/27/15 | 03/30/15 17:20 | 150327L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 80 | 33-120 | |
| 2-Fluorophenol | 76 | 24-120 | |
| Nitrobenzene-d5 | 71 | 38-120 | |
| p-Terphenyl-d14 | 73 | 41-137 | |
| Phenol-d6 | 78 | 16-120 | |
| 2,4,6-Tribromophenol | 74 | 27-159 | |


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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-03-2020-3-G | 03/25/15 11:17 | Aqueous | GC/MS CCC | 03/27/15 | 03/30/15 17:38 | 150327L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 80 | 33-120 | |
| 2-Fluorophenol | 77 | 24-120 | |
| Nitrobenzene-d5 | 72 | 38-120 | |
| p-Terphenyl-d14 | 75 | 41-137 | |
| Phenol-d6 | 79 | 16-120 | |
| 2,4,6-Tribromophenol | 74 | 27-159 | |


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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-03-2020-4-G | 03/25/15 11:30 | Aqueous | GC/MS CCC | 03/27/15 | 03/30/15 17:56 | 150327L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 82 | 33-120 | |
| 2-Fluorophenol | 78 | 24-120 | |
| Nitrobenzene-d5 | 73 | 38-120 | |
| p-Terphenyl-d14 | 75 | 41-137 | |
| Phenol-d6 | 81 | 16-120 | |
| 2,4,6-Tribromophenol | 76 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM150 | 15-03-2020-5-G | 03/25/15 12:25 | Aqueous | GC/MS CCC | 03/27/15 | 03/30/15 18:14 | 150327L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 90 | 33-120 | |
| 2-Fluorophenol | 83 | 24-120 | |
| Nitrobenzene-d5 | 81 | 38-120 | |
| p-Terphenyl-d14 | 86 | 41-137 | |
| Phenol-d6 | 84 | 16-120 | |
| 2,4,6-Tribromophenol | 83 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-03-2020-6-G | 03/25/15 12:50 | Aqueous | GC/MS CCC | 03/27/15 | 03/30/15 18:32 | 150327L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 85 | 33-120 | |
| 2-Fluorophenol | 82 | 24-120 | |
| Nitrobenzene-d5 | 77 | 38-120 | |
| p-Terphenyl-d14 | 79 | 41-137 | |
| Phenol-d6 | 84 | 16-120 | |
| 2,4,6-Tribromophenol | 80 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-03-2020-7-G | 03/25/15 13:18 | Aqueous | GC/MS CCC | 03/27/15 | 03/30/15 18:50 | 150327L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 90 | 33-120 | |
| 2-Fluorophenol | 87 | 24-120 | |
| Nitrobenzene-d5 | 81 | 38-120 | |
| p-Terphenyl-d14 | 81 | 41-137 | |
| Phenol-d6 | 88 | 16-120 | |
| 2,4,6-Tribromophenol | 81 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM110 | 15-03-2020-8-G | 03/25/15 10:28 | Aqueous | GC/MS CCC | 03/27/15 | 03/30/15 19:08 | 150327L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 83 | 33-120 | |
| 2-Fluorophenol | 78 | 24-120 | |
| Nitrobenzene-d5 | 75 | 38-120 | |
| p-Terphenyl-d14 | 78 | 41-137 | |
| Phenol-d6 | 79 | 16-120 | |
| 2,4,6-Tribromophenol | 77 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-4016 | N/A | Aqueous | GC/MS CCC | 03/27/15 | 03/27/15 20:24 | 150327L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 85 | 33-120 | |
| 2-Fluorophenol | 73 | 24-120 | |
| Nitrobenzene-d5 | 78 | 38-120 | |
| p-Terphenyl-d14 | 86 | 41-137 | |
| Phenol-d6 | 76 | 16-120 | |
| 2,4,6-Tribromophenol | 79 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 1 of 20

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-03-2020-1-A | 03/25/15 10:28 | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 11:49 | 150404L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 88 | 80-120 | |
| Dibromofluoromethane | 107 | 78-126 | |
| 1,2-Dichloroethane-d4 | 108 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM120 | 15-03-2020-2-A | 03/25/15 10:51 | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 15:23 | 150404L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 92 | 80-120 | |
| Dibromofluoromethane | 108 | 78-126 | |
| 1,2-Dichloroethane-d4 | 111 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 5 of 20

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-03-2020-3-A | 03/25/15 11:17 | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 15:50 | 150404L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 90 | 80-120 | |
| Dibromofluoromethane | 105 | 78-126 | |
| 1,2-Dichloroethane-d4 | 109 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-03-2020-4-A | 03/25/15 11:30 | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 16:17 | 150404L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 89 | 80-120 | |
| Dibromofluoromethane | 111 | 78-126 | |
| 1,2-Dichloroethane-d4 | 112 | 75-135 | |
| Toluene-d8 | 104 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 9 of 20

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM150 | 15-03-2020-5-A | 03/25/15 12:25 | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 16:43 | 150404L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 10 of 20

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 92 | 80-120 | |
| Dibromofluoromethane | 109 | 78-126 | |
| 1,2-Dichloroethane-d4 | 113 | 75-135 | |
| Toluene-d8 | 105 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-03-2020-6-A | 03/25/15 12:50 | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 17:10 | 150404L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 89 | 80-120 | |
| Dibromofluoromethane | 107 | 78-126 | |
| 1,2-Dichloroethane-d4 | 110 | 75-135 | |
| Toluene-d8 | 106 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 13 of 20

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-03-2020-7-A | 03/25/15 13:18 | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 17:37 | 150404L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 90 | 80-120 | |
| Dibromofluoromethane | 110 | 78-126 | |
| 1,2-Dichloroethane-d4 | 110 | 75-135 | |
| Toluene-d8 | 103 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 15 of 20

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM110 | 15-03-2020-8-A | 03/25/15 10:28 | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 18:04 | 150404L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 88 | 80-120 | |
| Dibromofluoromethane | 113 | 78-126 | |
| 1,2-Dichloroethane-d4 | 115 | 75-135 | |
| Toluene-d8 | 104 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 17 of 20

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| TB04 | 15-03-2020-9-A | 03/25/15 00:00 | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 18:30 | 150404L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 89 | 80-120 | |
| Dibromofluoromethane | 109 | 78-126 | |
| 1,2-Dichloroethane-d4 | 112 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-16776 | N/A | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 11:20 | 150404L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 110 | 78-126 | |
| 1,2-Dichloroethane-d4 | 110 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

03/26/15

Work Order:

15-03-2020

Page 1 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HSM110 | 15-03-2020-1 | 03/25/15 10:28 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.23 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Chloride (24) | 24 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Bromide (24) | 0.16 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Nitrate (as N) (24) | 0.36 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Sulfate (24) | 33 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 04/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 244 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 244 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Solids, Total Dissolved (24) | 365 | 1.00 | 0.870 | 1.00 | | mg/L | 03/31/15 | 03/31/15 | SM 2540 C |
| Solids, Total Suspended (24) | 5.9 | 1.0 | 0.83 | 1.00 | | mg/L | 03/30/15 | 03/30/15 | SM 2540 D |
| pH (24) | 7.29 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/26/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 04/03/15 | 04/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 1.4 | 1.0 | 0.47 | 2.00 | | mg/L | 04/08/15 | 04/09/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 1.9 | 0.50 | 0.24 | 1.00 | | mg/L | 03/31/15 | 04/01/15 | SM 5310 B |

| | | | |
|---------------|---------------------|-----------------------|----------------|
| HSM120 | 15-03-2020-2 | 03/25/15 10:51 | Aqueous |
|---------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.22 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Chloride (24) | 19 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Bromide (24) | 0.12 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.2 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Sulfate (24) | 27 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 04/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 248 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 248 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Solids, Total Dissolved (24) | 355 | 1.00 | 0.870 | 1.00 | | mg/L | 03/31/15 | 03/31/15 | SM 2540 C |
| Solids, Total Suspended (24) | 2.0 | 1.0 | 0.83 | 1.00 | | mg/L | 03/30/15 | 03/30/15 | SM 2540 D |
| pH (24) | 7.33 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/26/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 04/03/15 | 04/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 2.2 | 1.0 | 0.47 | 2.00 | | mg/L | 04/08/15 | 04/09/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.3 | 0.50 | 0.24 | 1.00 | | mg/L | 03/31/15 | 04/01/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

03/26/15

Work Order:

15-03-2020

Page 2 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HSM130 | 15-03-2020-3 | 03/25/15 11:17 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.22 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Chloride (24) | 23 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Bromide (24) | 0.12 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Nitrate (as N) (24) | 2.0 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Sulfate (24) | 33 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 04/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 273 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 273 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Solids, Total Dissolved (24) | 345 | 1.00 | 0.870 | 1.00 | | mg/L | 03/31/15 | 03/31/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 03/30/15 | 03/30/15 | SM 2540 D |
| pH (24) | 7.42 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/26/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 04/03/15 | 04/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 3.3 | 2.5 | 1.2 | 5.00 | | mg/L | 04/08/15 | 04/09/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 1.1 | 0.50 | 0.24 | 1.00 | | mg/L | 03/31/15 | 04/01/15 | SM 5310 B |

| | | | |
|---------------|---------------------|-----------------------|----------------|
| HSM140 | 15-03-2020-4 | 03/25/15 11:30 | Aqueous |
|---------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.22 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Chloride (24) | 19 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.1 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Sulfate (24) | 28 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 04/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 262 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 262 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Solids, Total Dissolved (24) | 315 | 1.00 | 0.870 | 1.00 | | mg/L | 03/31/15 | 03/31/15 | SM 2540 C |
| Solids, Total Suspended (24) | 2.2 | 1.0 | 0.83 | 1.00 | | mg/L | 03/30/15 | 03/30/15 | SM 2540 D |
| pH (24) | 7.62 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/26/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 04/03/15 | 04/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 1.0 | 0.47 | 2.00 | | mg/L | 04/08/15 | 04/09/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 1.9 | 0.50 | 0.24 | 1.00 | | mg/L | 03/31/15 | 04/01/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

03/26/15

Work Order:

15-03-2020

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HSM150 | 15-03-2020-5 | 03/25/15 12:25 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.23 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Chloride (24) | 19 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Bromide (24) | 0.11 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.1 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Sulfate (24) | 28 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 04/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 258 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 258 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Solids, Total Dissolved (24) | 345 | 1.00 | 0.870 | 1.00 | | mg/L | 03/31/15 | 03/31/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 03/30/15 | 03/30/15 | SM 2540 D |
| pH (24) | 7.71 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/26/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 04/03/15 | 04/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 2.8 | 1.0 | 0.47 | 2.00 | | mg/L | 04/08/15 | 04/09/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 03/31/15 | 04/01/15 | SM 5310 B |

| | | | |
|---------------|---------------------|-----------------------|----------------|
| HSM160 | 15-03-2020-6 | 03/25/15 12:50 | Aqueous |
|---------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.21 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Chloride (24) | 19 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Bromide (24) | 0.11 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.1 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Sulfate (24) | 28 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 04/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 256 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 256 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Solids, Total Dissolved (24) | 300 | 1.00 | 0.870 | 1.00 | | mg/L | 03/31/15 | 03/31/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.8 | 1.0 | 0.83 | 1.00 | | mg/L | 03/30/15 | 03/30/15 | SM 2540 D |
| pH (24) | 7.82 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/26/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 04/03/15 | 04/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 2.4 | 1.0 | 0.47 | 2.00 | | mg/L | 04/08/15 | 04/09/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 0.96 | 0.50 | 0.24 | 1.00 | | mg/L | 03/31/15 | 04/01/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

03/26/15

Work Order:

15-03-2020

Page 4 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HSM170 | 15-03-2020-7 | 03/25/15 13:18 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.22 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Chloride (24) | 19 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Bromide (24) | 0.11 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.1 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Sulfate (24) | 28 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 04/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 255 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 255 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Solids, Total Dissolved (24) | 355 | 1.00 | 0.870 | 1.00 | | mg/L | 03/31/15 | 03/31/15 | SM 2540 C |
| Solids, Total Suspended (24) | 2.7 | 1.0 | 0.83 | 1.00 | | mg/L | 03/30/15 | 03/30/15 | SM 2540 D |
| pH (24) | 7.87 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/26/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 04/03/15 | 04/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 1.5 | 1.0 | 0.47 | 2.00 | | mg/L | 04/08/15 | 04/09/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.2 | 0.50 | 0.24 | 1.00 | | mg/L | 03/31/15 | 04/01/15 | SM 5310 B |

| | | | |
|-----------------|---------------------|-----------------------|----------------|
| FDHSM110 | 15-03-2020-8 | 03/25/15 10:28 | Aqueous |
|-----------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.23 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Chloride (24) | 24 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Bromide (24) | 0.15 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Nitrate (as N) (24) | 0.36 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Sulfate (24) | 33 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/27/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.022 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 04/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 253 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 253 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Solids, Total Dissolved (24) | 385 | 1.00 | 0.870 | 1.00 | | mg/L | 03/31/15 | 03/31/15 | SM 2540 C |
| Solids, Total Suspended (24) | 9.0 | 1.0 | 0.83 | 1.00 | | mg/L | 03/30/15 | 03/30/15 | SM 2540 D |
| pH (24) | 7.57 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 03/26/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 04/03/15 | 04/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 4.6 | 2.5 | 1.2 | 5.00 | | mg/L | 04/08/15 | 04/09/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 3.1 | 0.50 | 0.24 | 1.00 | | mg/L | 03/31/15 | 04/01/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

03/26/15

Work Order:

15-03-2020

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|-------------------|---------------------|----------------|
| Method Blank | | N/A | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|-------|---------------|---------------|-----------------|
| Fluoride (24) | ND | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Chloride (24) | ND | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Nitrate (as N) (24) | ND | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Sulfate (24) | ND | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 03/26/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 04/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 04/08/15 | SM 2320B |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.87 | 1.00 | | mg/L | 03/31/15 | 03/31/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 03/30/15 | 03/30/15 | SM 2540 D |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 04/03/15 | 04/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 04/08/15 | 04/09/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 03/31/15 | 04/01/15 | SM 5310 B |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-03-2018-3 | Sample | Aqueous | IC 15 | N/A | 03/26/15 21:41 | 150326S02A |
| 15-03-2018-3 | Matrix Spike | Aqueous | IC 15 | N/A | 03/26/15 23:39 | 150326S02A |
| 15-03-2018-3 | Matrix Spike Duplicate | Aqueous | IC 15 | N/A | 03/26/15 23:56 | 150326S02A |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.1694 | 250.0 | 270.6 | 108 | 271.4 | 108 | 80-120 | 0 | 0-20 | |
| Chloride | 75.80 | 5000 | 5162 | 102 | 5139 | 101 | 80-120 | 0 | 0-20 | |
| Bromide | 0.1331 | 500.0 | 506.2 | 101 | 506.5 | 101 | 80-120 | 0 | 0-20 | |
| Nitrate (as N) | ND | 500.0 | 516.9 | 103 | 515.3 | 103 | 80-120 | 0 | 0-20 | |
| Sulfate | 217.8 | 5000 | 5289 | 101 | 5252 | 101 | 80-120 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM110 | Sample | Aqueous | ACA 1 | N/A | 04/09/15 14:45 | 150409SO1 |
| HSM110 | Matrix Spike | Aqueous | ACA 1 | N/A | 04/09/15 14:45 | 150409SO1 |
| HSM110 | Matrix Spike Duplicate | Aqueous | ACA 1 | N/A | 04/09/15 14:45 | 150409SO1 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | ND | 0.2000 | 0.2114 | 106 | 0.2133 | 107 | 90-110 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM110 | Sample | Aqueous | TOC 8 | 04/08/15 | 04/09/15 02:26 | F0408TOCS1 |
| HSM110 | Matrix Spike | Aqueous | TOC 8 | 04/08/15 | 04/09/15 02:26 | F0408TOCS1 |
| HSM110 | Matrix Spike Duplicate | Aqueous | TOC 8 | 04/08/15 | 04/09/15 02:26 | F0408TOCS1 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Carbon, Total Organic | 1.356 | 20.00 | 19.30 | 90 | 18.76 | 87 | 31-145 | 3 | 0-23 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|---------------------|-----------------|------------|---------------|-------------------|
| HSM110 | Sample | Aqueous | TOC 8 | 03/31/15 | 04/01/15 00:32 | F0331DOCS1 | | | | |
| HSM110 | Matrix Spike | Aqueous | TOC 8 | 03/31/15 | 04/01/15 00:32 | F0331DOCS1 | | | | |
| HSM110 | Matrix Spike Duplicate | Aqueous | TOC 8 | 03/31/15 | 04/01/15 00:32 | F0331DOCS1 | | | | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Carbon, Dissolved Organic | 1.900 | 10.00 | 11.20 | 93 | 10.60 | 87 | 31-145 | 6 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM110 | Sample | Aqueous | ICP 7300 | 03/27/15 | 04/09/15 22:10 | 150327SA4A |
| HSM110 | Matrix Spike | Aqueous | ICP 7300 | 03/27/15 | 04/09/15 22:12 | 150327SA4A |
| HSM110 | Matrix Spike Duplicate | Aqueous | ICP 7300 | 03/27/15 | 04/09/15 22:14 | 150327SA4A |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Calcium | 93.90 | 0.5000 | 91.60 | 4X | 91.56 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 17.97 | 0.5000 | 18.08 | 4X | 18.28 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 1.310 | 5.000 | 6.076 | 95 | 6.142 | 97 | 83-131 | 1 | 0-7 | |
| Sodium | 12.69 | 5.000 | 17.64 | 99 | 17.52 | 97 | 73-127 | 1 | 0-9 | |
| Strontium | 0.5683 | 0.5000 | 1.070 | 100 | 1.066 | 100 | 81-123 | 0 | 0-6 | |
| Silicon | 3.627 | 0.5000 | 3.891 | 4X | 3.987 | 4X | 24-180 | 4X | 0-15 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|---------------------------|------------------------|-------------|----------|------------|---------------|----------------|---------------------|-----|--------|------------|
| HSM110 | Sample | | Aqueous | ICP/MS 03 | 03/27/15 | 03/28/15 05:36 | 150327SA2 | | | |
| HSM110 | Matrix Spike | | Aqueous | ICP/MS 03 | 03/27/15 | 03/28/15 04:46 | 150327SA2 | | | |
| HSM110 | Matrix Spike Duplicate | | Aqueous | ICP/MS 03 | 03/27/15 | 03/28/15 04:50 | 150327SA2 | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.09764 | 98 | 0.09578 | 96 | 85-133 | 2 | 0-11 | |
| Arsenic | ND | 0.1000 | 0.09620 | 96 | 0.09635 | 96 | 73-127 | 0 | 0-11 | |
| Barium | 0.04094 | 0.1000 | 0.1431 | 102 | 0.1401 | 99 | 74-128 | 2 | 0-10 | |
| Beryllium | ND | 0.1000 | 0.09566 | 96 | 0.09402 | 94 | 56-122 | 2 | 0-11 | |
| Cadmium | ND | 0.1000 | 0.09273 | 93 | 0.09150 | 91 | 84-114 | 1 | 0-8 | |
| Chromium | ND | 0.1000 | 0.1084 | 108 | 0.1077 | 108 | 73-133 | 1 | 0-11 | |
| Copper | ND | 0.1000 | 0.09824 | 98 | 0.09708 | 97 | 72-108 | 1 | 0-10 | |
| Lead | ND | 0.1000 | 0.1090 | 109 | 0.1064 | 106 | 79-121 | 2 | 0-10 | |
| Nickel | 0.001933 | 0.1000 | 0.09848 | 97 | 0.09765 | 96 | 68-122 | 1 | 0-10 | |
| Selenium | ND | 0.1000 | 0.08043 | 80 | 0.08116 | 81 | 59-125 | 1 | 0-12 | |
| Silver | ND | 0.05000 | 0.04849 | 97 | 0.04799 | 96 | 68-128 | 1 | 0-14 | |
| Thallium | ND | 0.1000 | 0.1081 | 108 | 0.1049 | 105 | 73-121 | 3 | 0-11 | |
| Zinc | ND | 0.1000 | 0.08774 | 88 | 0.08583 | 86 | 43-145 | 2 | 0-39 | |
| Aluminum | ND | 0.1000 | 0.1102 | 110 | 0.1102 | 110 | 47-161 | 0 | 0-24 | |
| Iron | ND | 5.100 | 5.426 | 106 | 5.393 | 106 | 27-201 | 1 | 0-24 | |
| Manganese | 0.02222 | 0.1000 | 0.1247 | 103 | 0.1229 | 101 | 72-126 | 2 | 0-42 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 7470A Total
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-03-2120-1 | Sample | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 16:40 | 150401S04 |
| 15-03-2120-1 | Matrix Spike | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 16:43 | 150401S04 |
| 15-03-2120-1 | Matrix Spike Duplicate | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 16:45 | 150401S04 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | ND | 0.01000 | 0.01095 | 109 | 0.01059 | 106 | 57-141 | 3 | 0-10 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM170 | Sample | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 11:05 | 150328S13 |
| HSM170 | Matrix Spike | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 11:22 | 150328S13 |
| HSM170 | Matrix Spike Duplicate | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 11:30 | 150328S13 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Caffeine | ND | 100.0 | 107.6 | 108 | 109.9 | 110 | 70-130 | 2 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM110 | Sample | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 11:49 | 150404S002 |
| HSM110 | Matrix Spike | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 12:16 | 150404S002 |
| HSM110 | Matrix Spike Duplicate | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 12:42 | 150404S002 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Benzene | ND | 50.00 | 48.59 | 97 | 48.55 | 97 | 74-122 | 0 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 65.82 | 132 | 67.48 | 135 | 60-144 | 2 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 50.88 | 102 | 48.86 | 98 | 73-120 | 4 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 50.28 | 101 | 50.87 | 102 | 80-122 | 1 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 48.67 | 97 | 47.92 | 96 | 70-120 | 2 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 49.78 | 100 | 50.54 | 101 | 64-142 | 2 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 48.65 | 97 | 48.95 | 98 | 52-136 | 1 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 51.82 | 104 | 51.06 | 102 | 77-125 | 1 | 0-24 | |
| Toluene | ND | 50.00 | 50.04 | 100 | 50.69 | 101 | 72-126 | 1 | 0-23 | |
| Trichloroethene | ND | 50.00 | 48.31 | 97 | 49.30 | 99 | 74-128 | 2 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 46.71 | 93 | 49.37 | 99 | 67-133 | 6 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 102.1 | 102 | 99.99 | 100 | 63-129 | 2 | 0-25 | |
| o-Xylene | ND | 50.00 | 51.77 | 104 | 51.01 | 102 | 62-128 | 1 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 52.14 | 104 | 55.33 | 111 | 68-134 | 6 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|---------------------|--------------------|------------------|-----------------------|-----------------------|-----------------------|
| HSM110 | Sample | Aqueous | ICP/MS 03 | 03/27/15 00:00 | 03/28/15 05:36 | 150327SA2 |
| HSM110 | PDS | Aqueous | ICP/MS 03 | 03/27/15 00:00 | 03/30/15 13:56 | 150327SA2 |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>PDS Conc.</u> | <u>PDS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Antimony | ND | 0.1000 | 0.09903 | 99 | 75-125 | |
| Arsenic | ND | 0.1000 | 0.09484 | 95 | 75-125 | |
| Barium | 0.04094 | 0.1000 | 0.1422 | 101 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.09830 | 98 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.09563 | 96 | 75-125 | |
| Chromium | ND | 0.1000 | 0.1058 | 106 | 75-125 | |
| Copper | ND | 0.1000 | 0.09546 | 95 | 75-125 | |
| Lead | ND | 0.1000 | 0.1085 | 109 | 75-125 | |
| Nickel | 0.001933 | 0.1000 | 0.09740 | 95 | 75-125 | |
| Selenium | ND | 0.1000 | 0.08715 | 87 | 75-125 | |
| Silver | ND | 0.05000 | 0.04873 | 97 | 75-125 | |
| Thallium | ND | 0.1000 | 0.1071 | 107 | 75-125 | |
| Zinc | ND | 0.1000 | 0.09145 | 91 | 75-125 | |
| Aluminum | ND | 0.1000 | 0.1129 | 113 | 75-125 | |
| Iron | ND | 5.100 | 5.122 | 100 | 75-125 | |
| Manganese | 0.02222 | 0.1000 | 0.1272 | 105 | 75-125 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-04-0164-1 | Sample | Aqueous | PH1/BUR03 | N/A | 04/08/15 21:35 | F0408ALKD7 |
| 15-04-0164-1 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 04/08/15 21:35 | F0408ALKD7 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|---|---------------------|------------------|------------|---------------|-------------------|
| Alkalinity, Total (as CaCO ₃) | 349.0 | 350.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-04-0164-1 | Sample | Aqueous | PH1/BUR03 | N/A | 04/08/15 21:35 | F0408HCOD7 |
| 15-04-0164-1 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 04/08/15 21:35 | F0408HCOD7 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Bicarbonate (as CaCO ₃) | 349.0 | 350.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-04-0164-1 | Sample | Aqueous | PH1/BUR03 | N/A | 04/08/15 21:35 | F0408CO3D7 |
| 15-04-0164-1 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 04/08/15 21:35 | F0408CO3D7 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|---------------------|------------------|-----------------------|-----------------------|------------------------|
| HSM110 | Sample | Aqueous | SC 5 | 03/31/15 00:00 | 03/31/15 16:00 | F0331TDSD2 |
| HSM110 | Sample Duplicate | Aqueous | SC 5 | 03/31/15 00:00 | 03/31/15 16:00 | F0331TDSD2 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Solids, Total Dissolved | | 365.0 | 345.0 | 6 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-03-2045-1 | Sample | Aqueous | N/A | 03/30/15 00:00 | 03/30/15 14:00 | F0330TSSD1 |
| 15-03-2045-1 | Sample Duplicate | Aqueous | N/A | 03/30/15 00:00 | 03/30/15 14:00 | F0330TSSD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Suspended | 32.00 | 33.00 | 3 | 0-20 | |


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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|---------------------|------------------|---------------|-----------------------|------------------------|
| HSM110 | Sample | Aqueous | PH 1 | N/A | 03/26/15 18:08 | F0326PHD1 |
| HSM110 | Sample Duplicate | Aqueous | PH 1 | N/A | 03/26/15 18:08 | F0326PHD1 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| pH | | 7.290 | 7.360 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 4500 N Org B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-03-2116-5 | Sample | Aqueous | BUR05 | 04/03/15 00:00 | 04/03/15 17:40 | F0403TKND1 |
| 15-03-2116-5 | Sample Duplicate | Aqueous | BUR05 | 04/03/15 00:00 | 04/03/15 17:40 | F0403TKND1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Total Kjeldahl Nitrogen | 1.400 | 1.260 | 11 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-12-906-5550 | LCS | Aqueous | IC 15 | N/A | 03/26/15 20:00 | 150326L02 |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Fluoride | | 2.500 | 2.704 | 108 | 90-110 | |
| Chloride | | 50.00 | 49.16 | 98 | 90-110 | |
| Bromide | | 5.000 | 4.931 | 99 | 90-110 | |
| Nitrate (as N) | | 5.000 | 4.901 | 98 | 90-110 | |
| Sulfate | | 50.00 | 49.21 | 98 | 90-110 | |


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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-739-102 | LCS | Aqueous | ACA 1 | N/A | 04/09/15 14:45 | 150409LO1 | | | |
| 099-12-739-102 | LCSD | Aqueous | ACA 1 | N/A | 04/09/15 14:45 | 150409LO1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Phosphorus, Total | 0.2000 | 0.1942 | 97 | 0.1939 | 97 | 90-110 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-859-653 | LCS | Aqueous | PH1/BUR03 | N/A | 04/08/15 21:35 | F0408ALKB7 | | | |
| 099-15-859-653 | LCSD | Aqueous | PH1/BUR03 | N/A | 04/08/15 21:35 | F0408ALKB7 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 100.0 | 96.00 | 96 | 97.00 | 97 | 80-120 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4493 | LCS | Aqueous | SC 5 | 03/31/15 | 03/31/15 16:00 | F0331TDSL1 | | | |
| 099-12-180-4493 | LCSD | Aqueous | SC 5 | 03/31/15 | 03/31/15 16:00 | F0331TDSL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 100.0 | 100 | 105.0 | 105 | 80-120 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-09-010-7116 | LCS | Aqueous | N/A | 03/30/15 | 03/30/15 14:00 | F0330TSSL1 | | | |
| 099-09-010-7116 | LCSD | Aqueous | N/A | 03/30/15 | 03/30/15 14:00 | F0330TSSL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Suspended | 100.0 | 97.00 | 97 | 108.0 | 108 | 80-120 | 11 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-097-5582 | LCS | Aqueous | TOC 8 | 04/08/15 | 04/09/15 02:26 | F0408TOCL1 | | | |
| 099-05-097-5582 | LCSD | Aqueous | TOC 8 | 04/08/15 | 04/09/15 02:26 | F0408TOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 10.00 | 9.480 | 95 | 9.760 | 98 | 80-120 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-115-1415 | LCS | Aqueous | TOC 8 | 03/31/15 | 04/01/15 00:32 | F0331DOCL1 | | | |
| 099-05-115-1415 | LCSD | Aqueous | TOC 8 | 03/31/15 | 04/01/15 00:32 | F0331DOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Dissolved Organic | 10.00 | 9.390 | 94 | 9.630 | 96 | 80-120 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-15-683-1214 | LCS | Aqueous | ICP 7300 | 03/27/15 | 04/10/15 11:56 | 150327L4FF |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Calcium | | 0.5000 | 0.4970 | 99 | 80-120 | |
| Magnesium | | 0.5000 | 0.5718 | 114 | 80-120 | |
| Potassium | | 5.000 | 5.029 | 101 | 80-120 | |
| Sodium | | 5.000 | 5.056 | 101 | 80-120 | |
| Strontium | | 0.5000 | 0.4983 | 100 | 80-120 | |
| Silicon | | 0.5000 | 0.4713 | 94 | 80-120 | |


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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-771 | LCS | Aqueous | ICP/MS 03 | 03/27/15 | 03/30/15 12:35 | 150327LA2F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.09688 | 97 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.09482 | 95 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.09290 | 93 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.09953 | 100 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.09826 | 98 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.09945 | 99 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.09729 | 97 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.09527 | 95 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.09429 | 94 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.1013 | 101 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.04850 | 97 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.09357 | 94 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.09931 | 99 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.1154 | 115 | 80-120 | 73-127 | |
| Iron | 5.100 | 5.026 | 99 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.09703 | 97 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

Page 10 of 17

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|------------------|
| 099-15-763-527 | LCS | Aqueous | Mercury 04 | 04/01/15 | 04/01/15 16:38 | 150401L04F |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.01000 | 0.01105 | 110 | 85-121 | |


 Return to Contents



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|----------------|----------------|-----------------|-----------------------|------------------|
| 099-16-376-12 | LCS | Aqueous | LC/TQ 2 | 03/28/15 | 03/30/15 09:58 | 150328L13 |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Caffeine | 100.0 | 107.0 | 107 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-529-796 | LCS | Aqueous | | GC 44 | 03/27/15 | 03/30/15 12:46 | 150327L02A | | | |
| 099-12-529-796 | LCSD | Aqueous | | GC 44 | 03/27/15 | 03/30/15 13:00 | 150327L02A | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Alpha-BHC | 0.5000 | 0.4850 | 97 | 0.4828 | 97 | 50-135 | 36-149 | 0 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.5070 | 101 | 0.5042 | 101 | 50-135 | 36-149 | 1 | 0-25 | |
| Beta-BHC | 0.5000 | 0.4772 | 95 | 0.4782 | 96 | 50-135 | 36-149 | 0 | 0-25 | |
| Heptachlor | 0.5000 | 0.4821 | 96 | 0.4850 | 97 | 50-135 | 36-149 | 1 | 0-25 | |
| Delta-BHC | 0.5000 | 0.4634 | 93 | 0.4551 | 91 | 50-135 | 36-149 | 2 | 0-25 | |
| Aldrin | 0.5000 | 0.4749 | 95 | 0.4772 | 95 | 50-135 | 36-149 | 0 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.4811 | 96 | 0.4732 | 95 | 50-135 | 36-149 | 2 | 0-25 | |
| Endosulfan I | 0.5000 | 0.4744 | 95 | 0.4709 | 94 | 50-135 | 36-149 | 1 | 0-25 | |
| Dieldrin | 0.5000 | 0.4839 | 97 | 0.4751 | 95 | 50-135 | 36-149 | 2 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.4858 | 97 | 0.4703 | 94 | 50-135 | 36-149 | 3 | 0-25 | |
| Endrin | 0.5000 | 0.4426 | 89 | 0.4365 | 87 | 50-135 | 36-149 | 1 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.4914 | 98 | 0.4952 | 99 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.4758 | 95 | 0.4610 | 92 | 50-135 | 36-149 | 3 | 0-25 | |
| Endosulfan II | 0.5000 | 0.4640 | 93 | 0.4575 | 91 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.4236 | 85 | 0.4212 | 84 | 50-135 | 36-149 | 1 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.4443 | 89 | 0.4370 | 87 | 50-135 | 36-149 | 2 | 0-25 | |
| Methoxychlor | 0.5000 | 0.4198 | 84 | 0.4184 | 84 | 50-135 | 36-149 | 0 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-533-1019 | LCS | Aqueous | GC 31 | 03/27/15 | 03/30/15 12:35 | 150327L03 | | | |
| 099-12-533-1019 | LCSD | Aqueous | GC 31 | 03/27/15 | 03/30/15 12:54 | 150327L03 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Aroclor-1016 | 2.000 | 1.627 | 81 | 1.650 | 82 | 50-135 | 1 | 0-25 | |
| Aroclor-1260 | 2.000 | 1.787 | 89 | 1.774 | 89 | 50-135 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-963-91 | LCS | Aqueous | | GC 35 | 03/27/15 | 03/31/15 13:03 | 150327L04 | | | |
| 099-15-963-91 | LCSD | Aqueous | | GC 35 | 03/27/15 | 03/31/15 13:49 | 150327L04 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.03546 | 89 | 0.03499 | 87 | 30-130 | 13-147 | 1 | 0-30 | |
| Bolstar | 0.04000 | 0.03327 | 83 | 0.03426 | 86 | 30-130 | 13-147 | 3 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.03179 | 79 | 0.03357 | 84 | 30-130 | 13-147 | 5 | 0-30 | |
| Coumaphos | 0.04000 | 0.03464 | 87 | 0.03413 | 85 | 30-130 | 13-147 | 1 | 0-30 | |
| Diazinon | 0.04000 | 0.03167 | 79 | 0.03410 | 85 | 30-130 | 13-147 | 7 | 0-30 | |
| Disulfoton | 0.04000 | 0.03590 | 90 | 0.03761 | 94 | 30-130 | 13-147 | 5 | 0-30 | |
| Ethoprop | 0.04000 | 0.03292 | 82 | 0.03434 | 86 | 30-130 | 13-147 | 4 | 0-30 | |
| Fensulfothion | 0.04000 | 0.03543 | 89 | 0.03564 | 89 | 30-130 | 13-147 | 1 | 0-30 | |
| Fenthion | 0.04000 | 0.03284 | 82 | 0.03428 | 86 | 30-130 | 13-147 | 4 | 0-30 | |
| Merphos | 0.04000 | 0.03377 | 84 | 0.03489 | 87 | 30-130 | 13-147 | 3 | 0-30 | |
| Methyl Parathion | 0.04000 | 0.03267 | 82 | 0.03369 | 84 | 30-130 | 13-147 | 3 | 0-30 | |
| Phorate | 0.04000 | 0.03866 | 97 | 0.04052 | 101 | 30-130 | 13-147 | 5 | 0-30 | |
| Ronnel | 0.04000 | 0.03059 | 76 | 0.03228 | 81 | 30-130 | 13-147 | 5 | 0-30 | |
| Stirophos | 0.04000 | 0.02142 | 54 | 0.02241 | 56 | 30-130 | 13-147 | 5 | 0-30 | |
| Tokuthion | 0.04000 | 0.03195 | 80 | 0.03343 | 84 | 30-130 | 13-147 | 5 | 0-30 | |
| Trichloronate | 0.04000 | 0.03261 | 82 | 0.03435 | 86 | 30-130 | 13-147 | 5 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 095-01-034-640 | LCS | Aqueous | GC 40 | 03/27/15 | 03/31/15 16:50 | 150327L05 |
| 095-01-034-640 | LCSD | Aqueous | GC 40 | 03/27/15 | 03/31/15 17:13 | 150327L05 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| 2,4-D | 20.00 | 15.04 | 75 | 18.24 | 91 | 30-130 | 19 | 0-30 | |
| 2,4,5-T | 2.000 | 2.065 | 103 | 1.985 | 99 | 30-130 | 4 | 0-30 | |
| 2,4-DB | 20.00 | 15.68 | 78 | 19.28 | 96 | 30-130 | 21 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-4016 | LCS | Aqueous | | GC/MS CCC | 03/27/15 | 03/27/15 20:42 | 150327L06 | | | |
| 095-01-003-4016 | LCSD | Aqueous | | GC/MS CCC | 03/27/15 | 03/27/15 21:00 | 150327L06 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 200.0 | 173.6 | 87 | 175.8 | 88 | 61-120 | 51-130 | 1 | 0-20 | |
| Acenaphthylene | 200.0 | 173.1 | 87 | 173.2 | 87 | 55-120 | 44-131 | 0 | 0-20 | |
| Butyl Benzyl Phthalate | 200.0 | 182.5 | 91 | 180.6 | 90 | 56-122 | 45-133 | 1 | 0-20 | |
| 4-Chloro-3-Methylphenol | 200.0 | 147.8 | 74 | 143.8 | 72 | 52-120 | 41-131 | 3 | 0-20 | |
| 2-Chlorophenol | 200.0 | 151.7 | 76 | 149.2 | 75 | 47-120 | 35-132 | 2 | 0-20 | |
| 1,4-Dichlorobenzene | 200.0 | 148.7 | 74 | 149.7 | 75 | 36-120 | 22-134 | 1 | 0-20 | |
| Dimethyl Phthalate | 200.0 | 161.9 | 81 | 163.6 | 82 | 60-120 | 50-130 | 1 | 0-20 | |
| 2,4-Dinitrotoluene | 200.0 | 158.7 | 79 | 157.3 | 79 | 61-121 | 51-131 | 1 | 0-20 | |
| Fluorene | 200.0 | 173.1 | 87 | 172.3 | 86 | 67-120 | 58-129 | 0 | 0-20 | |
| N-Nitroso-di-n-propylamine | 200.0 | 147.4 | 74 | 146.9 | 73 | 39-123 | 25-137 | 0 | 0-20 | |
| Naphthalene | 200.0 | 154.2 | 77 | 152.0 | 76 | 54-120 | 43-131 | 1 | 0-20 | |
| 4-Nitrophenol | 200.0 | 154.6 | 77 | 151.2 | 76 | 14-120 | 0-138 | 2 | 0-20 | |
| Pentachlorophenol | 200.0 | 131.3 | 66 | 127.7 | 64 | 31-127 | 15-143 | 3 | 0-20 | |
| Phenol | 200.0 | 148.1 | 74 | 146.9 | 73 | 17-120 | 0-137 | 1 | 0-20 | |
| Pyrene | 200.0 | 169.4 | 85 | 167.6 | 84 | 58-124 | 47-135 | 1 | 0-20 | |
| 1,2,4-Trichlorobenzene | 200.0 | 152.4 | 76 | 152.9 | 76 | 49-120 | 37-132 | 0 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 03/26/15
Work Order: 15-03-2020
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-16776 | LCS | Aqueous | GC/MS QQ | 04/04/15 | 04/04/15 10:17 | 150404L002 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 47.60 | 95 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 63.50 | 127 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 47.52 | 95 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 48.22 | 96 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 46.49 | 93 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 47.55 | 95 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 47.83 | 96 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 50.50 | 101 | 80-123 | 73-130 | |
| Toluene | 50.00 | 48.57 | 97 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 47.31 | 95 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 48.27 | 97 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 97.32 | 97 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 48.81 | 98 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 51.88 | 104 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Sample Analysis Summary Report

Work Order: 15-03-2020

Page 1 of 1

| <u>Method</u> | <u>Extraction</u> | <u>Chemist ID</u> | <u>Instrument</u> | <u>Analytical Location</u> |
|-----------------------|-------------------|-------------------|-------------------|----------------------------|
| EPA 1694 (M) Caffeine | EPA 1694 | 262 | LC/TQ 2 | 1 |
| EPA 300.0 | N/A | 650 | IC 15 | 1 |
| EPA 365.1 | N/A | 735 | ACA 1 | 1 |
| EPA 6010B | EPA 3005A Filt. | 935 | ICP 7300 | 1 |
| EPA 6020 | EPA 3005A Filt. | 598 | ICP/MS 03 | 1 |
| EPA 7470A | EPA 7470A Filt. | 915 | Mercury 04 | 1 |
| EPA 8081A | EPA 3510C | 669 | GC 44 | 1 |
| EPA 8082 | EPA 3510C | 960 | GC 31 | 1 |
| EPA 8141A | EPA 3510C | 949 | GC 35 | 1 |
| EPA 8151A | EPA 8151A | 669 | GC 40 | 1 |
| EPA 8260B | EPA 5030C | 486 | GC/MS QQ | 2 |
| EPA 8270C | EPA 3510C | 923 | GC/MS CCC | 1 |
| SM 2320B | N/A | 885 | PH1/BUR03 | 1 |
| SM 2540 C | N/A | 689 | SC 5 | 1 |
| SM 2540 D | N/A | 689 | N/A | 1 |
| SM 4500 H+ B | N/A | 688 | PH 1 | 1 |
| SM 4500 N Org B | N/A | 685 | BUR05 | 1 |
| SM 5310 B | N/A | 735 | TOC 8 | 1 |

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Location 1: 7440 Lincoln Way, Garden Grove, CA 92841

Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841

Glossary of Terms and Qualifiers

Work Order: 15-03-2020

Page 1 of 1

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|--|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDS or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.



Calscience

CHAIN OF CUSTODY RECORD

DATE: March 25th, 2015
PAGE: 1 OF 1

WO # / LAB USE ONLY

15-03-2020

7440 Lincoln Way, Garden Grove, CA 92641-1427 • (714) 895-5494
For courier service / sample drop off information, contact us26_sales@eurofins.com or call us.

LABORATORY CLIENT:

SWCA Environmental Consultants

ADDRESS: 6200 UTSA Blvd. Suite 102

CITY: San Antonio

STATE: TX

ZIP: 78249-1618

TEL: 210.877.2847

E-MAIL: P Pearce@swca.com

CLIENT PROJECT NAME / NUMBER:

EAA 27122

PROJECT CONTACT:

Philip Pearce

P.O. NO.:

SAMPLER(S) (PRINT)

Jennifer Moreland
Brittany Reiss

REQUESTED ANALYSES

TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD"):
☐ SAME DAY ☐ 24 HR ☐ 48 HR ☐ 72 HR ☒ 140 Days (standard)
☐ COELT EDF
GLOBAL ID: _____ LOG CODE: _____

SPECIAL INSTRUCTIONS:

Samples for EPA 6010B, EPA 6020, and EPA 7470 Metals are field filtered
EPA 6020 Metals: Zn, Ti, Se, Sb, Pb, Ni, Cu, Cr, Cd, Be, Ba, As, Al, Ag, Fe, Mn
Please also analyze each sample for pH.

| Please check box or fill in blank as needed. | | | | | | | | | | | | | | | | | | | | | | | | | |
|--|-----------|--------------|----------|-------|--------|----------------|-----------|-------------|-----------------------------------|-------------------|-----------------------|-------------------------------------|-----------------------|---------------------------------|----------------------------------|----------------------------------|-----------------------------|---|----------------------------|--|--------------|--------------|----------------------------------|---------------------|---------------|
| LAB USE ONLY | SAMPLE ID | NO. OF CONT. | SAMPLING | | MATRIX | Field Filtered | Preserved | Unpreserved | EPA 6010B (Si, Ca, Mg, K, Na, Sr) | EPA 7470A Mercury | EPA 1694 (M) Caffeine | EPA 8081A Organochlorine Pesticides | EPA 8082 PCB Aroclors | EPA 8141A Organophosphorus Pest | EPA 8151A Chlorinated Herbicides | EPA 8270C Semi-Volatile Organics | EPA 8260B Volatile Organics | EPA 300 Anions (F, Cl, Br, NO3, SO4, CO3) | EPA 365.1 Total Phosphorus | SM 2320B Alkalinity: Total, Bicarb, Carb | SM 5310B TOC | SM 5310B DOC | SM 5310B Solids, Total Dissolved | SM 4500 N Org B TKN | EPA 160.2 TSS |
| | | | DATE | TIME | | | | | | | | | | | | | | | | | | | | | |
| 1 | HSM 110 | 9 | 3/25/15 | 10:28 | NPW | 1 | 7 | 2 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 2 | HSM 120 | 9 | 3/25/15 | 10:51 | NPW | 1 | 7 | 2 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 3 | HSM 130 | 9 | 3/25/15 | 11:17 | NPW | 1 | 7 | 2 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 4 | HSM 140 | 9 | 3/25/15 | 11:30 | NPW | 1 | 7 | 2 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 5 | HSM 150 | 9 | 3/25/15 | 12:25 | NPW | 1 | 7 | 2 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 6 | HSM 160 | 9 | 3/25/15 | 12:50 | NPW | 1 | 7 | 2 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 7 | HSM 170 | 9 | 3/25/15 | 13:18 | NPW | 1 | 7 | 2 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 8 | FDHSM 170 | 9 | 3/25/15 | 10:28 | NPW | 1 | 7 | 2 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 9 | TBQ4 | 1 | 3/25/15 | NA | NPW | | | | | | | | | | | | | | | | | | | | |

Relinquished by: (Signature) [Signature] Received by: (Signature/Affiliation)
Relinquished by: (Signature) [Signature] Received by: (Signature/Affiliation)
Relinquished by: (Signature) [Signature] Received by: (Signature/Affiliation)

Date: 3/25/15 Time: 16:00

Date: _____ Time: _____

Date: 3/26/15 Time: 1040

Michael Belvin
SWCA
6200 UTSA Blvd
Sta 102
San Antonio, TX 78249



ActWgt: 60.0 LB
CAD: 8347991/INET3610

Dims: 24 X 13 X 14 IN

SHIP TO: (714) 895-5494
Lab
Eurofins Calscience
7440 LINCOLN WAY
GARDEN GROVE, CA 92841

BILL SENDER

Delivery Address Bar Code



Ref # 27122.02.06
Invoice #
PO #
Dept #

1 of 8

THU - 26 MAR 3:00P
STANDARD OVERNIGHT

TRK# 7732 1480 6187

0201

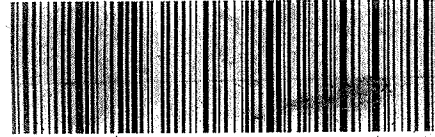
MASTER

A7 APVA

92841

CA-US

SNA



2020

Ship Date: 25MAR15
ActWgt: 60.0 LB
CAD: 8347991/INET3610

Dims: 24 X 13 X 14 IN

Delivery Address Bar Code



Ref # 27122.02.06
Invoice #
PO #
Dept #

Ship Date: 25MAR15
ActWgt: 60.0 LB
CAD: 8347991/INET3610

Dims: 24 X 13 X 14 IN

Delivery Address Bar Code



Ref # 27122.02.06
Invoice #
PO #
Dept #

Ship Date: 25MAR15
ActWgt: 60.0 LB
CAD: 8347991/INET3610

Dims: 24 X 13 X 14 IN

Delivery Address Bar Code



Ref # 27122.02.06
Invoice #
PO #
Dept #

2 of 8

THU - 26 MAR 3:00P
STANDARD OVERNIGHT

MPS# 7732 1480 6382

0263

Mstr# 7732 1480 6187

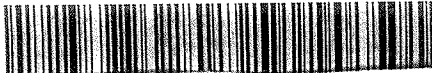
0201

A7 APVA

92841

CA-US

SNA



3 of 8

THU - 26 MAR 3:00P
STANDARD OVERNIGHT

MPS# 7732 1480 6268

0263

Mstr# 7732 1480 6187

0201

A7 APVA

92841

CA-US

SNA



4 of 8

THU - 26 MAR 3:00P
STANDARD OVERNIGHT

MPS# 7732 1480 6636

0263

Mstr# 7732 1480 6187

0201

A7 APVA

92841

CA-US

SNA



Ship Date: 25MAR15
ActWgt: 60.0 LB
CAD: 8347991/INET3610

Dims: 24 X 13 X 14 IN

Delivery Address Bar Code



Ref # 27122.02.06
Invoice #
PO #
Dept #

Ship Date: 25MAR15
ActWgt: 60.0 LB
CAD: 8347991/INET3610

Dims: 24 X 13 X 14 IN

Delivery Address Bar Code



Ref # 27122.02.06
Invoice #
PO #
Dept #

Ship Date: 25MAR15
ActWgt: 60.0 LB
CAD: 8347991/INET3610

Dims: 24 X 13 X 14 IN

Delivery Address Bar Code



Ref # 27122.02.06
Invoice #
PO #
Dept #

5 of 8

THU - 26 MAR 3:00P
STANDARD OVERNIGHT

MPS# 7732 1480 6740

0263

Mstr# 7732 1480 6187

0201

A7 APVA

92841

CA-US

SNA



6 of 8

THU - 26 MAR 3:00P
STANDARD OVERNIGHT

MPS# 7732 1480 7003

0263

Mstr# 7732 1480 6187

0201

A7 APVA

92841

CA-US

SNA



7 of 8

THU - 26 MAR 3:00P
STANDARD OVERNIGHT

MPS# 7732 1480 7036

0263

Mstr# 7732 1480 6187

0201

A7 APVA

92841

CA-US

SNA



Ship Date: 25MAR15
ActWgt: 60.0 LB
CAD: 8347991/INET3610

Dims: 24 X 13 X 14 IN

Delivery Address Bar Code



Ref # 27122.02.06

THU - 26 MAR 3:00P
STANDARD OVERNIGHT

8 of 8

7732 1480 7140

Mstr# 7732 1480 6187

0201

A7 APVA

92841

CA-US

SNA



537118/3AEE48



Calscience

WORK ORDER NUMBER: 15-03-2020

SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 8CLIENT: SWCADATE: 03/26/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 2.7 °C (w/ CF): 2.9 °C ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ FilterChecked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Not Intact ☐ Not Present ☐ N/AChecked by: 15Sample(s) ☐ Present and Intact ☐ Not Intact ☒ Not Present ☐ N/AChecked by: 689

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|--------------------------|--------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/ACONTAINER TYPE: 3(Trip Blank Lot Number: 150306A)Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn₁ ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal cube ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ 16ozPJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____ ☐ _____Air: ☐ Tedlar® ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ _____ ☐ _____

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,s=H₂SO₄, u=ultra-pure, z₂na=Zn(CH₃CO₂)₂ + NaOHLabeled/Checked by: 689Reviewed by: 816



Calscience

WORK ORDER NUMBER: 15-03-2020

SAMPLE RECEIPT CHECKLIST

COOLER 2 OF 8

CLIENT: SWCA

DATE: 03/26/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 3.3 °C (w/ CF): 3.5 °C ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Not Intact☐ Not Present☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact☐ Not Intact☒ Not Present☐ N/A

Checked by: 659

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: 3

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznnna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBna ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal cube ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ 16ozPJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____ ☐ _____Air: ☐ Tedlar® ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ _____ ☐ _____

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,s=H₂SO₄, u=ultra-pure, znna=Zn(CH₃CO₂)₂ + NaOH

Labeled/Checked by: 659

Reviewed by: 876



Calscience

WORK ORDER NUMBER: 15-03-2020

SAMPLE RECEIPT CHECKLIST

COOLER 3 OF 8

CLIENT: SWCA

DATE: 03/26/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 3.4 °C (w/ CF): 3.6 °C ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Not Intact☐ Not Present☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact☐ Not Intact☒ Not Present☐ N/A

Checked by: 659

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|--------------------------|--------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC..... ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC..... ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition..... ☒ Yes ☐ No ☐ N/AProper containers for analyses requested..... ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested..... ☒ Yes ☐ No ☐ N/ASamples received within holding time..... ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace..... ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation..... ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: 3

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBnf ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal cube ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ 16ozPJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____ ☐ _____Air: ☐ Tedlar® ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ _____ ☐ _____

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,s=H₂SO₄, u=ultra-pure, z₂na=Zn(CH₃CO₂)₂ + NaOH

Labeled/Checked by: 659

Reviewed by: 876



Calscience

WORK ORDER NUMBER: 15-03-2020

SAMPLE RECEIPT CHECKLIST

COOLER 4 OF 8

CLIENT: SWCA

DATE: 03/26/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 4.0 °C (w/ CF): 4.2 °C ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Not Intact☐ Not Present☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact☐ Not Intact☒ Not Present☐ N/A

Checked by: 659

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete..... ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC..... ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC..... ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition..... ☒ Yes ☐ No ☐ N/AProper containers for analyses requested..... ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested..... ☒ Yes ☐ No ☐ N/ASamples received within holding time..... ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace..... ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation..... ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: ? (Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBna ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal cube ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ 16ozPJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____ ☐ _____Air: ☐ Tedlar® ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ _____ ☐ _____

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,s=H₂SO₄, u=ultra-pure, z₂na=Zn(CH₃CO₂)₂ + NaOH

Labeled/Checked by: 659

Reviewed by: 806



Calscience

WORK ORDER NUMBER: 15-03-2020

SAMPLE RECEIPT CHECKLIST

COOLER 5 OF 8

CLIENT: SWCA

DATE: 03 / 26 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 3.5 °C (w/ CF): 3.7 °C ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Not Intact☐ Not Present☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact☐ Not Intact☒ Not Present☐ N/A

Checked by: 659

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: 3 (Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na2} ☐ 100PJ ☐ 100PJ_{na2} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☒ 250PB_{na} ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s☐ 500PB ☒ 1AGB ☐ 1AGB_{na2} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5 gal cube ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ 16ozPJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____ ☐ _____Air: ☐ Tedlar® ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ _____ ☐ _____

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,s=H₂SO₄, u=ultra-pure, znna=Zn(CH₃CO₂)₂ + NaOH

Labeled/Checked by: 659

Reviewed by: 896



Calscience

WORK ORDER NUMBER: 15-03-2020

SAMPLE RECEIPT CHECKLIST

COOLER 6 OF 8CLIENT: SWCADATE: 03 / 26 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 4.1 °C (w/ CF): 4.3 °C ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ FilterChecked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Not Intact☐ Not Present☐ N/AChecked by: 15Sample(s) ☐ Present and Intact☐ Not Intact☒ Not Present☐ N/AChecked by: 659

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/ACONTAINER TYPE: 3

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznnna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBna ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal cube ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ 16ozPJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____ ☐ _____Air: ☐ Tedlar® ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ _____ ☐ _____

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,s=H₂SO₄, u=ultra-pure, znnna=Zn(CH₃CO₂)₂ + NaOHLabeled/Checked by: 659Reviewed by: 8m



Calscience

WORK ORDER NUMBER: 15-03-2020

SAMPLE RECEIPT CHECKLIST

COOLER 7 OF 8CLIENT: SWCADATE: 03 / 26 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 3.9 °C (w/ CF): 4.1 °C ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ FilterChecked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Not Intact☐ Not Present☐ N/AChecked by: 15Sample(s) ☐ Present and Intact☐ Not Intact☒ Not Present☐ N/AChecked by: 659

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete..... ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC..... ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC..... ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition..... ☒ Yes ☐ No ☐ N/AProper containers for analyses requested..... ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested..... ☒ Yes ☐ No ☐ N/ASamples received within holding time..... ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace..... ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation..... ☐ Yes ☐ No ☒ N/ACONTAINER TYPE: 2

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznnna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBnf ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal cube ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ 16ozPJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____ ☐ _____Air: ☐ Tedlar® ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ _____ ☐ _____

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,s=H₂SO₄, u=ultra-pure, znnna=Zn(CH₃CO₂)₂ + NaOHLabeled/Checked by: 659Reviewed by: 896

SAMPLE RECEIPT CHECKLIST

COOLER 8 OF 8

CLIENT: SWCA

DATE: 03/26/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC4 (CF: +0.2°C) Temperature (w/o CF): 3.5 °C (w/ CF): 3.7 °C ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Not Intact☐ Not Present☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact☐ Not Intact☒ Not Present☐ N/A

Checked by: 659

SAMPLE CONDITION:

Yes No N/A

Chain-of-Custody (COC) document(s) received with samples ☒ ☐ ☐COC document(s) received complete..... ☒ ☐ ☐☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC..... ☒ ☐ ☐Sample container label(s) consistent with COC..... ☒ ☐ ☐Sample container(s) intact and in good condition..... ☒ ☐ ☐Proper containers for analyses requested..... ☒ ☐ ☐Sufficient volume/mass for analyses requested..... ☒ ☐ ☐Samples received within holding time..... ☒ ☐ ☐

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ ☒ ☐Proper preservation chemical(s) noted on COC and/or sample container ☒ ☐ ☐

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace..... ☒ ☐ ☐☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation..... ☐ ☐ ☒

CONTAINER TYPE: 2

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznn_a ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn_f ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal cube ☐ ☐ ☐Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ 16ozPJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ ☐ ☐Air: ☐ Tedlar® ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ ☐ ☐

Container: A=Amber, B=Bottle, C=Clear, E=Envelope, G=Glass, J=Jar, P=Plastic, and Z= Ziploc/Resealable Bag

Preservative: f=filtered, h=HCl, n=HNO₃, na=NaOH, na₂=Na₂S₂O₃, p=H₃PO₄,s=H₂SO₄, u=ultra-pure, znn_a=Zn(CH₃CO₂)₂ + NaOH

Labeled/Checked by: 659

Reviewed by: 659

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 11, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HCS110
Sample Number: AA96986
Sample Matrix: Non Potable Water

Collection Date/Time: 9/9/15 10:11
Receipt Date/Time: 9/9/15 14:07

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA96986-A | E. coli | √ | SM 9223B-2004 | 800 | MPN/100 mL | | 1 | 44141 | 9/9/15 | 16:11 | MD/PAL |
| AA96986-A | E. Coli Holding Time - IDEXX Colilert | | NA | 6.00 | hours | | 0.00 | 44140 | 9/9/15 | 16:11 | MD/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44141

QC Analyte Name

Initial Blank for E. coli

Log Range for E. coli

Result

Absent

0.4746

Units

Qualifier

Lower

0.0

Target

Absent

Upper

0.5



Patricia M. Carvajal
Quality Assurance Supervisor

9/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 11, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS120
Sample Number: AA96987
Sample Matrix: Non Potable Water

Collection Date/Time: 9/9/15 10:44
Receipt Date/Time: 9/9/15 14:07

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA96987-A | E. coli | √ | SM 9223B-2004 | 67 | MPN/100 mL | | 1 | 44141 | 9/9/15 | 16:11 | MD/PAL |
| AA96987-A | E. Coli Holding Time - IDEXX Colilert | | NA | 5.45 | hours | | 0.00 | 44140 | 9/9/15 | 16:11 | MD/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44141

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

9/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 11, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: FDHCS120
Sample Number: AA96988
Sample Matrix: Non Potable Water

Collection Date/Time: 9/9/15 10:44
Receipt Date/Time: 9/9/15 14:07

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA96988-A | E. coli | √ | SM 9223B-2004 | 54 | MPN/100 mL | | 1 | 44141 | 9/9/15 | 16:11 | MD/PAL |
| AA96988-A | E. Coli Holding Time - IDEXX Colilert | | NA | 5.45 | hours | | 0.00 | 44140 | 9/9/15 | 16:11 | MD/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44141

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

9/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 11, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS130
Sample Number: AA96989
Sample Matrix: Non Potable Water

Collection Date/Time: 9/9/15 09:30
Receipt Date/Time: 9/9/15 14:07

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA96989-A | E. coli | √ | SM 9223B-2004 | 460 | MPN/100 mL | | 1 | 44141 | 9/9/15 | 16:11 | MD/PAL |
| AA96989-A | E. Coli Holding Time - IDEXX Colilert | | NA | 6.68 | hours | | 0.00 | 44140 | 9/9/15 | 16:11 | MD/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44141

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

9/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 11, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS140
Sample Number: AA96990
Sample Matrix: Non Potable Water

Collection Date/Time: 9/9/15 11:12
Receipt Date/Time: 9/9/15 14:07

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA96990-A | E. coli | √ | SM 9223B-2004 | 140 | MPN/100 mL | | 1 | 44141 | 9/9/15 | 16:11 | MD/PAL |
| AA96990-A | E. Coli Holding Time - IDEXX Colilert | | NA | 4.98 | hours | | 0.00 | 44140 | 9/9/15 | 16:11 | MD/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44141

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

9/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



September 11, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS160
Sample Number: AA96991
Sample Matrix: Non Potable Water

Collection Date/Time: 9/9/15 11:39
Receipt Date/Time: 9/9/15 14:07

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA96991-A | E. coli | √ | SM 9223B-2004 | 120 | MPN/100 mL | | 1 | 44141 | 9/9/15 | 16:11 | MD/PAL |
| AA96991-A | E. Coli Holding Time - IDEXX Colilert | | NA | 4.53 | hours | | 0.00 | 44140 | 9/9/15 | 16:11 | MD/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44141

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

9/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable



Calscience



WORK ORDER NUMBER: 15-09-0733

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: SWCA Environmental Consultants

Client Project Name: EAA 27122

Attention: Philip Pearce
6200 UTSA Blvd.
Suite 102
San Antonio, TX 78249-1618

Approved for release on 10/04/2015 by:
Don Burley
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

Contents

Client Project Name: EAA 27122
Work Order Number: 15-09-0733

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| 1 | Work Order Narrative. | 3 |
| 2 | Sample Summary. | 4 |
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Work Order Narrative

Work Order: 15-09-0733

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Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 09/10/15. They were assigned to Work Order 15-09-0733.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.



Calscience

Sample Summary

| | | | |
|---------|--------------------------------|-----------------------|----------------|
| Client: | SWCA Environmental Consultants | Work Order: | 15-09-0733 |
| | 6200 UTSA Blvd., Suite 102 | Project Name: | EAA 27122 |
| | San Antonio, TX 78249-1618 | PO Number: | 27122.02.01 |
| | | Date/Time Received: | 09/10/15 11:00 |
| | | Number of Containers: | 55 |

Attn: Philip Pearce

| Sample Identification | Lab Number | Collection Date and Time | Number of Containers | Matrix |
|-----------------------|--------------|--------------------------|----------------------|---------|
| HCS110 | 15-09-0733-1 | 09/09/15 10:11 | 9 | Aqueous |
| HCS120 | 15-09-0733-2 | 09/09/15 10:44 | 9 | Aqueous |
| FDHCS120 | 15-09-0733-3 | 09/09/15 10:44 | 9 | Aqueous |
| HCS130 | 15-09-0733-4 | 09/09/15 09:30 | 9 | Aqueous |
| HCS140 | 15-09-0733-5 | 09/09/15 11:12 | 9 | Aqueous |
| HCS160 | 15-09-0733-6 | 09/09/15 11:39 | 9 | Aqueous |
| TB11 | 15-09-0733-7 | 09/09/15 00:00 | 1 | Aqueous |

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Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

Page 1 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS110 | 15-09-0733-1-I | 09/09/15 10:11 | Aqueous | IC 15 | N/A | 09/10/15 18:37 | 150910L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.23 | 0.10 | 0.025 | 1.00 | |
| Chloride | 18 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.098 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.2 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 26 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|-------|-----|----------------|-----------|
| HCS120 | 15-09-0733-2-I | 09/09/15 10:44 | Aqueous | IC 15 | N/A | 09/10/15 18:55 | 150910L01 |
|--------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.23 | 0.10 | 0.025 | 1.00 | |
| Chloride | 17 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.10 | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | 1.8 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 25 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|----------|----------------|----------------|---------|-------|-----|----------------|-----------|
| FDHCS120 | 15-09-0733-3-I | 09/09/15 10:44 | Aqueous | IC 15 | N/A | 09/10/15 19:14 | 150910L01 |
|----------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.22 | 0.10 | 0.025 | 1.00 | |
| Chloride | 17 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.096 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.8 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 25 | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

Page 2 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS130 | 15-09-0733-4-I | 09/09/15 09:30 | Aqueous | IC 15 | N/A | 09/10/15 19:32 | 150910L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.22 | 0.10 | 0.025 | 1.00 | |
| Chloride | 18 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.097 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.8 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 32 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|-------|-----|----------------|-----------|
| HCS140 | 15-09-0733-5-I | 09/09/15 11:12 | Aqueous | IC 15 | N/A | 09/10/15 19:51 | 150910L01 |
|--------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.21 | 0.10 | 0.025 | 1.00 | |
| Chloride | 17 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.11 | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | 1.7 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 26 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|-------|-----|----------------|-----------|
| HCS160 | 15-09-0733-6-I | 09/09/15 11:39 | Aqueous | IC 15 | N/A | 09/10/15 20:09 | 150910L01 |
|--------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.22 | 0.10 | 0.025 | 1.00 | |
| Chloride | 18 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.10 | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | 1.8 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 29 | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-906-6058 | N/A | Aqueous | IC 15 | N/A | 09/10/15 12:34 | 150910L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | ND | 0.10 | 0.025 | 1.00 | |
| Chloride | ND | 1.0 | 0.12 | 1.00 | |
| Bromide | ND | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | ND | 0.10 | 0.025 | 1.00 | |
| Sulfate | ND | 1.0 | 0.19 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS110 | 15-09-0733-1-D | 09/09/15 10:11 | Aqueous | ICP 7300 | 09/11/15 | 09/19/15 23:22 | 150911LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 77.9 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 17.5 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.75 | 0.500 | 0.103 | 1.00 | B |
| Sodium | 12.5 | 0.500 | 0.103 | 1.00 | B |
| Strontium | 0.729 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.32 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS120 | 15-09-0733-2-D | 09/09/15 10:44 | Aqueous | ICP 7300 | 09/11/15 | 09/19/15 23:32 | 150911LA5F |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 82.2 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 16.3 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 2.30 | 0.500 | 0.103 | 1.00 | B |
| Sodium | 12.7 | 0.500 | 0.103 | 1.00 | B |
| Strontium | 0.670 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.55 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|----------|----------------|----------------|---------|----------|----------|----------------|------------|
| FDHCS120 | 15-09-0733-3-D | 09/09/15 10:44 | Aqueous | ICP 7300 | 09/11/15 | 09/19/15 23:36 | 150911LA5F |
|----------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 81.3 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 15.4 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 2.04 | 0.500 | 0.103 | 1.00 | B |
| Sodium | 12.2 | 0.500 | 0.103 | 1.00 | B |
| Strontium | 0.665 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.20 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

Page 2 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS130 | 15-09-0733-4-D | 09/09/15 09:30 | Aqueous | ICP 7300 | 09/11/15 | 09/19/15 23:39 | 150911LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 82.7 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 16.0 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.81 | 0.500 | 0.103 | 1.00 | B |
| Sodium | 13.5 | 0.500 | 0.103 | 1.00 | B |
| Strontium | 0.666 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.24 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS140 | 15-09-0733-5-D | 09/09/15 11:12 | Aqueous | ICP 7300 | 09/11/15 | 09/19/15 23:42 | 150911LA5F |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 84.3 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 16.0 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.77 | 0.500 | 0.103 | 1.00 | B |
| Sodium | 12.6 | 0.500 | 0.103 | 1.00 | B |
| Strontium | 0.665 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.39 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS160 | 15-09-0733-6-D | 09/09/15 11:39 | Aqueous | ICP 7300 | 09/11/15 | 09/19/15 23:46 | 150911LA5F |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 81.2 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 15.5 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.73 | 0.500 | 0.103 | 1.00 | B |
| Sodium | 12.5 | 0.500 | 0.103 | 1.00 | B |
| Strontium | 0.655 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.09 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-683-1396 | N/A | Aqueous | ICP 7300 | 09/11/15 | 09/19/15 22:59 | 150911LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 0.0267 | 0.100 | 0.0118 | 1.00 | J |
| Magnesium | 0.0439 | 0.100 | 0.00336 | 1.00 | J |
| Potassium | 0.117 | 0.500 | 0.103 | 1.00 | J |
| Sodium | 0.116 | 0.500 | 0.103 | 1.00 | J |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 1 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS110 | 15-09-0733-1-D | 09/09/15 10:11 | Aqueous | ICP/MS 03 | 09/16/15 | 09/22/15 01:44 | 150916LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | 0.000505 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | 0.0582 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000492 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00457 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00199 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000503 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0314 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00923 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0424 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00481 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 2 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS120 | 15-09-0733-2-D | 09/09/15 10:44 | Aqueous | ICP/MS 03 | 09/16/15 | 09/22/15 01:48 | 150916LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | 0.000388 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | 0.0549 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000524 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00504 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00182 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000521 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | 0.000121 | 0.00100 | 0.000101 | 1.00 | J |
| Zinc | 0.00984 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00699 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0364 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.000820 | 0.00100 | 0.000139 | 1.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 3 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS120 | 15-09-0733-3-D | 09/09/15 10:44 | Aqueous | ICP/MS 03 | 09/16/15 | 09/22/15 01:51 | 150916LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0537 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000482 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.000780 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00187 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000932 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00773 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00954 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0406 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.000640 | 0.00100 | 0.000139 | 1.00 | J |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS130 | 15-09-0733-4-D | 09/09/15 09:30 | Aqueous | ICP/MS 03 | 09/16/15 | 09/22/15 01:55 | 150916LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0553 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000555 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00160 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00204 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000744 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | 0.000109 | 0.00100 | 0.000101 | 1.00 | J |
| Zinc | 0.0159 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00749 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0465 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00114 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS140 | 15-09-0733-5-D | 09/09/15 11:12 | Aqueous | ICP/MS 03 | 09/16/15 | 09/22/15 01:58 | 150916LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000106 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0555 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000583 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00276 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000112 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00228 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.00229 | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0350 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00923 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0451 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00137 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS160 | 15-09-0733-6-D | 09/09/15 11:39 | Aqueous | ICP/MS 03 | 09/16/15 | 09/22/15 02:02 | 150916LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0550 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000645 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00660 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00235 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.00193 | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00714 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00801 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0399 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00269 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-912 | N/A | Aqueous | ICP/MS 03 | 09/16/15 | 09/21/15 22:16 | 150916LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS110 | 15-09-0733-1-D | 09/09/15 10:11 | Aqueous | Mercury 04 | 09/15/15 | 09/15/15 20:36 | 150915LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS120 | 15-09-0733-2-D | 09/09/15 10:44 | Aqueous | Mercury 04 | 09/15/15 | 09/15/15 20:38 | 150915LA3F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|----------|----------------|----------------|---------|------------|----------|----------------|------------|
| FDHCS120 | 15-09-0733-3-D | 09/09/15 10:44 | Aqueous | Mercury 04 | 09/15/15 | 09/15/15 20:40 | 150915LA3F |
|----------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS130 | 15-09-0733-4-D | 09/09/15 09:30 | Aqueous | Mercury 04 | 09/15/15 | 09/15/15 20:47 | 150915LA3F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS140 | 15-09-0733-5-D | 09/09/15 11:12 | Aqueous | Mercury 04 | 09/15/15 | 09/15/15 20:49 | 150915LA3F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS160 | 15-09-0733-6-D | 09/09/15 11:39 | Aqueous | Mercury 04 | 09/15/15 | 09/15/15 20:51 | 150915LA3F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-763-614 | N/A | Aqueous | Mercury 04 | 09/15/15 | 09/15/15 20:09 | 150915LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS110 | 15-09-0733-1-I | 09/09/15 10:11 | Aqueous | GC 44 | 09/11/15 | 09/22/15 07:07 | 150911L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 75 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 73 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS120 | 15-09-0733-2-I | 09/09/15 10:44 | Aqueous | GC 44 | 09/11/15 | 09/22/15 07:21 | 150911L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 95 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 92 | 50-135 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| FDHCS120 | 15-09-0733-3-I | 09/09/15 10:44 | Aqueous | GC 44 | 09/11/15 | 09/22/15 07:35 | 150911L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 93 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 88 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS130 | 15-09-0733-4-I | 09/09/15 09:30 | Aqueous | GC 44 | 09/11/15 | 09/22/15 07:50 | 150911L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 94 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 95 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS140 | 15-09-0733-5-I | 09/09/15 11:12 | Aqueous | GC 44 | 09/11/15 | 09/22/15 08:04 | 150911L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 92 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 89 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS160 | 15-09-0733-6-I | 09/09/15 11:39 | Aqueous | GC 44 | 09/11/15 | 09/22/15 08:18 | 150911L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 93 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 89 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-840 | N/A | Aqueous | GC 44 | 09/19/15 | 09/22/15 06:53 | 150911L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 98 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 97 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS110 | 15-09-0733-1-I | 09/09/15 10:11 | Aqueous | GC 58 | 09/11/15 | 09/21/15 20:18 | 150911L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 81 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 78 | 50-135 | |

| HCS120 | 15-09-0733-2-I | 09/09/15 10:44 | Aqueous | GC 58 | 09/11/15 | 09/21/15 20:36 | 150911L05 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 102 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 98 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHCS120 | 15-09-0733-3-I | 09/09/15 10:44 | Aqueous | GC 58 | 09/11/15 | 09/21/15 20:54 | 150911L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 100 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 94 | 50-135 | |

| | | | | | | | |
|---------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HCS130 | 15-09-0733-4-I | 09/09/15 09:30 | Aqueous | GC 58 | 09/11/15 | 09/21/15 21:12 | 150911L05 |
|---------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 99 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 93 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS140 | 15-09-0733-5-I | 09/09/15 11:12 | Aqueous | GC 58 | 09/11/15 | 09/21/15 21:30 | 150911L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 99 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 94 | 50-135 | |

| HCS160 | 15-09-0733-6-I | 09/09/15 11:39 | Aqueous | GC 58 | 09/11/15 | 09/21/15 21:48 | 150911L05 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 101 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 95 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-533-1086 | N/A | Aqueous | GC 31 | 09/11/15 | 09/14/15 19:25 | 150911L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 94 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 87 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 1 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS110 | 15-09-0733-1-I | 09/09/15 10:11 | Aqueous | GC 35 | 09/11/15 | 09/17/15 20:44 | 150911L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 92 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS120 | 15-09-0733-2-I | 09/09/15 10:44 | Aqueous | GC 35 | 09/11/15 | 09/17/15 21:30 | 150911L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 92 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS120 | 15-09-0733-3-I | 09/09/15 10:44 | Aqueous | GC 35 | 09/11/15 | 09/17/15 22:15 | 150911L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 89 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS130 | 15-09-0733-4-I | 09/09/15 09:30 | Aqueous | GC 35 | 09/11/15 | 09/17/15 23:01 | 150911L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 91 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS140 | 15-09-0733-5-I | 09/09/15 11:12 | Aqueous | GC 35 | 09/11/15 | 09/17/15 23:47 | 150911L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 90 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS160 | 15-09-0733-6-I | 09/09/15 11:39 | Aqueous | GC 35 | 09/11/15 | 09/18/15 00:33 | 150911L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 88 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-111 | N/A | Aqueous | GC 35 | 09/11/15 | 09/17/15 16:54 | 150911L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 93 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS110 | 15-09-0733-1-I | 09/09/15 10:11 | Aqueous | GC 40 | 09/15/15 | 09/21/15 14:53 | 150915L16 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 62 | 0-123 | |

| HCS120 | 15-09-0733-2-I | 09/09/15 10:44 | Aqueous | GC 40 | 09/15/15 | 09/21/15 15:16 | 150915L16 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 53 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHCS120 | 15-09-0733-3-I | 09/09/15 10:44 | Aqueous | GC 40 | 09/15/15 | 09/21/15 15:39 | 150915L16 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 66 | 0-123 | |

| | | | | | | | |
|---------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HCS130 | 15-09-0733-4-I | 09/09/15 09:30 | Aqueous | GC 40 | 09/15/15 | 09/21/15 16:03 | 150915L16 |
|---------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 81 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS140 | 15-09-0733-5-I | 09/09/15 11:12 | Aqueous | GC 40 | 09/15/15 | 09/21/15 16:26 | 150915L16 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 65 | 0-123 | |

| HCS160 | 15-09-0733-6-I | 09/09/15 11:39 | Aqueous | GC 40 | 09/15/15 | 09/21/15 16:49 | 150915L16 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 69 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-034-659 | N/A | Aqueous | GC 40 | 09/15/15 | 09/21/15 14:30 | 150915L16 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 50 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS110 | 15-09-0733-1-H | 09/09/15 10:11 | Aqueous | GC/MS CCC | 09/11/15 | 09/14/15 15:06 | 150911L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

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Method: EPA 8270C
Units: ug/L

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 70 | 33-120 | |
| 2-Fluorophenol | 45 | 24-120 | |
| Nitrobenzene-d5 | 74 | 38-120 | |
| p-Terphenyl-d14 | 81 | 41-137 | |
| Phenol-d6 | 30 | 16-120 | |
| 2,4,6-Tribromophenol | 104 | 27-159 | |


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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 09/10/15
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Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS120 | 15-09-0733-2-H | 09/09/15 10:44 | Aqueous | GC/MS CCC | 09/11/15 | 09/14/15 15:25 | 150911L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 73 | 33-120 | |
| 2-Fluorophenol | 49 | 24-120 | |
| Nitrobenzene-d5 | 78 | 38-120 | |
| p-Terphenyl-d14 | 81 | 41-137 | |
| Phenol-d6 | 31 | 16-120 | |
| 2,4,6-Tribromophenol | 109 | 27-159 | |


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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS120 | 15-09-0733-3-H | 09/09/15 10:44 | Aqueous | GC/MS CCC | 09/11/15 | 09/14/15 15:43 | 150911L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 75 | 33-120 | |
| 2-Fluorophenol | 50 | 24-120 | |
| Nitrobenzene-d5 | 79 | 38-120 | |
| p-Terphenyl-d14 | 87 | 41-137 | |
| Phenol-d6 | 32 | 16-120 | |
| 2,4,6-Tribromophenol | 111 | 27-159 | |


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Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS130 | 15-09-0733-4-H | 09/09/15 09:30 | Aqueous | GC/MS CCC | 09/11/15 | 09/14/15 16:01 | 150911L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 74 | 33-120 | |
| 2-Fluorophenol | 48 | 24-120 | |
| Nitrobenzene-d5 | 78 | 38-120 | |
| p-Terphenyl-d14 | 88 | 41-137 | |
| Phenol-d6 | 31 | 16-120 | |
| 2,4,6-Tribromophenol | 115 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS140 | 15-09-0733-5-H | 09/09/15 11:12 | Aqueous | GC/MS CCC | 09/11/15 | 09/14/15 16:19 | 150911L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 72 | 33-120 | |
| 2-Fluorophenol | 44 | 24-120 | |
| Nitrobenzene-d5 | 76 | 38-120 | |
| p-Terphenyl-d14 | 82 | 41-137 | |
| Phenol-d6 | 29 | 16-120 | |
| 2,4,6-Tribromophenol | 104 | 27-159 | |



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 16 of 21

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS160 | 15-09-0733-6-H | 09/09/15 11:39 | Aqueous | GC/MS CCC | 09/11/15 | 09/14/15 16:37 | 150911L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 78 | 33-120 | |
| 2-Fluorophenol | 49 | 24-120 | |
| Nitrobenzene-d5 | 82 | 38-120 | |
| p-Terphenyl-d14 | 87 | 41-137 | |
| Phenol-d6 | 31 | 16-120 | |
| 2,4,6-Tribromophenol | 114 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-4103 | N/A | Aqueous | GC/MS CCC | 09/11/15 | 09/12/15 13:10 | 150911L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 71 | 33-120 | |
| 2-Fluorophenol | 50 | 24-120 | |
| Nitrobenzene-d5 | 77 | 38-120 | |
| p-Terphenyl-d14 | 86 | 41-137 | |
| Phenol-d6 | 32 | 16-120 | |
| 2,4,6-Tribromophenol | 98 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 1 of 16

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS110 | 15-09-0733-1-B | 09/09/15 10:11 | Aqueous | GC/MS UU | 09/19/15 | 09/19/15 14:57 | 150919L005 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 98 | 78-126 | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS120 | 15-09-0733-2-B | 09/09/15 10:44 | Aqueous | GC/MS UU | 09/19/15 | 09/19/15 15:25 | 150919L005 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 98 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS120 | 15-09-0733-3-B | 09/09/15 10:44 | Aqueous | GC/MS UU | 09/19/15 | 09/19/15 15:54 | 150919L005 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 97 | 78-126 | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS130 | 15-09-0733-4-B | 09/09/15 09:30 | Aqueous | GC/MS UU | 09/19/15 | 09/19/15 16:23 | 150919L005 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 98 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS140 | 15-09-0733-5-B | 09/09/15 11:12 | Aqueous | GC/MS UU | 09/19/15 | 09/19/15 16:51 | 150919L005 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 92 | 80-120 | |
| Dibromofluoromethane | 99 | 78-126 | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS160 | 15-09-0733-6-B | 09/09/15 11:39 | Aqueous | GC/MS UU | 09/19/15 | 09/19/15 17:20 | 150919L005 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 12 of 16

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 90 | 80-120 | |
| Dibromofluoromethane | 97 | 78-126 | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 13 of 16

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| TB11 | 15-09-0733-7-A | 09/09/15 00:00 | Aqueous | GC/MS UU | 09/19/15 | 09/19/15 14:28 | 150919L005 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | 0.25 | 1.0 | 0.24 | 1.00 | J |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 99 | 78-126 | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-18207 | N/A | Aqueous | GC/MS UU | 09/19/15 | 09/19/15 12:02 | 150919L005 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 97 | 78-126 | |
| 1,2-Dichloroethane-d4 | 99 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

09/10/15

Work Order:

15-09-0733

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS110 | 15-09-0733-1 | 09/09/15 10:11 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.085 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 09/15/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 235 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 235 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Solids, Total Dissolved (24) | 275 | 1.00 | 0.870 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 C |
| Solids, Total Suspended (24) | 21 | 1.0 | 0.83 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 D |
| pH (24) | 7.20 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/10/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.0 | 0.50 | 0.28 | 1.00 | | mg/L | 09/21/15 | 09/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 44 | 2.5 | 1.2 | 5.00 | | mg/L | 09/22/15 | 09/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 9.4 | 2.5 | 1.2 | 5.00 | | mg/L | 09/11/15 | 09/12/15 | SM 5310 B |

| | | | |
|---------------|---------------------|-----------------------|----------------|
| HCS120 | 15-09-0733-2 | 09/09/15 10:44 | Aqueous |
|---------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 09/15/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 221 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 221 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Solids, Total Dissolved (24) | 355 | 1.00 | 0.870 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 D |
| pH (24) | 7.23 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/10/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 09/21/15 | 09/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 8.0 | 2.5 | 1.2 | 5.00 | | mg/L | 09/22/15 | 09/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 12 | 2.5 | 1.2 | 5.00 | | mg/L | 09/11/15 | 09/12/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

09/10/15

Work Order:

15-09-0733

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| FDHCS120 | 15-09-0733-3 | 09/09/15 10:44 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 09/15/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 212 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 212 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Solids, Total Dissolved (24) | 340 | 1.00 | 0.870 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 D |
| pH (24) | 7.25 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/10/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 09/21/15 | 09/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.7 | 2.5 | 1.2 | 5.00 | | mg/L | 09/22/15 | 09/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 11 | 2.5 | 1.2 | 5.00 | | mg/L | 09/11/15 | 09/12/15 | SM 5310 B |

| | | | |
|---------------|---------------------|-----------------------|----------------|
| HCS130 | 15-09-0733-4 | 09/09/15 09:30 | Aqueous |
|---------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 09/15/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 186 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 186 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Solids, Total Dissolved (24) | 335 | 1.00 | 0.870 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.6 | 1.0 | 0.83 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 D |
| pH (24) | 7.29 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/10/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 09/21/15 | 09/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.5 | 2.5 | 1.2 | 5.00 | | mg/L | 09/22/15 | 09/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 3.7 | 2.5 | 1.2 | 5.00 | | mg/L | 09/11/15 | 09/12/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

09/10/15

Work Order:

15-09-0733

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS140 | 15-09-0733-5 | 09/09/15 11:12 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 09/15/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 210 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 210 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Solids, Total Dissolved (24) | 325 | 1.00 | 0.870 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 C |
| Solids, Total Suspended (24) | 5.9 | 1.0 | 0.83 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 D |
| pH (24) | 7.54 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/10/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 09/21/15 | 09/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.2 | 2.5 | 1.2 | 5.00 | | mg/L | 09/22/15 | 09/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 5.5 | 2.5 | 1.2 | 5.00 | | mg/L | 09/11/15 | 09/12/15 | SM 5310 B |

| | | | |
|---------------|---------------------|-----------------------|----------------|
| HCS160 | 15-09-0733-6 | 09/09/15 11:39 | Aqueous |
|---------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 09/15/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 210 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 210 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Solids, Total Dissolved (24) | 365 | 1.00 | 0.870 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 D |
| pH (24) | 7.59 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/10/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 09/21/15 | 09/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 4.1 | 2.5 | 1.2 | 5.00 | | mg/L | 09/22/15 | 09/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 8.4 | 2.5 | 1.2 | 5.00 | | mg/L | 09/11/15 | 09/12/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618
Project: EAA 27122

Date Received: 09/10/15
Work Order: 15-09-0733

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|-------------------|---------------------|----------------|
| Method Blank | | N/A | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|-------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 09/15/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/17/15 | SM 2320B |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.87 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 D |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 09/16/15 | 09/16/15 | SM 2540 D |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 09/21/15 | 09/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 09/22/15 | 09/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 09/11/15 | 09/12/15 | SM 5310 B |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-09-0649-1 | Sample | Aqueous | IC 15 | N/A | 09/10/15 14:36 | 150910S01 |
| 15-09-0649-1 | Matrix Spike | Aqueous | IC 15 | N/A | 09/10/15 15:49 | 150910S01 |
| 15-09-0649-1 | Matrix Spike Duplicate | Aqueous | IC 15 | N/A | 09/10/15 16:08 | 150910S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.2027 | 2.500 | 2.351 | 86 | 2.372 | 87 | 80-120 | 1 | 0-20 | |
| Chloride | 636.7 | 50.00 | 818.2 | 363 | 819.6 | 366 | 80-120 | 0 | 0-20 | 3 |
| Bromide | 6.201 | 5.000 | 10.93 | 95 | 11.05 | 97 | 80-120 | 1 | 0-20 | |
| Nitrate (as N) | ND | 5.000 | 4.454 | 89 | 4.583 | 92 | 80-120 | 3 | 0-20 | |
| Sulfate | 275.2 | 50.00 | 423.7 | 297 | 425.4 | 301 | 80-120 | 0 | 0-20 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|-----------|------------|---------------|----------------|---------------------|
| 15-09-0669-1 | Sample | Sea Water | ACA 1 | N/A | 09/15/15 15:53 | 150915S01 |
| 15-09-0669-1 | Matrix Spike | Sea Water | ACA 1 | N/A | 09/15/15 15:53 | 150915S01 |
| 15-09-0669-1 | Matrix Spike Duplicate | Sea Water | ACA 1 | N/A | 09/15/15 15:53 | 150915S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | ND | 0.2000 | 1.432 | 716 | 1.420 | 710 | 90-110 | 1 | 0-25 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-09-1130-1 | Sample | Aqueous | TOC 8 | 09/22/15 | 09/23/15 07:17 | F0922TOCS2 |
| 15-09-1130-1 | Matrix Spike | Aqueous | TOC 8 | 09/22/15 | 09/23/15 07:17 | F0922TOCS2 |
| 15-09-1130-1 | Matrix Spike Duplicate | Aqueous | TOC 8 | 09/22/15 | 09/23/15 07:17 | F0922TOCS2 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Carbon, Total Organic | 8.800 | 50.00 | 56.00 | 94 | 57.50 | 97 | 31-145 | 3 | 0-23 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HCS110 | Sample | Aqueous | TOC 8 | 09/11/15 | 09/12/15 01:31 | F0911DOCS1 | | | | |
| HCS110 | Matrix Spike | Aqueous | TOC 8 | 09/11/15 | 09/12/15 01:31 | F0911DOCS1 | | | | |
| HCS110 | Matrix Spike Duplicate | Aqueous | TOC 8 | 09/11/15 | 09/12/15 01:31 | F0911DOCS1 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Dissolved Organic | 9.400 | 50.00 | 50.00 | 81 | 58.50 | 98 | 31-145 | 16 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HCS110 | Sample | Aqueous | ICP 7300 | 09/11/15 | 09/19/15 23:22 | 150911SA5 |
| HCS110 | Matrix Spike | Aqueous | ICP 7300 | 09/11/15 | 09/19/15 23:25 | 150911SA5 |
| HCS110 | Matrix Spike Duplicate | Aqueous | ICP 7300 | 09/11/15 | 09/19/15 23:29 | 150911SA5 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Calcium | 77.89 | 0.5000 | 74.74 | 4X | 75.25 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 17.53 | 0.5000 | 16.67 | 4X | 16.73 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 1.746 | 5.000 | 7.108 | 107 | 7.289 | 111 | 83-131 | 3 | 0-7 | |
| Sodium | 12.48 | 5.000 | 16.90 | 88 | 17.28 | 96 | 73-127 | 2 | 0-9 | |
| Strontium | 0.7295 | 0.5000 | 1.186 | 91 | 1.218 | 98 | 81-123 | 3 | 0-6 | |
| Silicon | 5.315 | 0.5000 | 5.645 | 4X | 5.696 | 4X | 24-180 | 4X | 0-15 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3020A Total
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-09-0996-1 | Sample | Aqueous | ICP/MS 03 | 09/16/15 | 09/21/15 22:41 | 150916SA3 | | | | |
| 15-09-0996-1 | Matrix Spike | Aqueous | ICP/MS 03 | 09/16/15 | 09/21/15 22:23 | 150916SA3 | | | | |
| 15-09-0996-1 | Matrix Spike Duplicate | Aqueous | ICP/MS 03 | 09/16/15 | 09/21/15 22:27 | 150916SA3 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.1019 | 102 | 0.1039 | 104 | 85-133 | 2 | 0-11 | |
| Arsenic | ND | 0.1000 | 0.1105 | 111 | 0.1102 | 110 | 73-127 | 0 | 0-11 | |
| Barium | 0.1184 | 0.1000 | 0.2322 | 114 | 0.2232 | 105 | 74-128 | 4 | 0-10 | |
| Beryllium | ND | 0.1000 | 0.09766 | 98 | 0.09968 | 100 | 56-122 | 2 | 0-11 | |
| Cadmium | ND | 0.1000 | 0.09744 | 97 | 0.09862 | 99 | 84-114 | 1 | 0-8 | |
| Chromium | 0.03929 | 0.1000 | 0.1530 | 114 | 0.1504 | 111 | 73-133 | 2 | 0-11 | |
| Copper | ND | 0.1000 | 0.1028 | 103 | 0.1023 | 102 | 72-108 | 0 | 0-10 | |
| Lead | ND | 0.1000 | 0.1143 | 114 | 0.1162 | 116 | 79-121 | 2 | 0-10 | |
| Nickel | 0.007515 | 0.1000 | 0.1098 | 102 | 0.1077 | 100 | 68-122 | 2 | 0-10 | |
| Selenium | ND | 0.1000 | 0.09072 | 91 | 0.09158 | 92 | 59-125 | 1 | 0-12 | |
| Silver | ND | 0.05000 | 0.05360 | 107 | 0.05354 | 107 | 68-128 | 0 | 0-14 | |
| Thallium | ND | 0.1000 | 0.1092 | 109 | 0.1095 | 109 | 73-121 | 0 | 0-11 | |
| Zinc | ND | 0.1000 | 0.09848 | 98 | 0.09586 | 96 | 43-145 | 3 | 0-39 | |
| Aluminum | ND | 0.1000 | 0.1149 | 115 | 0.1097 | 110 | 47-161 | 5 | 0-24 | |
| Iron | ND | 5.100 | 5.816 | 114 | 5.668 | 111 | 27-201 | 3 | 0-24 | |
| Manganese | 0.01195 | 0.1000 | 0.1214 | 109 | 0.1200 | 108 | 72-126 | 1 | 0-42 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 7470A Total
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-09-0914-1 | Sample | Aqueous | Mercury 04 | 09/15/15 | 09/15/15 20:13 | 150915SA3 |
| 15-09-0914-1 | Matrix Spike | Aqueous | Mercury 04 | 09/15/15 | 09/15/15 20:20 | 150915SA3 |
| 15-09-0914-1 | Matrix Spike Duplicate | Aqueous | Mercury 04 | 09/15/15 | 09/15/15 20:22 | 150915SA3 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | ND | 0.01000 | 0.01019 | 102 | 0.009988 | 100 | 57-141 | 2 | 0-10 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3020A Total
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|--------------|-------------|------------|----------------|----------------|-----------------------|
| 15-09-0996-1 | Sample | Aqueous | ICP/MS 03 | 09/16/15 00:00 | 09/21/15 22:41 | 150916SA3 |
| 15-09-0996-1 | PDS | Aqueous | ICP/MS 03 | 09/16/15 00:00 | 09/21/15 22:30 | 150916SA3 |
| Parameter | Sample Conc. | Spike Added | PDS Conc. | PDS %Rec. | %Rec. CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.09812 | 98 | 75-125 | |
| Arsenic | ND | 0.1000 | 0.1027 | 103 | 75-125 | |
| Barium | 0.1184 | 0.1000 | 0.2189 | 101 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.09334 | 93 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.09263 | 93 | 75-125 | |
| Chromium | 0.03929 | 0.1000 | 0.1428 | 103 | 75-125 | |
| Copper | ND | 0.1000 | 0.09691 | 97 | 75-125 | |
| Lead | ND | 0.1000 | 0.1092 | 109 | 75-125 | |
| Nickel | 0.007515 | 0.1000 | 0.1032 | 96 | 75-125 | |
| Selenium | ND | 0.1000 | 0.08516 | 85 | 75-125 | |
| Silver | ND | 0.05000 | 0.04834 | 97 | 75-125 | |
| Thallium | ND | 0.1000 | 0.1033 | 103 | 75-125 | |
| Zinc | ND | 0.1000 | 0.08546 | 85 | 75-125 | |
| Aluminum | ND | 0.1000 | 0.1046 | 105 | 75-125 | |
| Iron | ND | 5.100 | 5.302 | 104 | 75-125 | |
| Manganese | 0.01195 | 0.1000 | 0.1129 | 101 | 75-125 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-09-1300-2 | Sample | Aqueous | PH1/BUR03 | N/A | 09/17/15 14:00 | F0917ALKD2 |
| 15-09-1300-2 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 09/17/15 14:00 | F0917ALKD2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|---|---------------------|------------------|------------|---------------|-------------------|
| Alkalinity, Total (as CaCO ₃) | 105.0 | 106.0 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|-------------------------------------|------------------|---------------------|------------------|---------------|----------------|------------------------|
| 15-09-1300-2 | Sample | Aqueous | PH1/BUR03 | N/A | 09/17/15 14:00 | F0917HCOD2 |
| 15-09-1300-2 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 09/17/15 14:00 | F0917HCOD2 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Bicarbonate (as CaCO ₃) | | 105.0 | 106.0 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-09-1300-2 | Sample | Aqueous | PH1/BUR03 | N/A | 09/17/15 14:00 | F0917CO3D2 |
| 15-09-1300-2 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 09/17/15 14:00 | F0917CO3D2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-09-0675-1 | Sample | Aqueous | N/A | 09/16/15 00:00 | 09/16/15 20:00 | F0916TDSD6 |
| 15-09-0675-1 | Sample Duplicate | Aqueous | N/A | 09/16/15 00:00 | 09/16/15 20:00 | F0916TDSD6 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Dissolved | 685.0 | 725.0 | 6 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| HCS110 | Sample | Aqueous | N/A | 09/16/15 00:00 | 09/16/15 18:00 | F0916TSSD2 |
| HCS110 | Sample Duplicate | Aqueous | N/A | 09/16/15 00:00 | 09/16/15 18:00 | F0916TSSD2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Suspended | 20.90 | 22.30 | 6 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-09-0680-1 | Sample | Aqueous | PH 1 | N/A | 09/10/15 19:57 | F0910PHD5 |
| 15-09-0680-1 | Sample Duplicate | Aqueous | PH 1 | N/A | 09/10/15 19:57 | F0910PHD5 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|------------------|---------------------|------------------|------------|---------------|-------------------|
| pH | 5.760 | 5.770 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 4500 N Org B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-09-1091-2 | Sample | Aqueous | BUR05 | 09/21/15 00:00 | 09/21/15 19:34 | F0921TKND2 |
| 15-09-1091-2 | Sample Duplicate | Aqueous | BUR05 | 09/21/15 00:00 | 09/21/15 19:34 | F0921TKND2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Total Kjeldahl Nitrogen | 2.520 | 2.380 | 6 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|--------------------|------------------------|------------------|-----------------|-------------------|
| 099-12-906-6058 | LCS | Aqueous | IC 15 | N/A | 09/10/15 12:52 | 150910L01 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Fluoride | | 2.500 | 2.485 | 99 | 90-110 | |
| Chloride | | 50.00 | 48.38 | 97 | 90-110 | |
| Bromide | | 5.000 | 4.928 | 99 | 90-110 | |
| Nitrate (as N) | | 5.000 | 4.912 | 98 | 90-110 | |
| Sulfate | | 50.00 | 48.63 | 97 | 90-110 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-739-116 | LCS | Aqueous | ACA 1 | N/A | 09/15/15 15:53 | 150915L01 | | | |
| 099-12-739-116 | LCSD | Aqueous | ACA 1 | N/A | 09/15/15 15:53 | 150915L01 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Phosphorus, Total | 0.2000 | 0.1967 | 98 | 0.2032 | 102 | 90-110 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-859-797 | LCS | Aqueous | PH1/BUR03 | N/A | 09/17/15 14:00 | F0917ALKB2 | | | |
| 099-15-859-797 | LCSD | Aqueous | PH1/BUR03 | N/A | 09/17/15 14:00 | F0917ALKB2 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 100.0 | 99.00 | 99 | 99.00 | 99 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4750 | LCS | Aqueous | N/A | 09/16/15 | 09/16/15 20:00 | F0916TDSL6 | | | |
| 099-12-180-4750 | LCSD | Aqueous | N/A | 09/16/15 | 09/16/15 20:00 | F0916TDSL6 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 105.0 | 105 | 110.0 | 110 | 80-120 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-09-010-7314 | LCS | Aqueous | N/A | 09/16/15 | 09/16/15 18:00 | F0916TSSL2 | | | |
| 099-09-010-7314 | LCSD | Aqueous | N/A | 09/16/15 | 09/16/15 18:00 | F0916TSSL2 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Suspended | 100.0 | 86.00 | 86 | 89.00 | 89 | 80-120 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-09-010-7321 | LCS | Aqueous | N/A | 09/16/15 | 09/16/15 18:00 | F0916TSSL2 | | | |
| 099-09-010-7321 | LCSD | Aqueous | N/A | 09/16/15 | 09/16/15 18:00 | F0916TSSL2 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Suspended | 100.0 | 86.00 | 86 | 89.00 | 89 | 80-120 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-097-5756 | LCS | Aqueous | TOC 8 | 09/22/15 | 09/23/15 07:17 | F0922TOCL2 | | | |
| 099-05-097-5756 | LCSD | Aqueous | TOC 8 | 09/22/15 | 09/23/15 07:17 | F0922TOCL2 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 10.00 | 10.10 | 101 | 10.40 | 104 | 80-120 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------------|------------|---------------|-------------------|
| 099-05-115-1455 | LCS | Aqueous | TOC 8 | 09/11/15 | 09/12/15 01:31 | F0911DOCL1 | | | |
| 099-05-115-1455 | LCSD | Aqueous | TOC 8 | 09/11/15 | 09/12/15 01:31 | F0911DOCL1 | | | |
| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Carbon, Dissolved Organic | 10.00 | 10.20 | 102 | 10.60 | 106 | 80-120 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-15-683-1396 | LCS | Aqueous | ICP 7300 | 09/11/15 | 09/22/15 19:12 | 150911LA5F |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Calcium | | 0.5000 | 0.5386 | 108 | 80-120 | |
| Magnesium | | 0.5000 | 0.4963 | 99 | 80-120 | |
| Potassium | | 5.000 | 5.550 | 111 | 80-120 | |
| Sodium | | 5.000 | 5.873 | 117 | 80-120 | |
| Strontium | | 0.5000 | 0.5977 | 120 | 80-120 | |
| Silicon | | 0.5000 | 0.5120 | 102 | 80-120 | |


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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-912 | LCS | Aqueous | ICP/MS 03 | 09/16/15 | 09/21/15 22:20 | 150916LA3F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.09960 | 100 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.1031 | 103 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.1002 | 100 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.09791 | 98 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.1012 | 101 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.1059 | 106 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.1010 | 101 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.1049 | 105 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.09570 | 96 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.1041 | 104 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.05052 | 101 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.09719 | 97 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.1005 | 100 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.1055 | 106 | 80-120 | 73-127 | |
| Iron | 5.100 | 4.996 | 98 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.1009 | 101 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| 099-15-763-614 | LCS | Aqueous | Mercury 04 | 09/15/15 | 09/15/15 20:11 | 150915LA3F |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.01000 | 0.01027 | 103 | 85-121 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-529-840 | LCS | Aqueous | | GC 44 | 09/19/15 | 09/22/15 13:34 | 150911L04 | | | |
| 099-12-529-840 | LCSD | Aqueous | | GC 44 | 09/19/15 | 09/22/15 13:48 | 150911L04 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Alpha-BHC | 0.5000 | 0.4451 | 89 | 0.4246 | 85 | 50-135 | 36-149 | 5 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.4727 | 95 | 0.4481 | 90 | 50-135 | 36-149 | 5 | 0-25 | |
| Beta-BHC | 0.5000 | 0.4224 | 84 | 0.4299 | 86 | 50-135 | 36-149 | 2 | 0-25 | |
| Heptachlor | 0.5000 | 0.4651 | 93 | 0.4583 | 92 | 50-135 | 36-149 | 1 | 0-25 | |
| Delta-BHC | 0.5000 | 0.4682 | 94 | 0.4324 | 86 | 50-135 | 36-149 | 8 | 0-25 | |
| Aldrin | 0.5000 | 0.4277 | 86 | 0.4393 | 88 | 50-135 | 36-149 | 3 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.4342 | 87 | 0.4539 | 91 | 50-135 | 36-149 | 4 | 0-25 | |
| Endosulfan I | 0.5000 | 0.4562 | 91 | 0.4576 | 92 | 50-135 | 36-149 | 0 | 0-25 | |
| Dieldrin | 0.5000 | 0.4759 | 95 | 0.4739 | 95 | 50-135 | 36-149 | 0 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.4651 | 93 | 0.4617 | 92 | 50-135 | 36-149 | 1 | 0-25 | |
| Endrin | 0.5000 | 0.5162 | 103 | 0.4822 | 96 | 50-135 | 36-149 | 7 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.4893 | 98 | 0.4487 | 90 | 50-135 | 36-149 | 9 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.4815 | 96 | 0.4660 | 93 | 50-135 | 36-149 | 3 | 0-25 | |
| Endosulfan II | 0.5000 | 0.4956 | 99 | 0.4746 | 95 | 50-135 | 36-149 | 4 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.5059 | 101 | 0.4420 | 88 | 50-135 | 36-149 | 13 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.4915 | 98 | 0.4503 | 90 | 50-135 | 36-149 | 9 | 0-25 | |
| Methoxychlor | 0.5000 | 0.5191 | 104 | 0.4463 | 89 | 50-135 | 36-149 | 15 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-533-1086 | LCS | Aqueous | GC 31 | 09/11/15 | 09/14/15 18:47 | 150911L05 | | | |
| 099-12-533-1086 | LCSD | Aqueous | GC 31 | 09/11/15 | 09/14/15 19:06 | 150911L05 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Aroclor-1016 | 2.000 | 1.563 | 78 | 1.903 | 95 | 50-135 | 20 | 0-25 | |
| Aroclor-1260 | 2.000 | 1.588 | 79 | 1.738 | 87 | 50-135 | 9 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-963-111 | LCS | Aqueous | | GC 35 | 09/11/15 | 09/17/15 19:12 | 150911L03 | | | |
| 099-15-963-111 | LCSD | Aqueous | | GC 35 | 09/11/15 | 09/17/15 19:58 | 150911L03 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.03912 | 98 | 0.04060 | 102 | 30-130 | 13-147 | 4 | 0-30 | |
| Bolstar | 0.04000 | 0.03662 | 92 | 0.03780 | 94 | 30-130 | 13-147 | 3 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.03560 | 89 | 0.03692 | 92 | 30-130 | 13-147 | 4 | 0-30 | |
| Coumaphos | 0.04000 | 0.03518 | 88 | 0.03616 | 90 | 30-130 | 13-147 | 3 | 0-30 | |
| Diazinon | 0.04000 | 0.03660 | 92 | 0.03782 | 95 | 30-130 | 13-147 | 3 | 0-30 | |
| Disulfoton | 0.04000 | 0.03686 | 92 | 0.03767 | 94 | 30-130 | 13-147 | 2 | 0-30 | |
| Ethoprop | 0.04000 | 0.04376 | 109 | 0.04544 | 114 | 30-130 | 13-147 | 4 | 0-30 | |
| Fensulfothion | 0.04000 | 0.03714 | 93 | 0.03823 | 96 | 30-130 | 13-147 | 3 | 0-30 | |
| Fenthion | 0.04000 | 0.03803 | 95 | 0.03950 | 99 | 30-130 | 13-147 | 4 | 0-30 | |
| Merphos | 0.04000 | 0.07192 | 180 | 0.07487 | 187 | 30-130 | 13-147 | 4 | 0-30 | X |
| Methyl Parathion | 0.04000 | 0.03678 | 92 | 0.03844 | 96 | 30-130 | 13-147 | 4 | 0-30 | |
| Phorate | 0.04000 | 0.04420 | 110 | 0.04605 | 115 | 30-130 | 13-147 | 4 | 0-30 | |
| Ronnel | 0.04000 | 0.03562 | 89 | 0.03720 | 93 | 30-130 | 13-147 | 4 | 0-30 | |
| Stirophos | 0.04000 | 0.03723 | 93 | 0.03876 | 97 | 30-130 | 13-147 | 4 | 0-30 | |
| Tokuthion | 0.04000 | 0.03568 | 89 | 0.03698 | 92 | 30-130 | 13-147 | 4 | 0-30 | |
| Trichloronate | 0.04000 | 0.03803 | 95 | 0.03932 | 98 | 30-130 | 13-147 | 3 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-034-659 | LCS | Aqueous | GC 40 | 09/15/15 | 09/21/15 13:43 | 150915L16 | | | |
| 095-01-034-659 | LCSD | Aqueous | GC 40 | 09/15/15 | 09/21/15 14:07 | 150915L16 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| 2,4-D | 20.00 | 8.370 | 42 | 9.565 | 48 | 30-130 | 13 | 0-30 | |
| 2,4,5-T | 2.000 | 1.090 | 54 | 1.255 | 63 | 30-130 | 14 | 0-30 | |
| 2,4-DB | 20.00 | 7.540 | 38 | 8.855 | 44 | 30-130 | 16 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-4103 | LCS | Aqueous | | GC/MS CCC | 09/11/15 | 09/12/15 13:28 | 150911L06 | | | |
| 095-01-003-4103 | LCSD | Aqueous | | GC/MS CCC | 09/11/15 | 09/12/15 13:46 | 150911L06 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 200.0 | 185.6 | 93 | 181.6 | 91 | 61-120 | 51-130 | 2 | 0-20 | |
| Acenaphthylene | 200.0 | 180.6 | 90 | 177.8 | 89 | 55-120 | 44-131 | 2 | 0-20 | |
| Butyl Benzyl Phthalate | 200.0 | 173.4 | 87 | 171.2 | 86 | 56-122 | 45-133 | 1 | 0-20 | |
| 4-Chloro-3-Methylphenol | 200.0 | 189.3 | 95 | 182.9 | 91 | 52-120 | 41-131 | 3 | 0-20 | |
| 2-Chlorophenol | 200.0 | 177.3 | 89 | 167.6 | 84 | 47-120 | 35-132 | 6 | 0-20 | |
| 1,4-Dichlorobenzene | 200.0 | 160.3 | 80 | 155.0 | 78 | 36-120 | 22-134 | 3 | 0-20 | |
| Dimethyl Phthalate | 200.0 | 195.1 | 98 | 194.5 | 97 | 60-120 | 50-130 | 0 | 0-20 | |
| 2,4-Dinitrotoluene | 200.0 | 212.6 | 106 | 211.2 | 106 | 61-121 | 51-131 | 1 | 0-20 | |
| Fluorene | 200.0 | 194.9 | 97 | 193.3 | 97 | 67-120 | 58-129 | 1 | 0-20 | |
| N-Nitroso-di-n-propylamine | 200.0 | 173.5 | 87 | 160.8 | 80 | 39-123 | 25-137 | 8 | 0-20 | |
| Naphthalene | 200.0 | 174.4 | 87 | 164.2 | 82 | 54-120 | 43-131 | 6 | 0-20 | |
| 4-Nitrophenol | 200.0 | 81.69 | 41 | 84.62 | 42 | 14-120 | 0-138 | 4 | 0-20 | |
| Pentachlorophenol | 200.0 | 180.4 | 90 | 182.5 | 91 | 31-127 | 15-143 | 1 | 0-20 | |
| Phenol | 200.0 | 78.24 | 39 | 75.35 | 38 | 17-120 | 0-137 | 4 | 0-20 | |
| Pyrene | 200.0 | 181.8 | 91 | 178.9 | 89 | 58-124 | 47-135 | 2 | 0-20 | |
| 1,2,4-Trichlorobenzene | 200.0 | 185.6 | 93 | 174.3 | 87 | 49-120 | 37-132 | 6 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/10/15
Work Order: 15-09-0733
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|-----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-14-001-18207 | LCS | Aqueous | | GC/MS UU | 09/19/15 | 09/19/15 10:33 | 150919L005 | | | |
| 099-14-001-18207 | LCSD | Aqueous | | GC/MS UU | 09/19/15 | 09/19/15 11:02 | 150919L005 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Benzene | 50.00 | 50.49 | 101 | 52.88 | 106 | 80-120 | 73-127 | 5 | 0-20 | |
| Carbon Tetrachloride | 50.00 | 50.48 | 101 | 54.72 | 109 | 67-139 | 55-151 | 8 | 0-20 | |
| Chlorobenzene | 50.00 | 50.23 | 100 | 51.12 | 102 | 78-120 | 71-127 | 2 | 0-20 | |
| 1,2-Dibromoethane | 50.00 | 49.71 | 99 | 50.43 | 101 | 80-120 | 73-127 | 1 | 0-20 | |
| 1,2-Dichlorobenzene | 50.00 | 51.16 | 102 | 52.27 | 105 | 63-129 | 52-140 | 2 | 0-20 | |
| 1,2-Dichloroethane | 50.00 | 51.70 | 103 | 51.52 | 103 | 70-130 | 60-140 | 0 | 0-20 | |
| 1,1-Dichloroethene | 50.00 | 46.27 | 93 | 51.12 | 102 | 66-126 | 56-136 | 10 | 0-20 | |
| Ethylbenzene | 50.00 | 53.51 | 107 | 55.83 | 112 | 80-123 | 73-130 | 4 | 0-20 | |
| Toluene | 50.00 | 49.48 | 99 | 52.24 | 104 | 80-120 | 73-127 | 5 | 0-20 | |
| Trichloroethene | 50.00 | 47.14 | 94 | 49.53 | 99 | 80-122 | 73-129 | 5 | 0-20 | |
| Vinyl Chloride | 50.00 | 46.71 | 93 | 51.15 | 102 | 70-130 | 60-140 | 9 | 0-20 | |
| p/m-Xylene | 100.0 | 108.7 | 109 | 112.1 | 112 | 75-123 | 67-131 | 3 | 0-20 | |
| o-Xylene | 50.00 | 53.71 | 107 | 54.92 | 110 | 74-122 | 66-130 | 2 | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 47.96 | 96 | 48.23 | 96 | 69-129 | 59-139 | 1 | 0-20 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Sample Analysis Summary Report

Work Order: 15-09-0733

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| <u>Method</u> | <u>Extraction</u> | <u>Chemist ID</u> | <u>Instrument</u> | <u>Analytical Location</u> |
|-----------------|-------------------|-------------------|-------------------|----------------------------|
| EPA 300.0 | N/A | 1027 | IC 15 | 1 |
| EPA 365.1 | N/A | 650 | ACA 1 | 1 |
| EPA 6010B | EPA 3005A Filt. | 935 | ICP 7300 | 1 |
| EPA 6020 | EPA 3005A Filt. | 598 | ICP/MS 03 | 1 |
| EPA 7470A | EPA 7470A Filt. | 915 | Mercury 04 | 1 |
| EPA 8081A | EPA 3510C | 669 | GC 44 | 1 |
| EPA 8082 | EPA 3510C | 944 | GC 58 | 1 |
| EPA 8141A | EPA 3510C | 960 | GC 35 | 1 |
| EPA 8151A | EPA 8151A | 944 | GC 40 | 1 |
| EPA 8260B | EPA 5030C | 823 | GC/MS UU | 2 |
| EPA 8270C | EPA 3510C | 923 | GC/MS CCC | 1 |
| SM 2320B | N/A | 650 | PH1/BUR03 | 1 |
| SM 2540 C | N/A | 1009 | N/A | 1 |
| SM 2540 D | N/A | 1009 | N/A | 1 |
| SM 4500 H+ B | N/A | 688 | PH 1 | 1 |
| SM 4500 N Org B | N/A | 685 | BUR05 | 1 |
| SM 5310 B | N/A | 735 | TOC 8 | 1 |


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Location 1: 7440 Lincoln Way, Garden Grove, CA 92841

Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841

Glossary of Terms and Qualifiers

Work Order: 15-09-0733

Page 1 of 1

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|---|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| CI | See case narrative. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |
| | Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time. |
| | A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations. |

SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 6

CLIENT: SWCA

DATE: 09/10/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF: -0.2°C); Temperature (w/o CF): 2.7 °C (w/ CF): 2.5 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by:)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A Checked by: 1013

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: 150121E)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOAn₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PBz_{na} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☒ 250PB_h ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGB_s ☐ 1PB ☐ 1PBna ☒ 3.5 gal cube ☐ ☐ ☐

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve () ☐ EnCores® () ☐ TerraCores® () ☐

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ Other Matrix (): ☐ ☐

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1013

s = H₂SO₄, u = ultra-pure, z_{na} = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 619

SAMPLE RECEIPT CHECKLIST

COOLER 2 OF 6

CLIENT: SWCA

DATE: 09 / 10 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF:-0.2°C); Temperature (w/o CF): 2.4 °C (w/ CF): 2.2 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1013

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE: (Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznn ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☒ 250PB ☐ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal. ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1013

s = H₂SO₄, u = ultra-pure, znn = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 1013

SAMPLE RECEIPT CHECKLIST

COOLER 3 OF 6

CLIENT: SWCA

DATE: 09 / 10 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF:-0.2°C); Temperature (w/o CF): 2.6 °C (w/ CF): 2.4 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1013

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☒ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBz_{na} ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal. Cube ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1013s = H₂SO₄, u = ultra-pure, z_{na} = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 679

SAMPLE RECEIPT CHECKLIST

COOLER 4 OF 6

CLIENT: SWCA

DATE: 09 / 10 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF: -0.2°C); Temperature (w/o CF): 2.2 °C (w/ CF): 2.0 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1013

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznn ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 25 gal. ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1013s = H₂SO₄, u = ultra-pure, znn = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 659

SAMPLE RECEIPT CHECKLIST

COOLER 5 OF 6

CLIENT: SWCA

DATE: 09/10/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF: -0.2°C); Temperature (w/o CF): 3.1 °C (w/ CF): 2.9 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1012

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☒ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: 3 (Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na2} ☐ 100PJ ☐ 100PJ_{na2} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☒ 250PB_h ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s☐ 500PB ☒ 1AGB ☐ 1AGB_{na2} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5 gnl. ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1013s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 619

SAMPLE RECEIPT CHECKLIST

COOLER 6 OF 6

CLIENT: SWCA

DATE: 09 / 10 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF:-0.2°C); Temperature (w/o CF): 2.0 °C (w/ CF): 1.8 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1013

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOAn₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB
☐ 125PBz_{nn} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☒ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s
☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGB_s ☐ 1PB ☐ 1PBna ☒ 2.5 gal. cube ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1013

s = H₂SO₄, u = ultra-pure, z_{nn}a = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 659



Calscience

Subcontractor Analysis Report

Work Order: 15-09-0733Page 1 of 1

One or more samples in this work order have tests that were subcontracted. The subcontract report(s) follows.

For subcontracted tests, please reference the laboratory information noted below.

1. ALS - Columbia Analytical Services, Inc. - Kelso, WA CA ELAP 2286, NELAP WA100010
Method 1694 Caffeine

A blue upward-pointing arrow icon.
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ALS Environmental
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September 30, 2015

Analytical Report for Service Request No: K1510108

Donald Burley
CalScience Environmental Laboratories, Incorporated
7440 Lincoln Way
Garden Grove, CA 92841

RE: 15-09-0733

Dear Donald,

Enclosed are the results of the sample(s) submitted to our laboratory September 12, 2015
For your reference, these analyses have been assigned our service request number **K1510108**.

Analyses were performed according to our laboratory's NELAP-approved quality assurance program. The test results meet requirements of the current NELAP standards, where applicable, and except as noted in the laboratory case narrative provided. For a specific list of NELAP-accredited analytes, refer to the certifications section at www.alsglobal.com. All results are intended to be considered in their entirety, and ALS Group USA Corp. dba ALS Environmental (ALS) is not responsible for use of less than the complete report. Results apply only to the items submitted to the laboratory for analysis and individual items (samples) analyzed, as listed in the report.

Please contact me if you have any questions. My extension is 3364. You may also contact me via email at howard.holmes@alsglobal.com.

Respectfully submitted,

ALS Group USA, Corp. dba ALS Environmental

Howard Holmes
Project Manager

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Acronyms

| | |
|------------|--|
| ASTM | American Society for Testing and Materials |
| A2LA | American Association for Laboratory Accreditation |
| CARB | California Air Resources Board |
| CAS Number | Chemical Abstract Service registry Number |
| CFC | Chlorofluorocarbon |
| CFU | Colony-Forming Unit |
| DEC | Department of Environmental Conservation |
| DEQ | Department of Environmental Quality |
| DHS | Department of Health Services |
| DOE | Department of Ecology |
| DOH | Department of Health |
| EPA | U. S. Environmental Protection Agency |
| ELAP | Environmental Laboratory Accreditation Program |
| GC | Gas Chromatography |
| GC/MS | Gas Chromatography/Mass Spectrometry |
| LOD | Limit of Detection |
| LOQ | Limit of Quantitation |
| LUFT | Leaking Underground Fuel Tank |
| M | Modified |
| MCL | Maximum Contaminant Level is the highest permissible concentration of a substance allowed in drinking water as established by the USEPA. |
| MDL | Method Detection Limit |
| MPN | Most Probable Number |
| MRL | Method Reporting Limit |
| NA | Not Applicable |
| NC | Not Calculated |
| NCASI | National Council of the Paper Industry for Air and Stream Improvement |
| ND | Not Detected |
| NIOSH | National Institute for Occupational Safety and Health |
| PQL | Practical Quantitation Limit |
| RCRA | Resource Conservation and Recovery Act |
| SIM | Selected Ion Monitoring |
| TPH | Total Petroleum Hydrocarbons |
| tr | Trace level is the concentration of an analyte that is less than the PQL but greater than or equal to the MDL. |

Inorganic Data Qualifiers

- * The result is an outlier. See case narrative.
- # The control limit criteria is not applicable. See case narrative.
- B The analyte was found in the associated method blank at a level that is significant relative to the sample result as defined by the DOD or NELAC standards.
- E The result is an estimate amount because the value exceeded the instrument calibration range.
- J The result is an estimated value.
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.
DOD-QSM 4.2 definition : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- i The MRL/MDL or LOQ/LOD is elevated due to a matrix interference.
- X See case narrative.
- Q See case narrative. One or more quality control criteria was outside the limits.
- H The holding time for this test is immediately following sample collection. The samples were analyzed as soon as possible after receipt by the laboratory.

Metals Data Qualifiers

- # The control limit criteria is not applicable. See case narrative.
- J The result is an estimated value.
- E The percent difference for the serial dilution was greater than 10%, indicating a possible matrix interference in the sample.
- M The duplicate injection precision was not met.
- N The Matrix Spike sample recovery is not within control limits. See case narrative.
- S The reported value was determined by the Method of Standard Additions (MSA).
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.
DOD-QSM 4.2 definition : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- W The post-digestion spike for furnace AA analysis is out of control limits, while sample absorbance is less than 50% of spike absorbance.
- i The MRL/MDL or LOQ/LOD is elevated due to a matrix interference.
- X See case narrative.
- + The correlation coefficient for the MSA is less than 0.995.
- Q See case narrative. One or more quality control criteria was outside the limits.

Organic Data Qualifiers

- * The result is an outlier. See case narrative.
- # The control limit criteria is not applicable. See case narrative.
- A A tentatively identified compound, a suspected aldol-condensation product.
- B The analyte was found in the associated method blank at a level that is significant relative to the sample result as defined by the DOD or NELAC standards.
- C The analyte was qualitatively confirmed using GC/MS techniques, pattern recognition, or by comparing to historical data.
- D The reported result is from a dilution.
- E The result is an estimated value.
- J The result is an estimated value.
- N The result is presumptive. The analyte was tentatively identified, but a confirmation analysis was not performed.
- P The GC or HPLC confirmation criteria was exceeded. The relative percent difference is greater than 40% between the two analytical results.
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.
DOD-QSM 4.2 definition : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- i The MRL/MDL or LOQ/LOD is elevated due to a chromatographic interference.
- X See case narrative.
- Q See case narrative. One or more quality control criteria was outside the limits.

Additional Petroleum Hydrocarbon Specific Qualifiers

- F The chromatographic fingerprint of the sample matches the elution pattern of the calibration standard.
- L The chromatographic fingerprint of the sample resembles a petroleum product, but the elution pattern indicates the presence of a greater amount of lighter molecular weight constituents than the calibration standard.
- H The chromatographic fingerprint of the sample resembles a petroleum product, but the elution pattern indicates the presence of a greater amount of heavier molecular weight constituents than the calibration standard.
- O The chromatographic fingerprint of the sample resembles an oil, but does not match the calibration standard.
- Y The chromatographic fingerprint of the sample resembles a petroleum product eluting in approximately the correct carbon range, but the elution pattern does not match the calibration standard.
- Z The chromatographic fingerprint does not resemble a petroleum product.

**ALS Group USA Corp. dba ALS Environmental (ALS) - Kelso
State Certifications, Accreditations, and Licenses**

| Agency | Web Site | Number |
|--------------------------|---|---------------|
| Alaska DEC UST | http://dec.alaska.gov/applications/eh/ehllabreports/USTLabs.aspx | UST-040 |
| Arizona DHS | http://www.azdhs.gov/lab/license/env.htm | AZ0339 |
| Arkansas - DEQ | http://www.adeq.state.ar.us/techsvs/labcert.htm | 88-0637 |
| California DHS (ELAP) | http://www.cdph.ca.gov/certlic/labs/Pages/ELAP.aspx | 2795 |
| DOD ELAP | http://www.denix.osd.mil/edqw/Accreditation/AccreditedLabs.cfm | L14-51 |
| Florida DOH | http://www.doh.state.fl.us/lab/EnvLabCert/WaterCert.htm | E87412 |
| Hawaii DOH | Not available | - |
| Idaho DHW | http://www.healthandwelfare.idaho.gov/Health/Labs/CertificationDrinkingWaterLabs/tabid/1833/Default.aspx | - |
| ISO 17025 | http://www.pjlabs.com/ | L14-50 |
| Louisiana DEQ | http://www.deq.louisiana.gov/portal/DIVISIONS/PublicParticipationandPermitSupport/LouisianaLaboratoryAccreditationProgram.aspx | 03016 |
| Maine DHS | Not available | WA01276 |
| Michigan DEQ | http://www.michigan.gov/deq/0,1607,7-135-3307_4131_4156---,00.html | 9949 |
| Minnesota DOH | http://www.health.state.mn.us/accreditation | 053-999-457 |
| Montana DPHHS | http://www.dphhs.mt.gov/publichealth/ | CERT0047 |
| Nevada DEP | http://ndep.nv.gov/bsdwlabservice.htm | WA01276 |
| New Jersey DEP | http://www.nj.gov/dep/oqa/ | WA005 |
| North Carolina DWQ | http://www.dwqlab.org/ | 605 |
| Oklahoma DEQ | http://www.deq.state.ok.us/CSDnew/labcert.htm | 9801 |
| Oregon – DEQ (NELAP) | http://public.health.oregon.gov/LaboratoryServices/EnvironmentalLaboratoryAccreditation/Pages/index.aspx | WA100010 |
| South Carolina DHEC | http://www.scdhec.gov/environment/envserv/ | 61002 |
| Texas CEQ | http://www.tceq.texas.gov/field/qa/env_lab_accreditation.html | T104704427 |
| Washington DOE | http://www.ecy.wa.gov/programs/eap/labs/lab-accreditation.html | C544 |
| Wisconsin DNR | http://dnr.wi.gov/ | 998386840 |
| Wyoming (EPA Region 8) | http://www.epa.gov/region8/water/dwhome/wyomingdi.html | - |
| Kelso Laboratory Website | www.alsglobal.com | NA |

Analyses were performed according to our laboratory's NELAP-approved quality assurance program. A complete listing of specific NELAP-certified analytes, can be found in the certification section at www.ALSGlobal.com or at the accreditation bodies web site.

Please refer to the certification and/or accreditation body's web site if samples are submitted for compliance purposes. The states highlighted above, require the analysis be listed on the state certification if used for compliance purposes and if the method/analyte is offered by that state.



Case Narrative

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ALS Environmental—Kelso Laboratory
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RIGHT SOLUTIONS | RIGHT PARTNER

ALS ENVIRONMENTAL

| | | | |
|-----------------------|--|-----------------------------|----------|
| Client: | Eurofins Calscience Environmental Laboratory | Service Request No.: | K1510108 |
| Project: | 15-09-0733 | Date Received: | 09/12/15 |
| Sample Matrix: | Water | | |

Case Narrative

All analyses were performed consistent with the quality assurance program of ALS Environmental. This report contains analytical results for samples designated for Tier II data deliverables. When appropriate to the method, method blank results have been reported with each analytical test. Surrogate recoveries have been reported for all applicable organic analyses. Additional quality control analyses reported herein include: Laboratory Control Sample (LCS), and Laboratory/Duplicate Laboratory Control Sample (LCS/DLCS).

Sample Receipt

Six water samples were received for analysis at ALS Environmental on 09/12/15. The samples were received in good condition and consistent with the accompanying chain of custody form. The samples were stored in a refrigerator at 4°C upon receipt at the laboratory.

Steroids and Endocrine Disrupting Compounds by Method 1694**Elevated Detection Limits:**

The detection limit for Caffeine for all samples was elevated due to less than optimal sample volume available for analysis.

No other anomalies associated with the analysis of these samples were observed.

Approved by _____





Chain of Custody

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RIGHT SOLUTIONS : RIGHT PARTNER

PC H2

Cooler Receipt and Preservation Form

Client / Project: CALSCIENCE Service Request K15 10108
 Received: 9/12/15 Opened: 9/12/15 By: L Unloaded: 9/12/15 By: L

1. Samples were received via? Mail Fed Ex UPS DIHL PDX Courier Hand Delivered
 2. Samples were received in: (circle) Cooler Box Envelope Other NA
 3. Were custody seals on coolers? NA Y N If yes, how many and where? 1 Front
 If present, were custody seals intact? Y N If present, were they signed and dated? Y N

| Raw Cooler Temp | Corrected Cooler Temp | Raw Temp Blank | Corrected Temp Blank | Corr. Factor | Thermometer ID | Cooler/COC ID | Tracking Number | NA | Filed |
|-----------------|-----------------------|----------------|----------------------|--------------|----------------|---------------|----------------------|----|-------|
| <u>0.4</u> | <u>0.7</u> | <u>0.0</u> | <u>0.3</u> | <u>0.3</u> | <u>360</u> | <u>NA</u> | <u>7744 92880491</u> | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |

4. Packing material: Inserts Baggies Bubble Wrap Gel Packs Wet Ice Dry Ice Sleeves
 5. Were custody papers properly filled out (ink, signed, etc.)? NA Y N
 6. Did all bottles arrive in good condition (unbroken)? *Indicate in the table below.* NA Y N
 7. Were all sample labels complete (i.e analysis, preservation, etc.)? NA Y N
 8. Did all sample labels and tags agree with custody papers? *Indicate major discrepancies in the table on page 2.* NA Y N
 9. Were appropriate bottles/containers and volumes received for the tests indicated? NA Y N
 10. Were the pH-preserved bottles (*see SMO GEN SOP*) received at the appropriate pH? *Indicate in the table below* NA Y N
 11. Were VOA vials received without headspace? *Indicate in the table below.* NA Y N
 12. Was C12/Res negative? NA Y N

| Sample ID on Bottle | Sample ID on COC | Identified by: |
|---------------------|------------------|----------------|
| | | |
| | | |
| | | |

| Sample ID | Bottle Count Bottle Type | Out of Temp | Head- space | Broke | pH | Reagent | Volume added | Reagent Lot Number | Initials | Time |
|-----------|-----------------------------|----------------|----------------|-------|----|---------|-----------------|-----------------------|----------|------|
| | | | | | | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |

Notes, Discrepancies, & Resolutions: _____



Steroids and Endocrine Disrupting Compounds

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ALS Group USA, Corp.
dba ALS Environmental

Analytical Report

Client: Eurofins Calscience Environmental Laboratory
Project: 15-09-0733
Sample Matrix: Water

Service Request: K1510108
Date Collected: 09/09/15 10:11
Date Received: 09/12/15 10:00

Sample Name: HCS110
Lab Code: K1510108-001

Units: ng/L
Basis: NA

Steroids and Endocrine Disrupting Compounds

Analysis Method: 1694
Prep Method: Method

| Analyte Name | Result | MRL | Dil. | Date Analyzed | Date Extracted | Q |
|--------------|--------|-----|------|----------------|----------------|---|
| Caffeine | 41 | 4.0 | 1 | 09/22/15 02:23 | 9/16/15 | |

| Surrogate Name | % Rec | Control Limits | Date Analyzed | Q |
|-------------------------|-------|----------------|----------------|---|
| Caffeine-trimethyl-13C3 | 76 | 46 - 161 | 09/22/15 02:23 | |

ALS Group USA, Corp.
dba ALS Environmental

Analytical Report

Client: Eurofins Calscience Environmental Laboratory
Project: 15-09-0733
Sample Matrix: Water

Service Request: K1510108
Date Collected: 09/09/15 10:44
Date Received: 09/12/15 10:00

Sample Name: HCS120
Lab Code: K1510108-002

Units: ng/L
Basis: NA

Steroids and Endocrine Disrupting Compounds

Analysis Method: 1694
Prep Method: Method

| Analyte Name | Result | MRL | Dil. | Date Analyzed | Date Extracted | Q |
|--------------|--------|-----|------|----------------|----------------|---|
| Caffeine | ND U | 4.0 | 1 | 09/22/15 02:44 | 9/16/15 | |

| Surrogate Name | % Rec | Control Limits | Date Analyzed | Q |
|-------------------------|-------|----------------|----------------|---|
| Caffeine-trimethyl-13C3 | 85 | 46 - 161 | 09/22/15 02:44 | |


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ALS Group USA, Corp.
dba ALS Environmental

Analytical Report

Client: Eurofins Calscience Environmental Laboratory
Project: 15-09-0733
Sample Matrix: Water
Sample Name: FDHCS120
Lab Code: K1510108-003

Service Request: K1510108
Date Collected: 09/09/15 10:44
Date Received: 09/12/15 10:00

Units: ng/L
Basis: NA

Steroids and Endocrine Disrupting Compounds

Analysis Method: 1694
Prep Method: Method

| Analyte Name | Result | MRL | Dil. | Date Analyzed | Date Extracted | Q |
|---------------------|---------------|------------|-------------|----------------------|-----------------------|----------|
| Caffeine | 16 | 4.0 | 1 | 09/22/15 03:05 | 9/16/15 | |

| Surrogate Name | % Rec | Control Limits | Date Analyzed | Q |
|-------------------------|--------------|-----------------------|----------------------|----------|
| Caffeine-trimethyl-13C3 | 96 | 46 - 161 | 09/22/15 03:05 | |


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ALS Group USA, Corp.
dba ALS Environmental

Analytical Report

Client: Eurofins Calscience Environmental Laboratory
Project: 15-09-0733
Sample Matrix: Water
Sample Name: HCS130
Lab Code: K1510108-004

Service Request: K1510108
Date Collected: 09/09/15 09:30
Date Received: 09/12/15 10:00

Units: ng/L
Basis: NA

Steroids and Endocrine Disrupting Compounds

Analysis Method: 1694
Prep Method: Method

| Analyte Name | Result | MRL | Dil. | Date Analyzed | Date Extracted | Q |
|--------------|--------|-----|------|----------------|----------------|---|
| Caffeine | ND U | 4.0 | 1 | 09/22/15 03:25 | 9/16/15 | |

| Surrogate Name | % Rec | Control Limits | Date Analyzed | Q |
|-------------------------|-------|----------------|----------------|---|
| Caffeine-trimethyl-13C3 | 92 | 46 - 161 | 09/22/15 03:25 | |


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ALS Group USA, Corp.
dba ALS Environmental

Analytical Report

Client: Eurofins Calscience Environmental Laboratory
Project: 15-09-0733
Sample Matrix: Water
Sample Name: HCS140
Lab Code: K1510108-005

Service Request: K1510108
Date Collected: 09/09/15 11:12
Date Received: 09/12/15 10:00

Units: ng/L
Basis: NA

Steroids and Endocrine Disrupting Compounds

Analysis Method: 1694
Prep Method: Method

| Analyte Name | Result | MRL | Dil. | Date Analyzed | Date Extracted | Q |
|--------------|--------|-----|------|----------------|----------------|---|
| Caffeine | ND U | 4.0 | 1 | 09/22/15 03:46 | 9/16/15 | |

| Surrogate Name | % Rec | Control Limits | Date Analyzed | Q |
|-------------------------|-------|----------------|----------------|---|
| Caffeine-trimethyl-13C3 | 115 | 46 - 161 | 09/22/15 03:46 | |


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ALS Group USA, Corp.
dba ALS Environmental

Analytical Report

Client: Eurofins Calscience Environmental Laboratory
Project: 15-09-0733
Sample Matrix: Water

Service Request: K1510108
Date Collected: 09/09/15 11:39
Date Received: 09/12/15 10:00

Sample Name: HCS160
Lab Code: K1510108-006

Units: ng/L
Basis: NA

Steroids and Endocrine Disrupting Compounds

Analysis Method: 1694
Prep Method: Method

| Analyte Name | Result | MRL | Dil. | Date Analyzed | Date Extracted | Q |
|--------------|--------|-----|------|----------------|----------------|---|
| Caffeine | ND U | 4.0 | 1 | 09/22/15 04:07 | 9/16/15 | |

| Surrogate Name | % Rec | Control Limits | Date Analyzed | Q |
|-------------------------|-------|----------------|----------------|---|
| Caffeine-trimethyl-13C3 | 94 | 46 - 161 | 09/22/15 04:07 | |

ALS Group USA, Corp.
dba ALS Environmental

Analytical Report

Client: Eurofins Calscience Environmental Laboratory
Project: 15-09-0733
Sample Matrix: Water

Service Request: K1510108**Date Collected:** NA**Date Received:** NA

Sample Name: Method Blank
Lab Code: KQ1510310-03

Units: ng/L**Basis:** NA**Steroids and Endocrine Disrupting Compounds**

Analysis Method: 1694
Prep Method: Method

| Analyte Name | Result | MRL | Dil. | Date Analyzed | Date Extracted | Q |
|--------------|--------|-----|------|----------------|----------------|---|
| Caffeine | ND U | 2.0 | 1 | 09/21/15 14:32 | 9/16/15 | |

| Surrogate Name | % Rec | Control Limits | Date Analyzed | Q |
|-------------------------|-------|----------------|----------------|---|
| Caffeine-trimethyl-13C3 | 93 | 46 - 161 | 09/21/15 14:32 | |


Return to Contents

QA/QC Report

Client: Eurofins Calscience Environmental Laboratory
Project: 15-09-0733
Sample Matrix: Water

Service Request: K1510108

SURROGATE RECOVERY SUMMARY
Steroids and Endocrine Disrupting Compounds

Analysis Method: 1694
Extraction Method: Method

| Caffeine-trimethyl-13C3 | | |
|------------------------------|--------------|----------|
| Sample Name | Lab Code | 46 - 161 |
| HCS110 | K1510108-001 | 76 |
| HCS120 | K1510108-002 | 85 |
| FDHCS120 | K1510108-003 | 96 |
| HCS130 | K1510108-004 | 92 |
| HCS140 | K1510108-005 | 115 |
| HCS160 | K1510108-006 | 94 |
| Lab Control Sample | KQ1510310-01 | 94 |
| Duplicate Lab Control Sample | KQ1510310-02 | 99 |
| Method Blank | KQ1510310-03 | 93 |


Return to Contents

ALS Group USA, Corp.
dba ALS Environmental

QA/QC Report

Client: Eurofins Calscience Environmental Laboratory
Project: 15-09-0733
Sample Matrix: Water

Service Request: K1510108
Date Analyzed: 09/21/15
Date Extracted: 09/16/15

Duplicate Lab Control Sample Summary
Steroids and Endocrine Disrupting Compounds

Analysis Method: 1694
Prep Method: Method

Units: ng/L
Basis: NA
Analysis Lot: 463091

Lab Control Sample
KQ1510310-01

Duplicate Lab Control Sample
KQ1510310-02

| Analyte Name | Result | Spike Amount | % Rec | Result | Spike Amount | % Rec | % Rec Limits | RPD | RPD Limit |
|---------------------|---------------|---------------------|--------------|---------------|---------------------|--------------|---------------------|------------|------------------|
| Caffeine | 99.9 | 100 | 100 | 97.1 | 100 | 97 | 50-140 | 3 | 30 |

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



September 22, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 110
Sample Number: AA97178
Sample Matrix: Non Potable Water

Collection Date/Time: 9/17/15 10:53
Receipt Date/Time: 9/17/15 15:27

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAP</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97178-A | E. coli | √ | SM 9223B-2004 | 45 | MPN/100 mL | | 1 | 44302 | 9/17/15 | 16:37 | HH/PAL |
| AA97178-A | E. Coli Holding Time - IDEXX Colilert | | NA | 5.73 | hours | | 0.00 | 44301 | 9/17/15 | 16:37 | HH/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44302

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

9/22/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



September 22, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
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Fax #: NA

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Sample Location: HSM 120
Sample Number: AA97179
Sample Matrix: Non Potable Water

Collection Date/Time: 9/17/15 11:30
Receipt Date/Time: 9/17/15 15:27

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAP</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97179-A | E. coli | √ | SM 9223B-2004 | 20 | MPN/100 mL | | 1 | 44302 | 9/17/15 | 16:37 | HH/PAL |
| AA97179-A | E. Coli Holding Time - IDEXX Colilert | | NA | 5.12 | hours | | 0.00 | 44301 | 9/17/15 | 16:37 | HH/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44302

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

9/22/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

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Sample Location: HSM 130
Sample Number: AA97180
Sample Matrix: Non Potable Water

Collection Date/Time: 9/17/15 11:57
Receipt Date/Time: 9/17/15 15:27

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAP</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97180-A | E. coli | √ | SM 9223B-2004 | 30 | MPN/100 mL | | 1 | 44302 | 9/17/15 | 16:37 | HH/PAL |
| AA97180-A | E. Coli Holding Time - IDEXX Colilert | | NA | 4.67 | hours | | 0.00 | 44301 | 9/17/15 | 16:37 | HH/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44302

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

9/22/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 140
Sample Number: AA97181
Sample Matrix: Non Potable Water

Collection Date/Time: 9/17/15 12:24
Receipt Date/Time: 9/17/15 15:27

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: No air space per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAP</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97181-A | E. coli | √ | SM 9223B-2004 | 37 | MPN/100 mL | | 1 | 44302 | 9/17/15 | 16:37 | HH/PAL |
| AA97181-A | E. Coli Holding Time - IDEXX Colilert | | NA | 4.22 | hours | | 0.00 | 44301 | 9/17/15 | 16:37 | HH/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44302

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

9/22/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 150
Sample Number: AA97182
Sample Matrix: Non Potable Water

Collection Date/Time: 9/17/15 12:51
Receipt Date/Time: 9/17/15 15:27

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAP</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97182-A | E. coli | √ | SM 9223B-2004 | 22 | MPN/100 mL | | 1 | 44302 | 9/17/15 | 16:37 | HH/PAL |
| AA97182-A | E. Coli Holding Time - IDEXX Colilert | | NA | 3.77 | hours | | 0.00 | 44301 | 9/17/15 | 16:37 | HH/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44302

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

9/22/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 160
Sample Number: AA97183
Sample Matrix: Non Potable Water

Collection Date/Time: 9/17/15 13:18
Receipt Date/Time: 9/17/15 15:27

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAP</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97183-A | E. coli | √ | SM 9223B-2004 | 28 | MPN/100 mL | | 1 | 44302 | 9/17/15 | 16:37 | HH/PAL |
| AA97183-A | E. Coli Holding Time - IDEXX Colilert | | NA | 3.32 | hours | | 0.00 | 44301 | 9/17/15 | 16:37 | HH/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44302

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

9/22/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



September 22, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 170
Sample Number: AA97184
Sample Matrix: Non Potable Water

Collection Date/Time: 9/17/15 13:47
Receipt Date/Time: 9/17/15 15:27

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAP</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97184-A | E. coli | √ | SM 9223B-2004 | 28 | MPN/100 mL | | 1 | 44302 | 9/17/15 | 16:37 | HH/PAL |
| AA97184-A | E. Coli Holding Time - IDEXX Colilert | | NA | 2.83 | hours | | 0.00 | 44301 | 9/17/15 | 16:37 | HH/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44302

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

9/22/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 1 of 2

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Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: FDHSM 110
Sample Number: AA97185
Sample Matrix: Non Potable Water

Collection Date/Time: 9/17/15 10:53
Receipt Date/Time: 9/17/15 15:27

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAP</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97185-A | E. coli | √ | SM 9223B-2004 | 75 | MPN/100 mL | | 1 | 44302 | 9/17/15 | 16:37 | HH/PAL |
| AA97185-A | E. Coli Holding Time - IDEXX Colilert | | NA | 5.73 | hours | | 0.00 | 44301 | 9/17/15 | 16:37 | HH/PAL |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



September 22, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44302

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

9/22/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable



WORK ORDER NUMBER: 15-09-1440

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: SWCA Environmental Consultants

Client Project Name: EAA 27122

Attention: Philip Pearce
6200 UTSA Blvd.
Suite 102
San Antonio, TX 78249-1618

A handwritten signature in black ink, appearing to read "Don Burley".

Approved for release on 10/20/2015 by:
Don Burley
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

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Work Order Number: 15-09-1440

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Work Order Narrative

Work Order: 15-09-1440

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Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 09/18/15. They were assigned to Work Order 15-09-1440.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.



Calscience

Sample Summary

| | |
|--|------------------------------------|
| Client: SWCA Environmental Consultants | Work Order: 15-09-1440 |
| 6200 UTSA Blvd., Suite 102 | Project Name: EAA 27122 |
| San Antonio, TX 78249-1618 | PO Number: 27122.02.06 |
| | Date/Time Received: 09/18/15 10:00 |
| | Number of Containers: 73 |

Attn: Philip Pearce

| Sample Identification | Lab Number | Collection Date and Time | Number of Containers | Matrix |
|-----------------------|--------------|--------------------------|----------------------|---------|
| HSM110 | 15-09-1440-1 | 09/17/15 10:53 | 9 | Aqueous |
| FDHSM110 | 15-09-1440-2 | 09/17/15 10:53 | 9 | Aqueous |
| HSM120 | 15-09-1440-3 | 09/17/15 11:30 | 9 | Aqueous |
| HSM130 | 15-09-1440-4 | 09/17/15 11:57 | 9 | Aqueous |
| HSM140 | 15-09-1440-5 | 09/17/15 12:24 | 9 | Aqueous |
| HSM150 | 15-09-1440-6 | 09/17/15 12:51 | 9 | Aqueous |
| HSM160 | 15-09-1440-7 | 09/17/15 13:18 | 9 | Aqueous |
| HSM170 | 15-09-1440-8 | 09/17/15 13:47 | 9 | Aqueous |
| TB12 | 15-09-1440-9 | 09/17/15 00:00 | 1 | Aqueous |

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Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-09-1440-1-K | 09/17/15 10:53 | Aqueous | IC 10 | N/A | 09/18/15 17:33 | 150918L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.28 | 0.10 | 0.025 | 1.00 | |
| Chloride | 29 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.17 | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | 0.56 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 36 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|----------|----------------|----------------|---------|-------|-----|----------------|-----------|
| FDHSM110 | 15-09-1440-2-K | 09/17/15 10:53 | Aqueous | IC 10 | N/A | 09/18/15 17:52 | 150918L01 |
|----------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.27 | 0.10 | 0.025 | 1.00 | |
| Chloride | 29 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.18 | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | 0.59 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 36 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|-------|-----|----------------|-----------|
| HSM120 | 15-09-1440-3-K | 09/17/15 11:30 | Aqueous | IC 10 | N/A | 09/18/15 18:11 | 150918L01 |
|--------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.21 | 0.10 | 0.025 | 1.00 | |
| Chloride | 19 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.11 | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | 1.3 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 28 | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-09-1440-4-K | 09/17/15 11:57 | Aqueous | IC 10 | N/A | 09/18/15 18:30 | 150918L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.24 | 0.10 | 0.025 | 1.00 | |
| Chloride | 23 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.10 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.7 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 32 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|--------|----------------|-------------------|---------|-------|-----|-------------------|-----------|
| HSM140 | 15-09-1440-5-K | 09/17/15 12:24 | Aqueous | IC 10 | N/A | 09/18/15 20:15 | 150918L01 |
|--------|----------------|-------------------|---------|-------|-----|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.27 | 0.10 | 0.025 | 1.00 | |
| Chloride | 19 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.086 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.3 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 28 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|--------|----------------|-------------------|---------|-------|-----|-------------------|-----------|
| HSM150 | 15-09-1440-6-K | 09/17/15 12:51 | Aqueous | IC 10 | N/A | 09/18/15 20:34 | 150918L01 |
|--------|----------------|-------------------|---------|-------|-----|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.23 | 0.10 | 0.025 | 1.00 | |
| Chloride | 19 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.069 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.3 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 28 | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-09-1440-7-K | 09/17/15 13:18 | Aqueous | IC 10 | N/A | 09/18/15 20:53 | 150918L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.22 | 0.10 | 0.025 | 1.00 | |
| Chloride | 19 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.098 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.3 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 28 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|-------|-----|----------------|-----------|
| HSM170 | 15-09-1440-8-K | 09/17/15 13:47 | Aqueous | IC 10 | N/A | 09/18/15 21:12 | 150918L01 |
|--------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.24 | 0.10 | 0.025 | 1.00 | |
| Chloride | 19 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.087 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.3 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 28 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|--------------|-----------------|-----|---------|-------|-----|----------------|-----------|
| Method Blank | 099-12-906-6100 | N/A | Aqueous | IC 10 | N/A | 09/18/15 14:41 | 150918L01 |
|--------------|-----------------|-----|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | ND | 0.10 | 0.025 | 1.00 | |
| Chloride | ND | 1.0 | 0.12 | 1.00 | |
| Bromide | ND | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | ND | 0.10 | 0.025 | 1.00 | |
| Sulfate | ND | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-09-1440-1-F | 09/17/15 10:53 | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 19:25 | 150922LA6F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 94.8 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 20.1 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 2.18 | 0.500 | 0.103 | 1.00 | |
| Sodium | 17.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.756 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.53 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|----------|----------------|----------------|---------|----------|----------|----------------|------------|
| FDHSM110 | 15-09-1440-2-F | 09/17/15 10:53 | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 19:28 | 150922LA6F |
|----------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 91.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 19.4 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 2.08 | 0.500 | 0.103 | 1.00 | |
| Sodium | 17.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.725 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.33 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|
| HSM120 | 15-09-1440-3-F | 09/17/15 11:30 | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 19:30 | 150922LA6F |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 95.1 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.1 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.54 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.576 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.37 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-09-1440-4-F | 09/17/15 11:57 | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 19:33 | 150922LA6F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 98.3 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 16.3 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.77 | 0.500 | 0.103 | 1.00 | |
| Sodium | 14.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.605 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.37 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|
| HSM140 | 15-09-1440-5-F | 09/17/15 12:24 | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 19:35 | 150922LA6F |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 94.9 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.0 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.52 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.584 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.19 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|
| HSM150 | 15-09-1440-6-F | 09/17/15 12:51 | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 19:38 | 150922LA6F |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 94.9 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 16.8 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.59 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.584 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.07 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-09-1440-7-F | 09/17/15 13:18 | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 19:40 | 150922LA6F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 94.9 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.2 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.55 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.6 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.588 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.29 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|
| HSM170 | 15-09-1440-8-F | 09/17/15 13:47 | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 19:43 | 150922LA6F |
|--------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 94.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.0 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.53 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.579 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.26 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|-----------------|-----|---------|----------|----------|----------------|------------|
| Method Blank | 099-15-683-1409 | N/A | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 18:51 | 150922LA6F |
|--------------|-----------------|-----|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|--------|---------|------|------------|
| Calcium | ND | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 0.00711 | 0.100 | 0.00336 | 1.00 | J |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-09-1440-1-F | 09/17/15 10:53 | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 16:09 | 150922LA3F |

Comment(s):
 - The reporting limit is elevated resulting from matrix interference.
 - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|---------|----------|------|------------|
| Antimony | ND | 0.00500 | 0.000498 | 5.00 | |
| Arsenic | ND | 0.00500 | 0.00193 | 5.00 | |
| Barium | 0.0407 | 0.00500 | 0.000493 | 5.00 | |
| Beryllium | ND | 0.00500 | 0.00145 | 5.00 | |
| Cadmium | ND | 0.00500 | 0.000642 | 5.00 | |
| Chromium | ND | 0.00500 | 0.00201 | 5.00 | |
| Copper | 0.00175 | 0.00500 | 0.000699 | 5.00 | J |
| Lead | ND | 0.00500 | 0.000449 | 5.00 | |
| Nickel | 0.00176 | 0.00500 | 0.000658 | 5.00 | J |
| Selenium | ND | 0.00500 | 0.000841 | 5.00 | |
| Silver | ND | 0.00500 | 0.000553 | 5.00 | |
| Thallium | ND | 0.00500 | 0.000504 | 5.00 | |
| Zinc | ND | 0.0250 | 0.00239 | 5.00 | |
| Aluminum | ND | 0.250 | 0.0165 | 5.00 | |
| Iron | ND | 0.250 | 0.0463 | 5.00 | |
| Manganese | 0.283 | 0.00500 | 0.000694 | 5.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM110 | 15-09-1440-2-F | 09/17/15 10:53 | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 16:12 | 150922LA3F |

Comment(s):
 - The reporting limit is elevated resulting from matrix interference.
 - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|---------|----------|------|------------|
| Antimony | ND | 0.00500 | 0.000498 | 5.00 | |
| Arsenic | ND | 0.00500 | 0.00193 | 5.00 | |
| Barium | 0.0394 | 0.00500 | 0.000493 | 5.00 | |
| Beryllium | ND | 0.00500 | 0.00145 | 5.00 | |
| Cadmium | ND | 0.00500 | 0.000642 | 5.00 | |
| Chromium | ND | 0.00500 | 0.00201 | 5.00 | |
| Copper | 0.00216 | 0.00500 | 0.000699 | 5.00 | J |
| Lead | ND | 0.00500 | 0.000449 | 5.00 | |
| Nickel | 0.00168 | 0.00500 | 0.000658 | 5.00 | J |
| Selenium | ND | 0.00500 | 0.000841 | 5.00 | |
| Silver | ND | 0.00500 | 0.000553 | 5.00 | |
| Thallium | ND | 0.00500 | 0.000504 | 5.00 | |
| Zinc | ND | 0.0250 | 0.00239 | 5.00 | |
| Aluminum | ND | 0.250 | 0.0165 | 5.00 | |
| Iron | ND | 0.250 | 0.0463 | 5.00 | |
| Manganese | 0.282 | 0.00500 | 0.000694 | 5.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM120 | 15-09-1440-3-F | 09/17/15 11:30 | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 16:16 | 150922LA3F |

Comment(s):
 - The reporting limit is elevated resulting from matrix interference.
 - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|---------|----------|------|------------|
| Antimony | ND | 0.00500 | 0.000498 | 5.00 | |
| Arsenic | ND | 0.00500 | 0.00193 | 5.00 | |
| Barium | 0.0377 | 0.00500 | 0.000493 | 5.00 | |
| Beryllium | ND | 0.00500 | 0.00145 | 5.00 | |
| Cadmium | ND | 0.00500 | 0.000642 | 5.00 | |
| Chromium | ND | 0.00500 | 0.00201 | 5.00 | |
| Copper | 0.00629 | 0.00500 | 0.000699 | 5.00 | |
| Lead | ND | 0.00500 | 0.000449 | 5.00 | |
| Nickel | 0.00180 | 0.00500 | 0.000658 | 5.00 | J |
| Selenium | ND | 0.00500 | 0.000841 | 5.00 | |
| Silver | ND | 0.00500 | 0.000553 | 5.00 | |
| Thallium | ND | 0.00500 | 0.000504 | 5.00 | |
| Zinc | ND | 0.0250 | 0.00239 | 5.00 | |
| Aluminum | ND | 0.250 | 0.0165 | 5.00 | |
| Iron | ND | 0.250 | 0.0463 | 5.00 | |
| Manganese | 0.00170 | 0.00500 | 0.000694 | 5.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 4 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-09-1440-4-F | 09/17/15 11:57 | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 16:19 | 150922LA3F |

Comment(s): - The reporting limit is elevated resulting from matrix interference.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|---------|----------|------|------------|
| Antimony | ND | 0.00500 | 0.000498 | 5.00 | |
| Arsenic | ND | 0.00500 | 0.00193 | 5.00 | |
| Barium | 0.0414 | 0.00500 | 0.000493 | 5.00 | |
| Beryllium | ND | 0.00500 | 0.00145 | 5.00 | |
| Cadmium | ND | 0.00500 | 0.000642 | 5.00 | |
| Chromium | ND | 0.00500 | 0.00201 | 5.00 | |
| Copper | 0.00204 | 0.00500 | 0.000699 | 5.00 | J |
| Lead | ND | 0.00500 | 0.000449 | 5.00 | |
| Nickel | 0.00162 | 0.00500 | 0.000658 | 5.00 | J |
| Selenium | ND | 0.00500 | 0.000841 | 5.00 | |
| Silver | ND | 0.00500 | 0.000553 | 5.00 | |
| Thallium | ND | 0.00500 | 0.000504 | 5.00 | |
| Zinc | ND | 0.0250 | 0.00239 | 5.00 | |
| Aluminum | ND | 0.250 | 0.0165 | 5.00 | |
| Iron | ND | 0.250 | 0.0463 | 5.00 | |
| Manganese | 0.00154 | 0.00500 | 0.000694 | 5.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-09-1440-5-F | 09/17/15 12:24 | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 16:23 | 150922LA3F |

Comment(s):
 - The reporting limit is elevated resulting from matrix interference.
 - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|---------|----------|------|------------|
| Antimony | ND | 0.00500 | 0.000498 | 5.00 | |
| Arsenic | ND | 0.00500 | 0.00193 | 5.00 | |
| Barium | 0.0371 | 0.00500 | 0.000493 | 5.00 | |
| Beryllium | ND | 0.00500 | 0.00145 | 5.00 | |
| Cadmium | ND | 0.00500 | 0.000642 | 5.00 | |
| Chromium | ND | 0.00500 | 0.00201 | 5.00 | |
| Copper | 0.00218 | 0.00500 | 0.000699 | 5.00 | J |
| Lead | ND | 0.00500 | 0.000449 | 5.00 | |
| Nickel | 0.00160 | 0.00500 | 0.000658 | 5.00 | J |
| Selenium | ND | 0.00500 | 0.000841 | 5.00 | |
| Silver | ND | 0.00500 | 0.000553 | 5.00 | |
| Thallium | ND | 0.00500 | 0.000504 | 5.00 | |
| Zinc | ND | 0.0250 | 0.00239 | 5.00 | |
| Aluminum | ND | 0.250 | 0.0165 | 5.00 | |
| Iron | ND | 0.250 | 0.0463 | 5.00 | |
| Manganese | 0.00148 | 0.00500 | 0.000694 | 5.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM150 | 15-09-1440-6-F | 09/17/15 12:51 | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 16:26 | 150922LA3F |

Comment(s):
 - The reporting limit is elevated resulting from matrix interference.
 - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|---------|----------|------|------------|
| Antimony | ND | 0.00500 | 0.000498 | 5.00 | |
| Arsenic | ND | 0.00500 | 0.00193 | 5.00 | |
| Barium | 0.0369 | 0.00500 | 0.000493 | 5.00 | |
| Beryllium | ND | 0.00500 | 0.00145 | 5.00 | |
| Cadmium | ND | 0.00500 | 0.000642 | 5.00 | |
| Chromium | ND | 0.00500 | 0.00201 | 5.00 | |
| Copper | 0.00165 | 0.00500 | 0.000699 | 5.00 | J |
| Lead | ND | 0.00500 | 0.000449 | 5.00 | |
| Nickel | 0.00157 | 0.00500 | 0.000658 | 5.00 | J |
| Selenium | ND | 0.00500 | 0.000841 | 5.00 | |
| Silver | ND | 0.00500 | 0.000553 | 5.00 | |
| Thallium | ND | 0.00500 | 0.000504 | 5.00 | |
| Zinc | ND | 0.0250 | 0.00239 | 5.00 | |
| Aluminum | ND | 0.250 | 0.0165 | 5.00 | |
| Iron | ND | 0.250 | 0.0463 | 5.00 | |
| Manganese | 0.00182 | 0.00500 | 0.000694 | 5.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-09-1440-7-F | 09/17/15 13:18 | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 16:30 | 150922LA3F |

Comment(s): - The reporting limit is elevated resulting from matrix interference.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|---------|----------|------|------------|
| Antimony | ND | 0.00500 | 0.000498 | 5.00 | |
| Arsenic | ND | 0.00500 | 0.00193 | 5.00 | |
| Barium | 0.0372 | 0.00500 | 0.000493 | 5.00 | |
| Beryllium | ND | 0.00500 | 0.00145 | 5.00 | |
| Cadmium | ND | 0.00500 | 0.000642 | 5.00 | |
| Chromium | ND | 0.00500 | 0.00201 | 5.00 | |
| Copper | 0.00691 | 0.00500 | 0.000699 | 5.00 | |
| Lead | ND | 0.00500 | 0.000449 | 5.00 | |
| Nickel | 0.00164 | 0.00500 | 0.000658 | 5.00 | J |
| Selenium | ND | 0.00500 | 0.000841 | 5.00 | |
| Silver | ND | 0.00500 | 0.000553 | 5.00 | |
| Thallium | ND | 0.00500 | 0.000504 | 5.00 | |
| Zinc | 0.0559 | 0.0250 | 0.00239 | 5.00 | |
| Aluminum | ND | 0.250 | 0.0165 | 5.00 | |
| Iron | ND | 0.250 | 0.0463 | 5.00 | |
| Manganese | 0.00320 | 0.00500 | 0.000694 | 5.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-09-1440-8-F | 09/17/15 13:47 | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 16:33 | 150922LA3F |

Comment(s):
- The reporting limit is elevated resulting from matrix interference.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|---------|----------|------|------------|
| Antimony | ND | 0.00500 | 0.000498 | 5.00 | |
| Arsenic | ND | 0.00500 | 0.00193 | 5.00 | |
| Barium | 0.0368 | 0.00500 | 0.000493 | 5.00 | |
| Beryllium | ND | 0.00500 | 0.00145 | 5.00 | |
| Cadmium | ND | 0.00500 | 0.000642 | 5.00 | |
| Chromium | ND | 0.00500 | 0.00201 | 5.00 | |
| Copper | 0.00856 | 0.00500 | 0.000699 | 5.00 | |
| Lead | ND | 0.00500 | 0.000449 | 5.00 | |
| Nickel | 0.00171 | 0.00500 | 0.000658 | 5.00 | J |
| Selenium | ND | 0.00500 | 0.000841 | 5.00 | |
| Silver | ND | 0.00500 | 0.000553 | 5.00 | |
| Thallium | ND | 0.00500 | 0.000504 | 5.00 | |
| Zinc | ND | 0.0250 | 0.00239 | 5.00 | |
| Aluminum | ND | 0.250 | 0.0165 | 5.00 | |
| Iron | ND | 0.250 | 0.0463 | 5.00 | |
| Manganese | 0.00325 | 0.00500 | 0.000694 | 5.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-914 | N/A | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 02:13 | 150922LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000138 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-09-1440-1-F | 09/17/15 10:53 | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 21:00 | 150923LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|----------|----------------|----------------|---------|------------|----------|----------------|------------|
| FDHSM110 | 15-09-1440-2-F | 09/17/15 10:53 | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 21:06 | 150923LA4F |
|----------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HSM120 | 15-09-1440-3-F | 09/17/15 11:30 | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 21:09 | 150923LA4F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HSM130 | 15-09-1440-4-F | 09/17/15 11:57 | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 21:11 | 150923LA4F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HSM140 | 15-09-1440-5-F | 09/17/15 12:24 | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 21:13 | 150923LA4F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HSM150 | 15-09-1440-6-F | 09/17/15 12:51 | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 21:15 | 150923LA4F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-09-1440-7-F | 09/17/15 13:18 | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 21:22 | 150923LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|
| HSM170 | 15-09-1440-8-F | 09/17/15 13:47 | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 21:24 | 150923LA4F |
|--------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|-----|---------|------------|----------|----------------|------------|
| Method Blank | 099-15-763-621 | N/A | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 20:55 | 150923LA4F |
|--------------|----------------|-----|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-09-1440-1-K | 09/17/15 10:53 | Aqueous | GC 44 | 09/21/15 | 09/26/15 09:22 | 150921L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 98 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 96 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM110 | 15-09-1440-2-K | 09/17/15 10:53 | Aqueous | GC 44 | 09/21/15 | 09/26/15 09:36 | 150921L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 98 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 96 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM120 | 15-09-1440-3-K | 09/17/15 11:30 | Aqueous | GC 44 | 09/21/15 | 09/26/15 09:50 | 150921L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 99 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 101 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-09-1440-4-K | 09/17/15 11:57 | Aqueous | GC 44 | 09/21/15 | 09/26/15 10:04 | 150921L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 102 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 102 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-09-1440-5-K | 09/17/15 12:24 | Aqueous | GC 44 | 09/21/15 | 09/26/15 10:19 | 150921L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 100 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 103 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM150 | 15-09-1440-6-K | 09/17/15 12:51 | Aqueous | GC 44 | 09/21/15 | 09/26/15 10:33 | 150921L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 101 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 101 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-09-1440-7-K | 09/17/15 13:18 | Aqueous | GC 44 | 09/21/15 | 09/26/15 10:47 | 150921L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 92 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 91 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-09-1440-8-K | 09/17/15 13:47 | Aqueous | GC 44 | 09/21/15 | 09/26/15 11:02 | 150921L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 100 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 94 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-843 | N/A | Aqueous | GC 44 | 09/21/15 | 09/26/15 09:07 | 150921L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 86 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 100 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-09-1440-1-K | 09/17/15 10:53 | Aqueous | GC 58 | 09/21/15 | 09/24/15 23:33 | 150921L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 97 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 91 | 50-135 | |

| FDHSM110 | 15-09-1440-2-K | 09/17/15 10:53 | Aqueous | GC 58 | 09/21/15 | 09/24/15 23:51 | 150921L12 |
|----------|----------------|----------------|---------|-------|----------|----------------|-----------|
|----------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 101 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 95 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM120 | 15-09-1440-3-K | 09/17/15 11:30 | Aqueous | GC 58 | 09/21/15 | 09/25/15 00:09 | 150921L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 99 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 96 | 50-135 | |

| HSM130 | 15-09-1440-4-K | 09/17/15 11:57 | Aqueous | GC 58 | 09/21/15 | 09/25/15 00:27 | 150921L12 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 101 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 95 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-09-1440-5-K | 09/17/15 12:24 | Aqueous | GC 58 | 09/21/15 | 09/25/15 00:45 | 150921L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 104 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 99 | 50-135 | |

| HSM150 | 15-09-1440-6-K | 09/17/15 12:51 | Aqueous | GC 58 | 09/21/15 | 09/25/15 01:02 | 150921L12 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 99 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 93 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-09-1440-7-K | 09/17/15 13:18 | Aqueous | GC 58 | 09/21/15 | 09/25/15 01:20 | 150921L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 96 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 91 | 50-135 | |

| | | | | | | | |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
| HSM170 | 15-09-1440-8-K | 09/17/15 13:47 | Aqueous | GC 58 | 09/21/15 | 09/25/15 01:38 | 150921L12 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 98 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 87 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-533-1091 | N/A | Aqueous | GC 58 | 09/21/15 | 09/24/15 23:15 | 150921L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 84 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 93 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 1 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-09-1440-1-K | 09/17/15 10:53 | Aqueous | GC 35 | 09/19/15 | 09/23/15 23:05 | 150919L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 96 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM110 | 15-09-1440-2-K | 09/17/15 10:53 | Aqueous | GC 35 | 09/19/15 | 09/23/15 23:51 | 150919L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 87 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM120 | 15-09-1440-3-K | 09/17/15 11:30 | Aqueous | GC 35 | 09/19/15 | 09/24/15 00:37 | 150919L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 98 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-09-1440-4-K | 09/17/15 11:57 | Aqueous | GC 35 | 09/19/15 | 09/24/15 01:23 | 150919L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 104 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-09-1440-5-K | 09/17/15 12:24 | Aqueous | GC 35 | 09/19/15 | 09/24/15 02:09 | 150919L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 109 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM150 | 15-09-1440-6-K | 09/17/15 12:51 | Aqueous | GC 35 | 09/19/15 | 09/24/15 02:55 | 150919L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 101 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-09-1440-7-K | 09/17/15 13:18 | Aqueous | GC 35 | 09/19/15 | 09/24/15 03:41 | 150919L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 105 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-09-1440-8-K | 09/17/15 13:47 | Aqueous | GC 35 | 09/19/15 | 09/24/15 04:26 | 150919L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 97 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-112 | N/A | Aqueous | GC 35 | 09/19/15 | 09/23/15 16:51 | 150919L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 92 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

Page 1 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-09-1440-1-K | 09/17/15 10:53 | Aqueous | GC 40 | 09/23/15 | 09/28/15 15:30 | 150923L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 91 | 0-123 | |

| FDHSM110 | 15-09-1440-2-K | 09/17/15 10:53 | Aqueous | GC 40 | 09/23/15 | 09/28/15 15:54 | 150923L05 |
|----------|----------------|----------------|---------|-------|----------|----------------|-----------|
|----------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 87 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM120 | 15-09-1440-3-K | 09/17/15 11:30 | Aqueous | GC 40 | 09/23/15 | 09/28/15 16:17 | 150923L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 78 | 0-123 | |

| HSM130 | 15-09-1440-4-K | 09/17/15 11:57 | Aqueous | GC 40 | 09/23/15 | 09/28/15 16:40 | 150923L05 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 90 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-09-1440-5-K | 09/17/15 12:24 | Aqueous | GC 40 | 09/23/15 | 09/28/15 17:03 | 150923L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 85 | 0-123 | |

| HSM150 | 15-09-1440-6-K | 09/17/15 12:51 | Aqueous | GC 40 | 09/23/15 | 09/28/15 17:26 | 150923L05 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 107 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-09-1440-7-K | 09/17/15 13:18 | Aqueous | GC 40 | 09/23/15 | 09/28/15 17:50 | 150923L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 90 | 0-123 | |

| HSM170 | 15-09-1440-8-K | 09/17/15 13:47 | Aqueous | GC 40 | 09/23/15 | 09/28/15 18:13 | 150923L05 |
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 101 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-034-660 | N/A | Aqueous | GC 40 | 09/23/15 | 09/25/15 13:49 | 150923L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 90 | 0-123 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-09-1440-1-I | 09/17/15 10:53 | Aqueous | GC/MS SS | 09/22/15 | 09/24/15 12:12 | 150922L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 77 | 33-120 | |
| 2-Fluorophenol | 51 | 24-120 | |
| Nitrobenzene-d5 | 86 | 38-120 | |
| p-Terphenyl-d14 | 77 | 41-137 | |
| Phenol-d6 | 33 | 16-120 | |
| 2,4,6-Tribromophenol | 78 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM110 | 15-09-1440-2-I | 09/17/15 10:53 | Aqueous | GC/MS SS | 09/22/15 | 09/24/15 12:32 | 150922L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 80 | 33-120 | |
| 2-Fluorophenol | 57 | 24-120 | |
| Nitrobenzene-d5 | 89 | 38-120 | |
| p-Terphenyl-d14 | 82 | 41-137 | |
| Phenol-d6 | 35 | 16-120 | |
| 2,4,6-Tribromophenol | 81 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM120 | 15-09-1440-3-I | 09/17/15 11:30 | Aqueous | GC/MS SS | 09/22/15 | 09/24/15 12:52 | 150922L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 81 | 33-120 | |
| 2-Fluorophenol | 55 | 24-120 | |
| Nitrobenzene-d5 | 86 | 38-120 | |
| p-Terphenyl-d14 | 78 | 41-137 | |
| Phenol-d6 | 35 | 16-120 | |
| 2,4,6-Tribromophenol | 77 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 10 of 27

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-09-1440-4-I | 09/17/15 11:57 | Aqueous | GC/MS SS | 09/22/15 | 09/24/15 13:11 | 150922L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 82 | 33-120 | |
| 2-Fluorophenol | 58 | 24-120 | |
| Nitrobenzene-d5 | 89 | 38-120 | |
| p-Terphenyl-d14 | 81 | 41-137 | |
| Phenol-d6 | 36 | 16-120 | |
| 2,4,6-Tribromophenol | 80 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-09-1440-5-I | 09/17/15 12:24 | Aqueous | GC/MS SS | 09/22/15 | 09/24/15 13:31 | 150922L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 71 | 33-120 | |
| 2-Fluorophenol | 53 | 24-120 | |
| Nitrobenzene-d5 | 76 | 38-120 | |
| p-Terphenyl-d14 | 71 | 41-137 | |
| Phenol-d6 | 34 | 16-120 | |
| 2,4,6-Tribromophenol | 68 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM150 | 15-09-1440-6-I | 09/17/15 12:51 | Aqueous | GC/MS SS | 09/22/15 | 09/24/15 13:51 | 150922L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 65 | 33-120 | |
| 2-Fluorophenol | 49 | 24-120 | |
| Nitrobenzene-d5 | 71 | 38-120 | |
| p-Terphenyl-d14 | 65 | 41-137 | |
| Phenol-d6 | 31 | 16-120 | |
| 2,4,6-Tribromophenol | 64 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-09-1440-7-I | 09/17/15 13:18 | Aqueous | GC/MS SS | 09/22/15 | 09/24/15 14:11 | 150922L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 60 | 33-120 | |
| 2-Fluorophenol | 45 | 24-120 | |
| Nitrobenzene-d5 | 62 | 38-120 | |
| p-Terphenyl-d14 | 61 | 41-137 | |
| Phenol-d6 | 29 | 16-120 | |
| 2,4,6-Tribromophenol | 56 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-09-1440-8-I | 09/17/15 13:47 | Aqueous | GC/MS SS | 09/22/15 | 09/24/15 14:30 | 150922L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 78 | 33-120 | |
| 2-Fluorophenol | 54 | 24-120 | |
| Nitrobenzene-d5 | 83 | 38-120 | |
| p-Terphenyl-d14 | 78 | 41-137 | |
| Phenol-d6 | 33 | 16-120 | |
| 2,4,6-Tribromophenol | 75 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-4108 | N/A | Aqueous | GC/MS SS | 09/22/15 | 09/24/15 11:08 | 150922L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 84 | 33-120 | |
| 2-Fluorophenol | 62 | 24-120 | |
| Nitrobenzene-d5 | 91 | 38-120 | |
| p-Terphenyl-d14 | 85 | 41-137 | |
| Phenol-d6 | 40 | 16-120 | |
| 2,4,6-Tribromophenol | 83 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 1 of 22

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM110 | 15-09-1440-1-A | 09/17/15 10:53 | Aqueous | GC/MS JJ | 09/22/15 | 09/22/15 17:06 | 150922L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 110 | 78-126 | |
| 1,2-Dichloroethane-d4 | 112 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM110 | 15-09-1440-2-A | 09/17/15 10:53 | Aqueous | GC/MS JJ | 09/22/15 | 09/22/15 17:34 | 150922L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 109 | 78-126 | |
| 1,2-Dichloroethane-d4 | 110 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 5 of 22

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM120 | 15-09-1440-3-A | 09/17/15 11:30 | Aqueous | GC/MS JJ | 09/22/15 | 09/22/15 18:02 | 150922L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 99 | 80-120 | |
| Dibromofluoromethane | 114 | 78-126 | |
| 1,2-Dichloroethane-d4 | 114 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 7 of 22

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM130 | 15-09-1440-4-A | 09/17/15 11:57 | Aqueous | GC/MS JJ | 09/22/15 | 09/22/15 18:30 | 150922L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 8 of 22

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 99 | 80-120 | |
| Dibromofluoromethane | 111 | 78-126 | |
| 1,2-Dichloroethane-d4 | 111 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 9 of 22

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM140 | 15-09-1440-5-D | 09/17/15 12:24 | Aqueous | GC/MS LL | 09/24/15 | 09/24/15 18:41 | 150924L008 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 99 | 78-126 | |
| 1,2-Dichloroethane-d4 | 94 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 11 of 22

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM150 | 15-09-1440-6-B | 09/17/15 12:51 | Aqueous | GC/MS LL | 09/24/15 | 09/24/15 16:19 | 150924L008 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 96 | 78-126 | |
| 1,2-Dichloroethane-d4 | 91 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM160 | 15-09-1440-7-B | 09/17/15 13:18 | Aqueous | GC/MS LL | 09/24/15 | 09/24/15 16:55 | 150924L008 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 92 | 80-120 | |
| Dibromofluoromethane | 98 | 78-126 | |
| 1,2-Dichloroethane-d4 | 91 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 15 of 22

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM170 | 15-09-1440-8-B | 09/17/15 13:47 | Aqueous | GC/MS LL | 09/24/15 | 09/24/15 17:30 | 150924L008 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 16 of 22

| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 99 | 78-126 | |
| 1,2-Dichloroethane-d4 | 93 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| TB12 | 15-09-1440-9-A | 09/17/15 00:00 | Aqueous | GC/MS JJ | 09/22/15 | 09/22/15 16:40 | 150922L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 108 | 78-126 | |
| 1,2-Dichloroethane-d4 | 110 | 75-135 | |
| Toluene-d8 | 103 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-18227 | N/A | Aqueous | GC/MS JJ | 09/22/15 | 09/22/15 11:22 | 150922L002 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 100 | 80-120 | |
| Dibromofluoromethane | 106 | 78-126 | |
| 1,2-Dichloroethane-d4 | 105 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 21 of 22

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-18261 | N/A | Aqueous | GC/MS LL | 09/24/15 | 09/24/15 13:57 | 150924L008 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | 0.50 | 10 | 0.41 | 1.00 | J |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 22 of 22

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 91 | 80-120 | |
| Dibromofluoromethane | 95 | 78-126 | |
| 1,2-Dichloroethane-d4 | 89 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

09/18/15

Work Order:

15-09-1440

Page 1 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HSM110 | 15-09-1440-1 | 09/17/15 10:53 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.038 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 10/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 266 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 266 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 420 | 1.00 | 0.870 | 1.00 | | mg/L | 09/24/15 | 09/24/15 | SM 2540 C |
| Solids, Total Suspended (24) | 4.5 | 1.0 | 0.83 | 1.00 | | mg/L | 09/22/15 | 09/22/15 | SM 2540 D |
| pH (24) | 7.03 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/18/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 09/26/15 | 09/26/15 | SM 4500 N Org B |

| FDHSM110 | 15-09-1440-2 | 09/17/15 10:53 | Aqueous |
|----------|--------------|----------------|---------|
|----------|--------------|----------------|---------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.039 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 10/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 271 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 271 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 400 | 1.00 | 0.870 | 1.00 | | mg/L | 09/24/15 | 09/24/15 | SM 2540 C |
| Solids, Total Suspended (24) | 4.6 | 1.0 | 0.83 | 1.00 | | mg/L | 09/22/15 | 09/22/15 | SM 2540 D |
| pH (24) | 7.10 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/18/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.35 | 0.50 | 0.28 | 1.00 | J | mg/L | 09/26/15 | 09/26/15 | SM 4500 N Org B |

| HSM120 | 15-09-1440-3 | 09/17/15 11:30 | Aqueous |
|--------|--------------|----------------|---------|
|--------|--------------|----------------|---------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.027 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 10/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 267 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 267 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 375 | 1.00 | 0.870 | 1.00 | | mg/L | 09/24/15 | 09/24/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 09/22/15 | 09/22/15 | SM 2540 D |
| pH (24) | 7.07 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/18/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.35 | 0.50 | 0.28 | 1.00 | J | mg/L | 09/26/15 | 09/26/15 | SM 4500 N Org B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

09/18/15

Work Order:

15-09-1440

Page 2 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HSM130 | 15-09-1440-4 | 09/17/15 11:57 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.029 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 10/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 268 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 268 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 335 | 1.00 | 0.870 | 1.00 | | mg/L | 09/24/15 | 09/24/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 09/22/15 | 09/22/15 | SM 2540 D |
| pH (24) | 7.15 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/18/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 09/26/15 | 09/26/15 | SM 4500 N Org B |

| HSM140 | 15-09-1440-5 | 09/17/15 12:24 | Aqueous |
|---------------|---------------------|-----------------------|----------------|
|---------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.027 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 10/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 268 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 268 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 285 | 1.00 | 0.870 | 1.00 | | mg/L | 09/24/15 | 09/24/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 09/22/15 | 09/22/15 | SM 2540 D |
| pH (24) | 7.32 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/18/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 09/26/15 | 09/26/15 | SM 4500 N Org B |

| HSM150 | 15-09-1440-6 | 09/17/15 12:51 | Aqueous |
|---------------|---------------------|-----------------------|----------------|
|---------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.027 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 10/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 265 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 265 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 325 | 1.00 | 0.870 | 1.00 | | mg/L | 09/24/15 | 09/24/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.8 | 1.0 | 0.83 | 1.00 | | mg/L | 09/22/15 | 09/22/15 | SM 2540 D |
| pH (24) | 7.37 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/18/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 09/26/15 | 09/26/15 | SM 4500 N Org B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

09/18/15

Work Order:

15-09-1440

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HSM160 | 15-09-1440-7 | 09/17/15 13:18 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.029 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 10/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 264 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 264 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 380 | 1.00 | 0.870 | 1.00 | | mg/L | 09/24/15 | 09/24/15 | SM 2540 C |
| Solids, Total Suspended (24) | 3.5 | 1.0 | 0.83 | 1.00 | | mg/L | 09/22/15 | 09/22/15 | SM 2540 D |
| pH (24) | 7.46 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/18/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 09/26/15 | 09/26/15 | SM 4500 N Org B |

| HSM170 | 15-09-1440-8 | 09/17/15 13:47 | Aqueous |
|---------------|---------------------|-----------------------|----------------|
|---------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.030 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 10/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 265 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 265 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 320 | 1.00 | 0.870 | 1.00 | | mg/L | 09/24/15 | 09/24/15 | SM 2540 C |
| Solids, Total Suspended (24) | 6.7 | 1.0 | 0.83 | 1.00 | | mg/L | 09/22/15 | 09/22/15 | SM 2540 D |
| pH (24) | 7.50 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 09/18/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 09/26/15 | 09/26/15 | SM 4500 N Org B |

| Method Blank | N/A | Aqueous |
|---------------------|------------|----------------|
|---------------------|------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|-------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 10/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 09/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.87 | 1.00 | | mg/L | 09/24/15 | 09/24/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 09/22/15 | 09/22/15 | SM 2540 D |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 09/26/15 | 09/26/15 | SM 4500 N Org B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-09-1412-1 | Sample | Aqueous | IC 10 | N/A | 09/18/15 15:59 | 150918S01 |
| 15-09-1412-1 | Matrix Spike | Aqueous | IC 10 | N/A | 09/18/15 22:08 | 150918S01 |
| 15-09-1412-1 | Matrix Spike Duplicate | Aqueous | IC 10 | N/A | 09/18/15 22:27 | 150918S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.1544 | 2.500 | 2.759 | 104 | 2.728 | 103 | 80-120 | 1 | 0-20 | |
| Chloride | 126.3 | 50.00 | 186.3 | 120 | 186.2 | 120 | 80-120 | 0 | 0-20 | |
| Bromide | 0.5047 | 5.000 | 5.369 | 97 | 5.368 | 97 | 80-120 | 0 | 0-20 | |
| Nitrate (as N) | 24.03 | 5.000 | 30.54 | 130 | 30.62 | 132 | 80-120 | 0 | 0-20 | 3 |
| Sulfate | 126.1 | 50.00 | 190.2 | 128 | 189.3 | 126 | 80-120 | 0 | 0-20 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|-----------|------------|---------------|----------------|---------------------|
| 15-09-1719-4 | Sample | Sea Water | ACA 1 | N/A | 10/02/15 15:04 | 151002S01 |
| 15-09-1719-4 | Matrix Spike | Sea Water | ACA 1 | N/A | 10/02/15 15:04 | 151002S01 |
| 15-09-1719-4 | Matrix Spike Duplicate | Sea Water | ACA 1 | N/A | 10/02/15 15:04 | 151002S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | 0.05681 | 0.2000 | 0.2857 | 114 | 0.2908 | 117 | 90-110 | 2 | 0-25 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM150 | Sample | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 19:38 | 150922SA6A |
| HSM150 | Matrix Spike | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 19:20 | 150922SA6A |
| HSM150 | Matrix Spike Duplicate | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 19:23 | 150922SA6A |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Calcium | 94.93 | 0.5000 | 94.91 | 4X | 95.08 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 16.80 | 0.5000 | 17.73 | 4X | 17.67 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 1.586 | 5.000 | 6.631 | 101 | 6.671 | 102 | 83-131 | 1 | 0-7 | |
| Sodium | 12.42 | 5.000 | 17.82 | 108 | 17.73 | 106 | 73-127 | 1 | 0-9 | |
| Strontium | 0.5843 | 0.5000 | 1.136 | 110 | 1.125 | 108 | 81-123 | 1 | 0-6 | |
| Silicon | 5.067 | 0.5000 | 5.757 | 4X | 5.736 | 4X | 24-180 | 4X | 0-15 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-09-1427-3 | Sample | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 02:52 | 150922SA3A |
| 15-09-1427-3 | Matrix Spike | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 02:20 | 150922SA3A |
| 15-09-1427-3 | Matrix Spike Duplicate | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 02:24 | 150922SA3A |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Antimony | ND | 0.1000 | 0.1072 | 107 | 0.1050 | 105 | 85-133 | 2 | 0-11 | |
| Arsenic | 0.07947 | 0.1000 | 0.1795 | 100 | 0.1773 | 98 | 73-127 | 1 | 0-11 | |
| Barium | 0.03030 | 0.1000 | 0.1396 | 109 | 0.1290 | 99 | 74-128 | 8 | 0-10 | |
| Beryllium | ND | 0.1000 | 0.1094 | 109 | 0.1063 | 106 | 56-122 | 3 | 0-11 | |
| Cadmium | ND | 0.1000 | 0.1009 | 101 | 0.09872 | 99 | 84-114 | 2 | 0-8 | |
| Chromium | 0.001546 | 0.1000 | 0.1120 | 110 | 0.1133 | 112 | 73-133 | 1 | 0-11 | |
| Copper | 0.001308 | 0.1000 | 0.1040 | 103 | 0.1031 | 102 | 72-108 | 1 | 0-10 | |
| Lead | ND | 0.1000 | 0.1129 | 113 | 0.1121 | 112 | 79-121 | 1 | 0-10 | |
| Nickel | 0.003385 | 0.1000 | 0.1042 | 101 | 0.1043 | 101 | 68-122 | 0 | 0-10 | |
| Selenium | ND | 0.1000 | 0.09469 | 95 | 0.09164 | 92 | 59-125 | 3 | 0-12 | |
| Silver | ND | 0.05000 | 0.05441 | 109 | 0.04931 | 99 | 68-128 | 10 | 0-14 | |
| Thallium | ND | 0.1000 | 0.1107 | 111 | 0.1098 | 110 | 73-121 | 1 | 0-11 | |
| Zinc | 0.02463 | 0.1000 | 0.1011 | 76 | 0.09564 | 71 | 43-145 | 6 | 0-39 | |
| Aluminum | ND | 0.1000 | 0.1271 | 127 | 0.1114 | 111 | 47-161 | 13 | 0-24 | |
| Iron | 12.65 | 5.100 | 17.65 | 98 | 18.04 | 106 | 27-201 | 2 | 0-24 | |
| Manganese | 1.733 | 0.1000 | 1.759 | 4X | 1.818 | 4X | 72-126 | 4X | 0-42 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM110 | Sample | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 21:00 | 150923SA4 |
| HSM110 | Matrix Spike | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 21:02 | 150923SA4 |
| HSM110 | Matrix Spike Duplicate | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 21:04 | 150923SA4 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | ND | 0.01000 | 0.009067 | 91 | 0.009003 | 90 | 55-133 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-09-1804-1 | Sample | Aqueous | GC/MS LL | 09/24/15 | 09/24/15 14:33 | 150924S006 | | | | |
| 15-09-1804-1 | Matrix Spike | Aqueous | GC/MS LL | 09/24/15 | 09/24/15 11:35 | 150924S006 | | | | |
| 15-09-1804-1 | Matrix Spike Duplicate | Aqueous | GC/MS LL | 09/24/15 | 09/24/15 12:11 | 150924S006 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 50.78 | 102 | 50.34 | 101 | 74-122 | 1 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 47.38 | 95 | 45.72 | 91 | 60-144 | 4 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 54.13 | 108 | 52.98 | 106 | 73-120 | 2 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 51.67 | 103 | 50.05 | 100 | 80-122 | 3 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 56.03 | 112 | 54.75 | 110 | 70-120 | 2 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 45.63 | 91 | 44.99 | 90 | 64-142 | 1 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 50.42 | 101 | 48.29 | 97 | 52-136 | 4 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 56.69 | 113 | 56.04 | 112 | 77-125 | 1 | 0-24 | |
| Toluene | ND | 50.00 | 54.72 | 109 | 53.59 | 107 | 72-126 | 2 | 0-23 | |
| Trichloroethene | ND | 50.00 | 52.40 | 105 | 51.58 | 103 | 74-128 | 2 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 54.92 | 110 | 54.81 | 110 | 67-133 | 0 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 109.6 | 110 | 107.9 | 108 | 63-129 | 2 | 0-25 | |
| o-Xylene | ND | 50.00 | 54.89 | 110 | 53.69 | 107 | 62-128 | 2 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 50.91 | 102 | 51.11 | 102 | 68-134 | 0 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|--------------|-------------|------------|----------------|----------------|-----------------------|
| 15-09-1427-3 | Sample | Aqueous | ICP/MS 03 | 09/22/15 00:00 | 09/24/15 02:52 | 150922SA3A |
| 15-09-1427-3 | PDS | Aqueous | ICP/MS 03 | 09/22/15 00:00 | 09/24/15 02:27 | 150922SA3A |
| Parameter | Sample Conc. | Spike Added | PDS Conc. | PDS %Rec. | %Rec. CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.1021 | 102 | 75-125 | |
| Arsenic | 0.07947 | 0.1000 | 0.1749 | 95 | 75-125 | |
| Barium | 0.03030 | 0.1000 | 0.1310 | 101 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.1016 | 102 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.09412 | 94 | 75-125 | |
| Chromium | 0.001546 | 0.1000 | 0.1063 | 105 | 75-125 | |
| Copper | 0.001308 | 0.1000 | 0.09770 | 96 | 75-125 | |
| Lead | ND | 0.1000 | 0.1064 | 106 | 75-125 | |
| Nickel | 0.003385 | 0.1000 | 0.09904 | 96 | 75-125 | |
| Selenium | ND | 0.1000 | 0.09106 | 91 | 75-125 | |
| Silver | ND | 0.05000 | 0.04634 | 93 | 75-125 | |
| Thallium | ND | 0.1000 | 0.1034 | 103 | 75-125 | |
| Zinc | 0.02463 | 0.1000 | 0.1130 | 88 | 75-125 | |
| Aluminum | ND | 0.1000 | 0.1086 | 109 | 75-125 | |
| Iron | 12.65 | 5.100 | 17.74 | 100 | 75-125 | |
| Manganese | 1.733 | 0.1000 | 1.799 | 4X | 75-125 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---|-------------------------|---------------------|------------------|---------------|-----------------------|------------------------|
| HSM110 | Sample | Aqueous | PH1/BUR03 | N/A | 09/29/15 15:23 | F0929ALKD1 |
| HSM110 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 09/29/15 15:23 | F0929ALKD1 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Alkalinity, Total (as CaCO ₃) | | 266.0 | 266.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|----------------|------------------|---------------|-----------------------|------------------------|
| HSM110 | Sample | Aqueous | PH1/BUR03 | N/A | 09/29/15 15:23 | F0929HCOD1 |
| HSM110 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 09/29/15 15:23 | F0929HCOD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Bicarbonate (as CaCO ₃) | 266.0 | 266.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|----------------|------------------|---------------|-----------------------|------------------------|
| HSM110 | Sample | Aqueous | PH1/BUR03 | N/A | 09/29/15 15:23 | F0929CO3D1 |
| HSM110 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 09/29/15 15:23 | F0929CO3D1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-09-1854-1 | Sample | Aqueous | SC 2 | 09/24/15 00:00 | 09/24/15 17:00 | F0924TDSD1 |
| 15-09-1854-1 | Sample Duplicate | Aqueous | SC 2 | 09/24/15 00:00 | 09/24/15 17:00 | F0924TDSD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Dissolved | 14220 | 14600 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-09-1314-2 | Sample | Aqueous | N/A | 09/22/15 00:00 | 09/22/15 20:00 | F0922TSSD4 |
| 15-09-1314-2 | Sample Duplicate | Aqueous | N/A | 09/22/15 00:00 | 09/22/15 20:00 | F0922TSSD4 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Suspended | 712.0 | 662.0 | 7 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-09-1456-1 | Sample | Aqueous | PH 1 | N/A | 09/18/15 20:03 | F0918PHD1 |
| 15-09-1456-1 | Sample Duplicate | Aqueous | PH 1 | N/A | 09/18/15 20:03 | F0918PHD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|------------------|---------------------|------------------|------------|---------------|-------------------|
| pH | 6.710 | 6.680 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: SM 4500 N Org B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-09-1885-1 | Sample | Aqueous | BUR05 | 09/26/15 00:00 | 09/26/15 17:39 | F0926TKND1 |
| 15-09-1885-1 | Sample Duplicate | Aqueous | BUR05 | 09/26/15 00:00 | 09/26/15 17:39 | F0926TKND1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Total Kjeldahl Nitrogen | 79.52 | 80.64 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|------------------|
| 099-12-906-6100 | LCS | Aqueous | IC 10 | N/A | 09/18/15 15:00 | 150918L01 |

| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
|----------------|-------------|-----------------|-----------|----------|------------|
| Fluoride | 2.500 | 2.425 | 97 | 90-110 | |
| Chloride | 50.00 | 51.63 | 103 | 90-110 | |
| Bromide | 5.000 | 5.085 | 102 | 90-110 | |
| Nitrate (as N) | 5.000 | 5.138 | 103 | 90-110 | |
| Sulfate | 50.00 | 51.15 | 102 | 90-110 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-739-118 | LCS | Aqueous | ACA 1 | N/A | 10/02/15 15:04 | 151002L01 | | | |
| 099-12-739-118 | LCSD | Aqueous | ACA 1 | N/A | 10/02/15 15:04 | 151002L01 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Phosphorus, Total | 0.2000 | 0.1936 | 97 | 0.1985 | 99 | 90-110 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-859-810 | LCS | Aqueous | PH1/BUR03 | N/A | 09/29/15 15:23 | F0929ALKB1 | | | |
| 099-15-859-810 | LCSD | Aqueous | PH1/BUR03 | N/A | 09/29/15 15:23 | F0929ALKB1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 100.0 | 98.00 | 98 | 98.00 | 98 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4755 | LCS | Aqueous | SC 2 | 09/24/15 | 09/24/15 17:00 | F0924TDSL1 | | | |
| 099-12-180-4755 | LCSD | Aqueous | SC 2 | 09/24/15 | 09/24/15 17:00 | F0924TDSL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 85.00 | 85 | 90.00 | 90 | 80-120 | 6 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-09-010-7327 | LCS | Aqueous | N/A | 09/22/15 | 09/22/15 20:00 | F0922TSSL4 | | | |
| 099-09-010-7327 | LCSD | Aqueous | N/A | 09/22/15 | 09/22/15 20:00 | F0922TSSL4 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Suspended | 100.0 | 107.0 | 107 | 109.0 | 109 | 80-120 | 2 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-15-683-1409 | LCS | Aqueous | ICP 7300 | 09/22/15 | 09/29/15 18:53 | 150922LA6F |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Calcium | | 0.5000 | 0.5027 | 101 | 80-120 | |
| Magnesium | | 0.5000 | 0.4298 | 86 | 80-120 | |
| Potassium | | 5.000 | 5.173 | 103 | 80-120 | |
| Sodium | | 5.000 | 5.153 | 103 | 80-120 | |
| Strontium | | 0.5000 | 0.5226 | 105 | 80-120 | |
| Silicon | | 0.5000 | 0.3795 | 76 | 80-120 | X |


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RPD: Relative Percent Difference. CL: Control Limits

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-914 | LCS | Aqueous | ICP/MS 03 | 09/22/15 | 09/24/15 02:17 | 150922LA3F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.1012 | 101 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.09927 | 99 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.09713 | 97 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.1078 | 108 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.1028 | 103 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.1022 | 102 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.09981 | 100 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.1008 | 101 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.09571 | 96 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.1032 | 103 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.04933 | 99 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.09890 | 99 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.1020 | 102 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.1038 | 104 | 80-120 | 73-127 | |
| Iron | 5.100 | 4.940 | 97 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.09943 | 99 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass



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Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|------------------|
| 099-15-763-621 | LCS | Aqueous | Mercury 04 | 09/23/15 | 09/23/15 20:58 | 150923LA4F |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.01000 | 0.008671 | 87 | 80-120 | |


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Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|--------|-----|--------|------------|
| 099-12-529-843 | LCS | Aqueous | GC 44 | 09/21/15 | 09/26/15 08:39 | 150921L11 | | | | |
| 099-12-529-843 | LCSD | Aqueous | GC 44 | 09/21/15 | 09/26/15 08:53 | 150921L11 | | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Alpha-BHC | 0.5000 | 0.4263 | 85 | 0.4453 | 89 | 50-135 | 36-149 | 4 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.4506 | 90 | 0.4694 | 94 | 50-135 | 36-149 | 4 | 0-25 | |
| Beta-BHC | 0.5000 | 0.4293 | 86 | 0.4537 | 91 | 50-135 | 36-149 | 6 | 0-25 | |
| Heptachlor | 0.5000 | 0.4298 | 86 | 0.4113 | 82 | 50-135 | 36-149 | 4 | 0-25 | |
| Delta-BHC | 0.5000 | 0.4937 | 99 | 0.5280 | 106 | 50-135 | 36-149 | 7 | 0-25 | |
| Aldrin | 0.5000 | 0.4004 | 80 | 0.3687 | 74 | 50-135 | 36-149 | 8 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.4287 | 86 | 0.4403 | 88 | 50-135 | 36-149 | 3 | 0-25 | |
| Endosulfan I | 0.5000 | 0.4440 | 89 | 0.4578 | 92 | 50-135 | 36-149 | 3 | 0-25 | |
| Dieldrin | 0.5000 | 0.4506 | 90 | 0.4658 | 93 | 50-135 | 36-149 | 3 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.4374 | 87 | 0.4433 | 89 | 50-135 | 36-149 | 1 | 0-25 | |
| Endrin | 0.5000 | 0.3748 | 75 | 0.3789 | 76 | 50-135 | 36-149 | 1 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.5100 | 102 | 0.5388 | 108 | 50-135 | 36-149 | 5 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.4376 | 88 | 0.4525 | 90 | 50-135 | 36-149 | 3 | 0-25 | |
| Endosulfan II | 0.5000 | 0.4390 | 88 | 0.4568 | 91 | 50-135 | 36-149 | 4 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.4495 | 90 | 0.4384 | 88 | 50-135 | 36-149 | 3 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.4345 | 87 | 0.4506 | 90 | 50-135 | 36-149 | 4 | 0-25 | |
| Methoxychlor | 0.5000 | 0.4271 | 85 | 0.4426 | 89 | 50-135 | 36-149 | 4 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-533-1091 | LCS | Aqueous | GC 58 | 09/21/15 | 09/24/15 22:38 | 150921L12 | | | |
| 099-12-533-1091 | LCSD | Aqueous | GC 58 | 09/21/15 | 09/24/15 22:57 | 150921L12 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Aroclor-1016 | 2.000 | 2.432 | 122 | 2.517 | 126 | 50-135 | 3 | 0-25 | |
| Aroclor-1260 | 2.000 | 2.486 | 124 | 2.585 | 129 | 50-135 | 4 | 0-25 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-963-112 | LCS | Aqueous | | GC 35 | 09/19/15 | 09/23/15 14:33 | 150919L09 | | | |
| 099-15-963-112 | LCSD | Aqueous | | GC 35 | 09/19/15 | 09/23/15 15:19 | 150919L09 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.04073 | 102 | 0.04139 | 103 | 30-130 | 13-147 | 2 | 0-30 | |
| Bolstar | 0.04000 | 0.03779 | 94 | 0.03829 | 96 | 30-130 | 13-147 | 1 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.03598 | 90 | 0.03645 | 91 | 30-130 | 13-147 | 1 | 0-30 | |
| Coumaphos | 0.04000 | 0.03619 | 90 | 0.03684 | 92 | 30-130 | 13-147 | 2 | 0-30 | |
| Diazinon | 0.04000 | 0.03756 | 94 | 0.03788 | 95 | 30-130 | 13-147 | 1 | 0-30 | |
| Disulfoton | 0.04000 | 0.03861 | 97 | 0.03886 | 97 | 30-130 | 13-147 | 1 | 0-30 | |
| Ethoprop | 0.04000 | 0.04308 | 108 | 0.04449 | 111 | 30-130 | 13-147 | 3 | 0-30 | |
| Fensulfothion | 0.04000 | 0.03812 | 95 | 0.03892 | 97 | 30-130 | 13-147 | 2 | 0-30 | |
| Fenthion | 0.04000 | 0.03851 | 96 | 0.03923 | 98 | 30-130 | 13-147 | 2 | 0-30 | |
| Merphos | 0.04000 | 0.06826 | 171 | 0.07033 | 176 | 30-130 | 13-147 | 3 | 0-30 | X |
| Methyl Parathion | 0.04000 | 0.03454 | 86 | 0.03737 | 93 | 30-130 | 13-147 | 8 | 0-30 | |
| Phorate | 0.04000 | 0.04601 | 115 | 0.04677 | 117 | 30-130 | 13-147 | 2 | 0-30 | |
| Ronnel | 0.04000 | 0.03550 | 89 | 0.03624 | 91 | 30-130 | 13-147 | 2 | 0-30 | |
| Stirophos | 0.04000 | 0.03865 | 97 | 0.03944 | 99 | 30-130 | 13-147 | 2 | 0-30 | |
| Tokuthion | 0.04000 | 0.03599 | 90 | 0.03640 | 91 | 30-130 | 13-147 | 1 | 0-30 | |
| Trichloronate | 0.04000 | 0.03823 | 96 | 0.03872 | 97 | 30-130 | 13-147 | 1 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-034-660 | LCS | Aqueous | GC 40 | 09/23/15 | 09/25/15 13:03 | 150923L05 | | | |
| 095-01-034-660 | LCSD | Aqueous | GC 40 | 09/23/15 | 09/25/15 13:26 | 150923L05 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| 2,4-D | 20.00 | 23.52 | 118 | 24.08 | 120 | 30-130 | 2 | 0-30 | |
| 2,4,5-T | 2.000 | 2.195 | 110 | 2.275 | 114 | 30-130 | 4 | 0-30 | |
| 2,4-DB | 20.00 | 21.98 | 110 | 22.76 | 114 | 30-130 | 3 | 0-30 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-4108 | LCS | Aqueous | | GC/MS SS | 09/22/15 | 09/24/15 10:28 | 150922L04 | | | |
| 095-01-003-4108 | LCSD | Aqueous | | GC/MS SS | 09/22/15 | 09/24/15 10:48 | 150922L04 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 200.0 | 177.4 | 89 | 178.3 | 89 | 61-120 | 51-130 | 1 | 0-20 | |
| Acenaphthylene | 200.0 | 176.7 | 88 | 177.8 | 89 | 55-120 | 44-131 | 1 | 0-20 | |
| Butyl Benzyl Phthalate | 200.0 | 186.0 | 93 | 187.4 | 94 | 56-122 | 45-133 | 1 | 0-20 | |
| 4-Chloro-3-Methylphenol | 200.0 | 197.2 | 99 | 195.0 | 97 | 52-120 | 41-131 | 1 | 0-20 | |
| 2-Chlorophenol | 200.0 | 194.9 | 97 | 195.3 | 98 | 47-120 | 35-132 | 0 | 0-20 | |
| 1,4-Dichlorobenzene | 200.0 | 164.0 | 82 | 166.8 | 83 | 36-120 | 22-134 | 2 | 0-20 | |
| Dimethyl Phthalate | 200.0 | 184.5 | 92 | 184.7 | 92 | 60-120 | 50-130 | 0 | 0-20 | |
| 2,4-Dinitrotoluene | 200.0 | 209.3 | 105 | 208.9 | 104 | 61-121 | 51-131 | 0 | 0-20 | |
| Fluorene | 200.0 | 179.9 | 90 | 179.8 | 90 | 67-120 | 58-129 | 0 | 0-20 | |
| N-Nitroso-di-n-propylamine | 200.0 | 184.0 | 92 | 178.7 | 89 | 39-123 | 25-137 | 3 | 0-20 | |
| Naphthalene | 200.0 | 167.7 | 84 | 168.1 | 84 | 54-120 | 43-131 | 0 | 0-20 | |
| 4-Nitrophenol | 200.0 | 95.21 | 48 | 93.28 | 47 | 14-120 | 0-138 | 2 | 0-20 | |
| Pentachlorophenol | 200.0 | 187.7 | 94 | 182.2 | 91 | 31-127 | 15-143 | 3 | 0-20 | |
| Phenol | 200.0 | 94.78 | 47 | 92.49 | 46 | 17-120 | 0-137 | 2 | 0-20 | |
| Pyrene | 200.0 | 180.9 | 90 | 181.4 | 91 | 58-124 | 47-135 | 0 | 0-20 | |
| 1,2,4-Trichlorobenzene | 200.0 | 168.0 | 84 | 170.8 | 85 | 49-120 | 37-132 | 2 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|-----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-14-001-18227 | LCS | Aqueous | | GC/MS JJ | 09/22/15 | 09/22/15 09:26 | 150922L002 | | | |
| 099-14-001-18227 | LCSD | Aqueous | | GC/MS JJ | 09/22/15 | 09/22/15 09:54 | 150922L002 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Benzene | 50.00 | 52.15 | 104 | 49.24 | 98 | 80-120 | 73-127 | 6 | 0-20 | |
| Carbon Tetrachloride | 50.00 | 55.86 | 112 | 51.11 | 102 | 67-139 | 55-151 | 9 | 0-20 | |
| Chlorobenzene | 50.00 | 51.12 | 102 | 49.64 | 99 | 78-120 | 71-127 | 3 | 0-20 | |
| 1,2-Dibromoethane | 50.00 | 49.28 | 99 | 50.60 | 101 | 80-120 | 73-127 | 3 | 0-20 | |
| 1,2-Dichlorobenzene | 50.00 | 50.68 | 101 | 49.33 | 99 | 63-129 | 52-140 | 3 | 0-20 | |
| 1,2-Dichloroethane | 50.00 | 54.76 | 110 | 54.49 | 109 | 70-130 | 60-140 | 0 | 0-20 | |
| 1,1-Dichloroethene | 50.00 | 50.86 | 102 | 46.28 | 93 | 66-126 | 56-136 | 9 | 0-20 | |
| Ethylbenzene | 50.00 | 52.85 | 106 | 50.23 | 100 | 80-123 | 73-130 | 5 | 0-20 | |
| Toluene | 50.00 | 54.66 | 109 | 51.45 | 103 | 80-120 | 73-127 | 6 | 0-20 | |
| Trichloroethene | 50.00 | 53.61 | 107 | 50.73 | 101 | 80-122 | 73-129 | 6 | 0-20 | |
| Vinyl Chloride | 50.00 | 41.24 | 82 | 36.88 | 74 | 70-130 | 60-140 | 11 | 0-20 | |
| p/m-Xylene | 100.0 | 108.0 | 108 | 102.1 | 102 | 75-123 | 67-131 | 6 | 0-20 | |
| o-Xylene | 50.00 | 53.51 | 107 | 51.25 | 102 | 74-122 | 66-130 | 4 | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 52.43 | 105 | 54.32 | 109 | 69-129 | 59-139 | 4 | 0-20 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 09/18/15
Work Order: 15-09-1440
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-18261 | LCS | Aqueous | GC/MS LL | 09/24/15 | 09/24/15 11:00 | 150924L008 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 46.16 | 92 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 41.87 | 84 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 49.23 | 98 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 48.25 | 96 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 50.45 | 101 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 43.32 | 87 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 43.81 | 88 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 51.36 | 103 | 80-123 | 73-130 | |
| Toluene | 50.00 | 49.48 | 99 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 46.13 | 92 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 46.70 | 93 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 99.79 | 100 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 49.95 | 100 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 48.41 | 97 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Sample Analysis Summary Report

Work Order: 15-09-1440

Page 1 of 1

| <u>Method</u> | <u>Extraction</u> | <u>Chemist ID</u> | <u>Instrument</u> | <u>Analytical Location</u> |
|-----------------|-------------------|-------------------|-------------------|----------------------------|
| EPA 300.0 | N/A | 1027 | IC 10 | 1 |
| EPA 365.1 | N/A | 650 | ACA 1 | 1 |
| EPA 6010B | EPA 3005A Filt. | 935 | ICP 7300 | 1 |
| EPA 6020 | EPA 3005A Filt. | 598 | ICP/MS 03 | 1 |
| EPA 7470A | EPA 7470A Filt. | 915 | Mercury 04 | 1 |
| EPA 8081A | EPA 3510C | 669 | GC 44 | 1 |
| EPA 8082 | EPA 3510C | 669 | GC 58 | 1 |
| EPA 8141A | EPA 3510C | 960 | GC 35 | 1 |
| EPA 8151A | EPA 8151A | 944 | GC 40 | 1 |
| EPA 8260B | EPA 5030C | 966 | GC/MS LL | 2 |
| EPA 8260B | EPA 5030C | 996 | GC/MS JJ | 2 |
| EPA 8270C | EPA 3510C | 923 | GC/MS SS | 1 |
| SM 2320B | N/A | 688 | PH1/BUR03 | 1 |
| SM 2540 C | N/A | 1009 | SC 2 | 1 |
| SM 2540 D | N/A | 689 | N/A | 1 |
| SM 4500 H+ B | N/A | 688 | PH 1 | 1 |
| SM 4500 N Org B | N/A | 685 | BUR05 | 1 |

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Location 1: 7440 Lincoln Way, Garden Grove, CA 92841

Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841

Glossary of Terms and Qualifiers

Work Order: 15-09-1440

Page 1 of 1

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|---|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| CI | See case narrative. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |
| | Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time. |
| | A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations. |

ORIGIN ID: SATA (210) 877-2847
SWCA INC

6200 UTSA BLVD STE 102

SAN ANTONIO, TX 782491618
UNITED STATES US

SHIP DATE: 17SEP15
ACTWGT: 55 LB
CAD: 7OFFC1601
DIMS: 24x13x14 IN
BILL SENDER

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1440

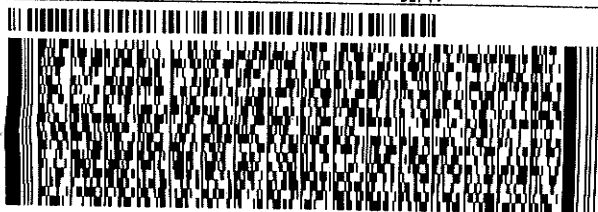
TO **DON BURLEY**
EUROFINS CALSCIENCE INC
7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 896-6994

REF:

DEPT:



FedEx
Express



J151215030301uy

5 of 8

MPS# 7813 5320 5609

0681

Mstr# 8079 4368 0339

0215

A7 APVA

FRI - 18 SEP AA
STANDARD OVERNIGHT

DSR

92841

CA-US SNA



4 of 8

MPS# 7813 5320 5594

0681

Mstr# 8079 4368 0339

0215

A7 APVA

FRI - 18 SEP AA
STANDARD OVERNIGHT

DSR

92841

CA-US SN



6 of 8

MPS# 7813 5320 5610

0681

Mstr# 8079 4368 0339

0215

A7 APVA

FRI - 18 SEP AA
STANDARD OVERNIGHT

DSR

92841

CA-US SNA



7 of 8

MPS# 7813 5320 5620

0681

Mstr# 8079 4368 0339

0215

A7 APVA

FRI - 18 SEP AA
STANDARD OVERNIGHT

DSR

92841

CA-US SNA



2 of 8

MPS# 7813 5320 5572

0681

Mstr# 8079 4368 0339

0215

A7 APVA

FRI - 18 SEP AA
STANDARD OVERNIGHT

DSR

92841

CA-US SNA



1 of 8
TRK# 8079 4368 0339
0215

MASTER

A7 APVA

FRI - 18 SEP AA
STANDARD OVERNIGHT

DSR

92841

CA-US SNA



3 of 8

MPS# 7813 5320 5583

0681

Mstr# 8079 4368 0339

0215

A7 APVA

FRI - 18 SEP AA
STANDARD OVERNIGHT

DSR

92841

CA-US SNA



8 of 8

MPS# 7813 5320 5631

0681

Mstr# 8079 4368 0339

0215

A7 APVA

FRI - 18 SEP AA
STANDARD OVERNIGHT

DSR

92841

CA-US SNA



SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 8

CLIENT: SWCA

DATE: 09 / 18 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF:-0.2°C); Temperature (w/o CF): 2.4 °C (w/ CF): 2.2 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A Checked by: 10/13

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☐ Yes ☒ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☐ No ☒ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☐ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: 5 (Trip Blank Lot Number: 150831A)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBh ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 g/L cube ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 10/13

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 6/29

SAMPLE RECEIPT CHECKLIST

COOLER 2 OF 8

CLIENT: SWCA

DATE: 09 / 18 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF:-0.2°C); Temperature (w/o CF): 2.7 °C (w/ CF): 2.5 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact

☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact

☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 10/1

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☐ No ☒ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☐ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: 5 (Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBz_{anna} ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal cuke ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) : ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 10/13

s = H₂SO₄, u = ultra-pure, z_{anna} = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 053

SAMPLE RECEIPT CHECKLIST

COOLER 3 OF 8

CLIENT: SWCA

DATE: 09 / 18 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF:-0.2°C); Temperature (w/o CF): 2.8 °C (w/ CF): 2.6 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact

☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact

☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 10/13

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE: 5 (Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal cuke ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 10/13

s = H₂SO₄, u = ultra-pure, z₂na = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 659

SAMPLE RECEIPT CHECKLIST

COOLER 4 OF 8

CLIENT: SWCA

DATE: 09 / 18 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF:-0.2°C); Temperature (w/o CF): 3.1 °C (w/ CF): 2.9 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A Checked by: 1013

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☐ No ☒ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☐ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: 5 (Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznnna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal cuke ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1013

s = H₂SO₄, u = ultra-pure, znnna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 619

SAMPLE RECEIPT CHECKLIST

COOLER 5 OF 8

CLIENT: SWCA

DATE: 09 / 18 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF: -0.2°C); Temperature (w/o CF): 3.3 °C (w/ CF): 3.1 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1013

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☐ No ☒ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☐ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: 5 (Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal Cuke ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1013s = H₂SO₄, u = ultra-pure, z₂na = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 659

SAMPLE RECEIPT CHECKLIST

COOLER 6 OF 8

CLIENT: SWCA

DATE: 09 / 18 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF: -0.2°C); Temperature (w/o CF): 2.5 °C (w/ CF): 2.3 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1013

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☐ No ☒ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☐ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: 5 (Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal Cuke ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1013

s = H₂SO₄, u = ultra-pure, z₂na = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 659

SAMPLE RECEIPT CHECKLIST

COOLER 7 OF 8

CLIENT: SWCA

DATE: 09 / 18 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF:-0.2°C); Temperature (w/o CF): 2.7 °C (w/ CF): 2.5 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 10/12

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☐ No ☒ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☐ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: 5 (Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna2 ☐ 100PJ ☐ 100PJna2 ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznnna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna2 ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal Cuke ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO3, na = NaOH, na2 = Na2S2O3, p = H3PO4, Labeled/Checked by: 10/12

s = H2SO4, u = ultra-pure, znnna = Zn(CH3CO2)2 + NaOH

Reviewed by: 6/9

SAMPLE RECEIPT CHECKLIST

COOLER 8 OF 8

CLIENT: SWCA

DATE: 09 / 18 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC5 (CF:-0.2°C); Temperature (w/o CF): 3.4 °C (w/ CF): 3.2 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 10/3

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☐ No ☒ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☐ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznn ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 10/3

s = H₂SO₄, u = ultra-pure, znn = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 6/9

SAMPLE ANOMALY REPORT

DATE: 09 / 18 / 2015

SAMPLES, CONTAINERS, AND LABELS:

- ☐ Sample(s) NOT RECEIVED but listed on COC
- ☐ Sample(s) received but NOT LISTED on COC
- ☐ Holding time expired (list client or ECI sample ID and analysis)
- ☐ Insufficient sample amount for requested analysis (list analysis)
- ☐ Improper container(s) used (list analysis)
- ☐ Improper preservative used (list analysis)
- ☐ No preservative noted on COC or label (list analysis and notify lab)
- ☐ Sample container(s) not labeled
- ☐ Client sample label(s) illegible (list container type and analysis)
- ☒ Client sample label(s) do not match COC (comment)
 - ☐ Project information
 - ☒ Client sample ID
 - ☐ Sampling date and/or time
 - ☒ Number of container(s)
 - ☐ Requested analysis
- ☐ Sample container(s) compromised (comment)
 - ☐ Broken
 - ☐ Water present in sample container
- ☐ Air sample container(s) compromised (comment)
 - ☐ Flat
 - ☐ Very low in volume
 - ☐ Leaking (not transferred; duplicate bag submitted)
 - ☐ Leaking (transferred into ECI Tedlar™ bags*)
 - ☐ Leaking (transferred into client's Tedlar™ bags*)

* Transferred at client's request.

MISCELLANEOUS: (Describe)

HEADSPACE:

(Containers with bubble > 6 mm or ¼ inch for volatile organic or dissolved gas analysis)

| ECI Sample ID | ECI Container ID | Total Number** | ECI Sample ID | ECI Container ID | Total Number** |
|---------------|------------------|----------------|---------------|------------------|----------------|
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

Comments

(-2) 25 Gallon cube labeled as
FDHSM 210

(-1) to (-8) Received 11 containers
instead of 9

Comments

(Containers with bubble for other analysis)

| ECI Sample ID | ECI Container ID | Total Number** | Requested Analysis |
|---------------|------------------|----------------|--------------------|
| | | | |
| | | | |
| | | | |
| | | | |

Comments: _____

** Record the total number of containers (i.e., vials or bottles) for the affected sample.

Reported by: 1013
Reviewed by: 649

Subcontractor Analysis Report

Work Order: 15-09-1440Page 1 of 1

One or more samples in this work order have tests that were subcontracted. The subcontract report(s) follows.

For subcontracted tests, please reference the laboratory information noted below.

1. McCampbell Analytical, Inc. - Pittsburg,CA CA ELAP 1644
TOC
2. ALS - Columbia Analytical Services, Inc. - Kelso,WA CA ELAP 2286, NELAP WA100010
Method 1694 Caffeine



McC Campbell Analytical, Inc.

"When Quality Counts"

Analytical Report

WorkOrder: 1509817

Report Created for: Eurofins Calscience, Inc.

7440 Lincoln Way
Garden Grove, CA 92841

Project Contact: Don Burley

Project P.O.:

Project Name: 15-09-1440

Project Received: 09/22/2015

Analytical Report reviewed & approved for release on 09/25/2015 by:

Angela Rydelius,
Laboratory Manager

The report shall not be reproduced except in full, without the written approval of the laboratory. The analytical results relate only to the items tested. Results reported conform to the most current NELAP standards, where applicable, unless otherwise stated in the case narrative.





Glossary of Terms & Qualifier Definitions

Client: Eurofins Calscience, Inc.

Project: 15-09-1440

WorkOrder: 1509817

Glossary Abbreviation

| | |
|--------------|--|
| 95% Interval | 95% Confident Interval |
| DF | Dilution Factor |
| DI WET | (DISTLC) Waste Extraction Test using DI water |
| DISS | Dissolved (direct analysis of 0.45 µm filtered and acidified water sample) |
| DUP | Duplicate |
| EDL | Estimated Detection Limit |
| ITEF | International Toxicity Equivalence Factor |
| LCS | Laboratory Control Sample |
| MB | Method Blank |
| MB % Rec | % Recovery of Surrogate in Method Blank, if applicable |
| MDL | Method Detection Limit |
| ML | Minimum Level of Quantitation |
| MS | Matrix Spike |
| MSD | Matrix Spike Duplicate |
| N/A | Not Applicable |
| ND | Not detected at or above the indicated MDL or RL |
| NR | Data Not Reported due to matrix interference or insufficient sample amount. |
| PF | Prep Factor |
| RD | Relative Difference |
| RL | Reporting Limit (The RL is the lowest calibration standard in a multipoint calibration.) |
| RPD | Relative Percent Deviation |
| RRT | Relative Retention Time |
| SPK Val | Spike Value |
| SPKRef Val | Spike Reference Value |
| SPLP | Synthetic Precipitation Leachate Procedure |
| TCLP | Toxicity Characteristic Leachate Procedure |
| TEQ | Toxicity Equivalents |
| WET (STLC) | Waste Extraction Test (Soluble Threshold Limit Concentration) |



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Analytical Report

Client: Eurofins Calscience, Inc.

WorkOrder: 1509817

Date Received: 9/22/15 12:55

Extraction Method: SM5310 B-2000

Date Prepared: 9/22/15

Analytical Method: SM5310 B-2000

Project: 15-09-1440

Unit: mg/L

Dissolved Organic Carbon (DOC)

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM110 | 1509817-001B | Water | 09/17/2015 10:53 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|--------------------------|--------|------|------|----|------------------|
| Dissolved Organic Carbon | 1.1 | 0.70 | 0.70 | 1 | 09/22/2015 20:02 |

Analyst(s): AV

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| FDHSM110 | 1509817-002B | Water | 09/17/2015 10:53 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|--------------------------|--------|------|------|----|------------------|
| Dissolved Organic Carbon | 1.2 | 0.70 | 0.70 | 1 | 09/22/2015 20:30 |

Analyst(s): AV

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM120 | 1509817-003B | Water | 09/17/2015 11:30 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|--------------------------|--------|------|------|----|------------------|
| Dissolved Organic Carbon | ND | 0.70 | 0.70 | 1 | 09/22/2015 22:51 |

Analyst(s): AV

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM130 | 1509817-004B | Water | 09/17/2015 11:57 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|--------------------------|--------|------|------|----|------------------|
| Dissolved Organic Carbon | ND | 0.70 | 0.70 | 1 | 09/22/2015 23:21 |

Analyst(s): AV

(Cont.)

CDPH ELAP 1644 ♦ NELAP 4033ORELAP

 Angela Rydelius, Lab Manager



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Analytical Report

Client: Eurofins Calscience, Inc.

WorkOrder: 1509817

Date Received: 9/22/15 12:55

Extraction Method: SM5310 B-2000

Date Prepared: 9/22/15

Analytical Method: SM5310 B-2000

Project: 15-09-1440

Unit: mg/L

Dissolved Organic Carbon (DOC)

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM140 | 1509817-005B | Water | 09/17/2015 12:24 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|--------------------------|--------|------|------|----|------------------|
| Dissolved Organic Carbon | ND | 0.70 | 0.70 | 1 | 09/23/2015 00:02 |

Analyst(s): AV

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM150 | 1509817-006B | Water | 09/17/2015 12:51 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|--------------------------|--------|------|------|----|------------------|
| Dissolved Organic Carbon | ND | 0.70 | 0.70 | 1 | 09/23/2015 00:31 |

Analyst(s): AV

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM160 | 1509817-007B | Water | 09/17/2015 13:18 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|--------------------------|--------|------|------|----|------------------|
| Dissolved Organic Carbon | ND | 0.70 | 0.70 | 1 | 09/23/2015 01:14 |

Analyst(s): AV

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM170 | 1509817-008B | Water | 09/17/2015 13:47 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|--------------------------|--------|------|------|----|------------------|
| Dissolved Organic Carbon | ND | 0.70 | 0.70 | 1 | 09/23/2015 03:20 |

Analyst(s): AV


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Analytical Report

Client: Eurofins Calscience, Inc.

WorkOrder: 1509817

Date Received: 9/22/15 12:55

Extraction Method: SM5310 Bm-2000

Date Prepared: 9/22/15

Analytical Method: SM5310 Bm-2000

Project: 15-09-1440

Unit: mg/L

Total Organic Carbon (TOC)

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM110 | 1509817-001A | Water | 09/17/2015 10:53 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|----------|--------|-------|------|----|------------------|
| TOC | 1.4 | 0.070 | 0.30 | 1 | 09/22/2015 20:17 |

Analyst(s): AV

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| FDHSM110 | 1509817-002A | Water | 09/17/2015 10:53 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|----------|--------|-------|------|----|------------------|
| TOC | 1.4 | 0.070 | 0.30 | 1 | 09/22/2015 20:43 |

Analyst(s): AV

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM120 | 1509817-003A | Water | 09/17/2015 11:30 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|----------|--------|-------|------|----|------------------|
| TOC | 0.86 | 0.070 | 0.30 | 1 | 09/22/2015 23:08 |

Analyst(s): AV

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM130 | 1509817-004A | Water | 09/17/2015 11:57 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|----------|--------|-------|------|----|------------------|
| TOC | 0.68 | 0.070 | 0.30 | 1 | 09/22/2015 23:35 |

Analyst(s): AV

(Cont.)


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Analytical Report

Client: Eurofins Calscience, Inc.

WorkOrder: 1509817

Date Received: 9/22/15 12:55

Extraction Method: SM5310 Bm-2000

Date Prepared: 9/22/15

Analytical Method: SM5310 Bm-2000

Project: 15-09-1440

Unit: mg/L

Total Organic Carbon (TOC)

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM140 | 1509817-005A | Water | 09/17/2015 12:24 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|----------|--------|-------|------|----|------------------|
| TOC | 0.81 | 0.070 | 0.30 | 1 | 09/23/2015 00:18 |

Analyst(s): AV

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM150 | 1509817-006A | Water | 09/17/2015 12:51 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|----------|--------|-------|------|----|------------------|
| TOC | 0.92 | 0.070 | 0.30 | 1 | 09/23/2015 00:46 |

Analyst(s): AV

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM160 | 1509817-007A | Water | 09/17/2015 13:18 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|----------|--------|-------|------|----|------------------|
| TOC | 0.93 | 0.070 | 0.30 | 1 | 09/23/2015 01:30 |

Analyst(s): AV

| Client ID | Lab ID | Matrix | Date Collected | Instrument | Batch ID |
|-----------|--------------|--------|------------------|------------|----------|
| HSM170 | 1509817-008A | Water | 09/17/2015 13:47 | WC_CNS | 110537 |

| Analytes | Result | MDL | RL | DF | Date Analyzed |
|----------|--------|-------|------|----|------------------|
| TOC | 1.2 | 0.070 | 0.30 | 1 | 09/23/2015 03:37 |

Analyst(s): AV



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Quality Control Report

Client: Eurofins Calscience, Inc.

WorkOrder: 1509817

Date Prepared: 9/22/15

BatchID: 110537

Date Analyzed: 9/22/15

Extraction Method: SM5310 B-2000

Instrument: WC_CNS

Analytical Method: SM5310 B-2000

Matrix: Water

Unit: mg/L

Project: 15-09-1440

Sample ID: MB/LCS-110537

1509726-007AMS/MSD

QC Summary Report for SM5310B

| Analyte | MB Result | LCS Result | MDL | RL | SPK Val | MB SS %REC | LCS %REC | LCS Limits |
|---------|-----------|------------|------|------|---------|------------|----------|------------|
| TOC | ND | 48.8 | 0.18 | 0.30 | 50 | - | 98 | 80-120 |

| Analyte | MS Result | MSD Result | SPK Val | SPKRef Val | MS %REC | MSD %REC | MS/MSD Limits | RPD | RPD Limit |
|---------|-----------|------------|---------|------------|---------|----------|---------------|------|-----------|
| TOC | 49.2 | 51.6 | 50 | ND | 97 | 102 | 70-130 | 4.68 | 20 |



CHAIN-OF-CUSTODY RECORD

WorkOrder: 1509817

ClientCode: CSEL

☐ WaterTrax☐ WriteOn☐ EDF☐ Excel☐ EQUIS☒ Email☐ HardCopy☐ ThirdParty☒ J-flag

Report to:

Don Burley
Eurofins Calscience, Inc.
7440 Lincoln Way
Garden Grove, CA 92841
(714) 895-5494 FAX: (714) 894-7501

Email: donaldburley@eurofinsUS.com

cc/3rd Party:

PO:

ProjectNo: 15-09-1440

Bill to:

Accounts Payable
Eurofins Calscience, Inc.
7440 Lincoln Way
Garden Grove, CA 92841

Requested TAT: 5 days;

Date Received: 09/22/2015

Date Printed: 09/22/2015

| Lab ID | Client ID | Matrix | Collection Date | Hold | Requested Tests (See legend below) | | | | | | | | | | | |
|-------------|-----------|--------|-----------------|--------------------------|------------------------------------|---|---|---|---|---|---|---|---|----|----|----|
| | | | | | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 |
| 1509817-001 | HSM110 | Water | 9/17/2015 10:53 | <input type="checkbox"/> | B | A | B | | | | | | | | | |
| 1509817-002 | FDHSM110 | Water | 9/17/2015 10:53 | <input type="checkbox"/> | B | A | B | | | | | | | | | |
| 1509817-003 | HSM120 | Water | 9/17/2015 11:30 | <input type="checkbox"/> | B | A | B | | | | | | | | | |
| 1509817-004 | HSM130 | Water | 9/17/2015 11:57 | <input type="checkbox"/> | B | A | B | | | | | | | | | |
| 1509817-005 | HSM140 | Water | 9/17/2015 12:24 | <input type="checkbox"/> | B | A | B | | | | | | | | | |
| 1509817-006 | HSM150 | Water | 9/17/2015 12:51 | <input type="checkbox"/> | B | A | B | | | | | | | | | |
| 1509817-007 | HSM160 | Water | 9/17/2015 13:18 | <input type="checkbox"/> | B | A | B | | | | | | | | | |
| 1509817-008 | HSM170 | Water | 9/17/2015 13:47 | <input type="checkbox"/> | B | A | B | | | | | | | | | |

Test Legend:

| | |
|---|----------------|
| 1 | cnsDOC_5310B_W |
| 5 | |
| 9 | |

| | |
|----|----------------|
| 2 | cnsTOC_5310B_W |
| 6 | |
| 10 | |

| | |
|----|-------------|
| 3 | PRDISSOLVED |
| 7 | |
| 11 | |

| | |
|----|--|
| 4 | |
| 8 | |
| 12 | |

Prepared by: Maria Venegas

Comments:

NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days).
Hazardous samples will be returned to client or disposed of at client expense.



WORK ORDER SUMMARY

Client Name: EUROFINS CALSCIENCE, INC.

QC Level: LEVEL 2

Work Order: 1509817

Project: 15-09-1440

Client Contact: Don Burley

Date Received: 9/22/2015

Comments:

Contact's Email: donaldburley@eurofinsUS.com

☐ WaterTrax ☐ WriteOn ☐ EDF ☐ Excel ☐ Fax ☒ Email ☐ HardCopy ☐ ThirdParty ☒ J-flag

| Lab ID | Client ID | Matrix | Test Name | Containers /Composites | Bottle & Preservative | De- chlorinated | Collection Date & Time | TAT | Sediment Content | Hold | SubOut |
|--------------|-----------|--------|----------------|---------------------------|-----------------------|--------------------------|---------------------------|--------|---------------------|--------------------------|--------|
| 1509817-001A | HSM110 | Water | SM5310Bm (TOC) | 1 | 250mL G w/ H2SO4 | <input type="checkbox"/> | 9/17/2015 10:53 | 5 days | Present | <input type="checkbox"/> | |
| 1509817-001B | HSM110 | Water | SM5310B (DOC) | 1 | 250mL G | <input type="checkbox"/> | 9/17/2015 10:53 | 5 days | Present | <input type="checkbox"/> | |
| 1509817-002A | FDHSM110 | Water | SM5310Bm (TOC) | 1 | 250mL G w/ H2SO4 | <input type="checkbox"/> | 9/17/2015 10:53 | 5 days | Present | <input type="checkbox"/> | |
| 1509817-002B | FDHSM110 | Water | SM5310B (DOC) | 1 | 250mL G | <input type="checkbox"/> | 9/17/2015 10:53 | 5 days | Present | <input type="checkbox"/> | |
| 1509817-003A | HSM120 | Water | SM5310Bm (TOC) | 1 | 250mL G w/ H2SO4 | <input type="checkbox"/> | 9/17/2015 11:30 | 5 days | Trace | <input type="checkbox"/> | |
| 1509817-003B | HSM120 | Water | SM5310B (DOC) | 1 | 250mL G | <input type="checkbox"/> | 9/17/2015 11:30 | 5 days | Trace | <input type="checkbox"/> | |
| 1509817-004A | HSM130 | Water | SM5310Bm (TOC) | 1 | 250mL G w/ H2SO4 | <input type="checkbox"/> | 9/17/2015 11:57 | 5 days | None | <input type="checkbox"/> | |
| 1509817-004B | HSM130 | Water | SM5310B (DOC) | 1 | 250mL G | <input type="checkbox"/> | 9/17/2015 11:57 | 5 days | None | <input type="checkbox"/> | |
| 1509817-005A | HSM140 | Water | SM5310Bm (TOC) | 1 | 250mL G w/ H2SO4 | <input type="checkbox"/> | 9/17/2015 12:24 | 5 days | Present | <input type="checkbox"/> | |
| 1509817-005B | HSM140 | Water | SM5310B (DOC) | 1 | 250mL G | <input type="checkbox"/> | 9/17/2015 12:24 | 5 days | Present | <input type="checkbox"/> | |
| 1509817-006A | HSM150 | Water | SM5310Bm (TOC) | 1 | 250mL G w/ H2SO4 | <input type="checkbox"/> | 9/17/2015 12:51 | 5 days | Present | <input type="checkbox"/> | |
| 1509817-006B | HSM150 | Water | SM5310B (DOC) | 1 | 250mL G | <input type="checkbox"/> | 9/17/2015 12:51 | 5 days | Present | <input type="checkbox"/> | |
| 1509817-007A | HSM160 | Water | SM5310Bm (TOC) | 1 | 250mL G w/ H2SO4 | <input type="checkbox"/> | 9/17/2015 13:18 | 5 days | Present | <input type="checkbox"/> | |
| 1509817-007B | HSM160 | Water | SM5310B (DOC) | 1 | 250mL G | <input type="checkbox"/> | 9/17/2015 13:18 | 5 days | Present | <input type="checkbox"/> | |
| 1509817-008A | HSM170 | Water | SM5310Bm (TOC) | 1 | 250mL G w/ H2SO4 | <input type="checkbox"/> | 9/17/2015 13:47 | 5 days | Present | <input type="checkbox"/> | |
| 1509817-008B | HSM170 | Water | SM5310B (DOC) | 1 | 250mL G | <input type="checkbox"/> | 9/17/2015 13:47 | 5 days | Present | <input type="checkbox"/> | |

NOTES: - STLC and TCLP extractions require 2 days to complete; therefore, all TATs begin after the extraction is completed (i.e., One-day TAT yields results in 3 days from sample submission).

- MAI assumes that all material present in the provided sampling container is considered part of the sample - MAI does not exclude any material from the sample prior to sample preparation unless requested in writing by the client.

To: McCampbell

DATE: 09/21/15

PAGE: 1 OF 1

7440 Lincoln Way, Garden Grove, CA 92841-1427 • (714) 895-5494

For courier service / sample drop off information, contact us26_sales@eurofinsus.com or call us.

[illegible]


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"When Quality Counts"

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 http://www.mcccampbell.com / E-mail: main@mcccampbell.com

Sample Receipt Checklist

 Client Name: **Eurofins Calscience, Inc.**

 Date and Time Received: **9/22/2015 12:55:08 PM**

 Project Name: **15-09-1440**

 LogIn Reviewed by: **Maria Venegas**

 WorkOrder №: **1509817**

 Matrix: Water

 Carrier: Golden State Overnight

Chain of Custody (COC) Information

| | | |
|---|---|-----------------------------|
| Chain of custody present? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> |
| Chain of custody signed when relinquished and received? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> |
| Chain of custody agrees with sample labels? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> |
| Sample IDs noted by Client on COC? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> |
| Date and Time of collection noted by Client on COC? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> |
| Sampler's name noted on COC? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> |

Sample Receipt Information

| | | | |
|--|---|-----------------------------|--|
| Custody seals intact on shipping container/cooler? | Yes <input type="checkbox"/> | No <input type="checkbox"/> | NA <input checked="" type="checkbox"/> |
| Shipping container/cooler in good condition? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| Samples in proper containers/bottles? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| Sample containers intact? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| Sufficient sample volume for indicated test? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |

Sample Preservation and Hold Time (HT) Information

| | | | |
|---|---|-----------------------------|--|
| All samples received within holding time? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| Sample/Temp Blank temperature | Temp: 1.2°C | | NA <input type="checkbox"/> |
| Water - VOA vials have zero headspace / no bubbles? | Yes <input type="checkbox"/> | No <input type="checkbox"/> | NA <input checked="" type="checkbox"/> |
| Sample labels checked for correct preservation? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| pH acceptable upon receipt (Metal: <2; 522: <4; 218.7: >8)? | Yes <input type="checkbox"/> | No <input type="checkbox"/> | NA <input checked="" type="checkbox"/> |
| Samples Received on Ice? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |

(Ice Type: WET ICE)

UCMR3 Samples:

| | | | |
|--|------------------------------|-----------------------------|--|
| Total Chlorine tested and acceptable upon receipt for EPA 522? | Yes <input type="checkbox"/> | No <input type="checkbox"/> | NA <input checked="" type="checkbox"/> |
| Free Chlorine tested and acceptable upon receipt for EPA 218.7, 300.1, 537, 539? | Yes <input type="checkbox"/> | No <input type="checkbox"/> | NA <input checked="" type="checkbox"/> |

* NOTE: If the "No" box is checked, see comments below.

Comments:

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ALS Environmental
ALS Group USA, Corp
1317 South 13th Avenue
Kelso, WA 98626
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: +1 360 636 1068
www.alsglobal.com

October 19, 2015

Analytical Report for Service Request No: K1510500

Donald Burley
Calscience Environmental Laboratories, Incorporated
7440 Lincoln Way
Garden Grove, CA 92841

RE: 15-09-1440

Dear Donald,

Enclosed are the results of the sample(s) submitted to our laboratory September 22, 2015
For your reference, these analyses have been assigned our service request number **K1510500**.

Analyses were performed according to our laboratory's NELAP-approved quality assurance program. The test results meet requirements of the current NELAP standards, where applicable, and except as noted in the laboratory case narrative provided. For a specific list of NELAP-accredited analytes, refer to the certifications section at www.alsglobal.com. All results are intended to be considered in their entirety, and ALS Group USA Corp. dba ALS Environmental (ALS) is not responsible for use of less than the complete report. Results apply only to the items submitted to the laboratory for analysis and individual items (samples) analyzed, as listed in the report.

Please contact me if you have any questions. My extension is 3364. You may also contact me via email at howard.holmes@alsglobal.com.

Respectfully submitted,

ALS Group USA, Corp. dba ALS Environmental

Howard Holmes
Project Manager



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Acronyms

Qualifiers

State Certifications, Accreditations, And Licenses

Chain of Custody

Steroids and Endocrine Disrupting Compounds

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Acronyms

| | |
|------------|--|
| ASTM | American Society for Testing and Materials |
| A2LA | American Association for Laboratory Accreditation |
| CARB | California Air Resources Board |
| CAS Number | Chemical Abstract Service registry Number |
| CFC | Chlorofluorocarbon |
| CFU | Colony-Forming Unit |
| DEC | Department of Environmental Conservation |
| DEQ | Department of Environmental Quality |
| DHS | Department of Health Services |
| DOE | Department of Ecology |
| DOH | Department of Health |
| EPA | U. S. Environmental Protection Agency |
| ELAP | Environmental Laboratory Accreditation Program |
| GC | Gas Chromatography |
| GC/MS | Gas Chromatography/Mass Spectrometry |
| LOD | Limit of Detection |
| LOQ | Limit of Quantitation |
| LUFT | Leaking Underground Fuel Tank |
| M | Modified |
| MCL | Maximum Contaminant Level is the highest permissible concentration of a substance allowed in drinking water as established by the USEPA. |
| MDL | Method Detection Limit |
| MPN | Most Probable Number |
| MRL | Method Reporting Limit |
| NA | Not Applicable |
| NC | Not Calculated |
| NCASI | National Council of the Paper Industry for Air and Stream Improvement |
| ND | Not Detected |
| NIOSH | National Institute for Occupational Safety and Health |
| PQL | Practical Quantitation Limit |
| RCRA | Resource Conservation and Recovery Act |
| SIM | Selected Ion Monitoring |
| TPH | Total Petroleum Hydrocarbons |
| tr | Trace level is the concentration of an analyte that is less than the PQL but greater than or equal to the MDL. |

Inorganic Data Qualifiers

- * The result is an outlier. See case narrative.
- # The control limit criteria is not applicable. See case narrative.
- B The analyte was found in the associated method blank at a level that is significant relative to the sample result as defined by the DOD or NELAC standards.
- E The result is an estimate amount because the value exceeded the instrument calibration range.
- J The result is an estimated value.
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.
DOD-QSM 4.2 definition : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- i The MRL/MDL or LOQ/LOD is elevated due to a matrix interference.
- X See case narrative.
- Q See case narrative. One or more quality control criteria was outside the limits.
- H The holding time for this test is immediately following sample collection. The samples were analyzed as soon as possible after receipt by the laboratory.

Metals Data Qualifiers

- # The control limit criteria is not applicable. See case narrative.
- J The result is an estimated value.
- E The percent difference for the serial dilution was greater than 10%, indicating a possible matrix interference in the sample.
- M The duplicate injection precision was not met.
- N The Matrix Spike sample recovery is not within control limits. See case narrative.
- S The reported value was determined by the Method of Standard Additions (MSA).
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.
DOD-QSM 4.2 definition : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- W The post-digestion spike for furnace AA analysis is out of control limits, while sample absorbance is less than 50% of spike absorbance.
- i The MRL/MDL or LOQ/LOD is elevated due to a matrix interference.
- X See case narrative.
- + The correlation coefficient for the MSA is less than 0.995.
- Q See case narrative. One or more quality control criteria was outside the limits.

Organic Data Qualifiers

- * The result is an outlier. See case narrative.
- # The control limit criteria is not applicable. See case narrative.
- A A tentatively identified compound, a suspected aldol-condensation product.
- B The analyte was found in the associated method blank at a level that is significant relative to the sample result as defined by the DOD or NELAC standards.
- C The analyte was qualitatively confirmed using GC/MS techniques, pattern recognition, or by comparing to historical data.
- D The reported result is from a dilution.
- E The result is an estimated value.
- J The result is an estimated value.
- N The result is presumptive. The analyte was tentatively identified, but a confirmation analysis was not performed.
- P The GC or HPLC confirmation criteria was exceeded. The relative percent difference is greater than 40% between the two analytical results.
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.
DOD-QSM 4.2 definition : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- i The MRL/MDL or LOQ/LOD is elevated due to a chromatographic interference.
- X See case narrative.
- Q See case narrative. One or more quality control criteria was outside the limits.

Additional Petroleum Hydrocarbon Specific Qualifiers

- F The chromatographic fingerprint of the sample matches the elution pattern of the calibration standard.
- L The chromatographic fingerprint of the sample resembles a petroleum product, but the elution pattern indicates the presence of a greater amount of lighter molecular weight constituents than the calibration standard.
- H The chromatographic fingerprint of the sample resembles a petroleum product, but the elution pattern indicates the presence of a greater amount of heavier molecular weight constituents than the calibration standard.
- O The chromatographic fingerprint of the sample resembles an oil, but does not match the calibration standard.
- Y The chromatographic fingerprint of the sample resembles a petroleum product eluting in approximately the correct carbon range, but the elution pattern does not match the calibration standard.
- Z The chromatographic fingerprint does not resemble a petroleum product.

ALS Group USA Corp. dba ALS Environmental (ALS) - Kelso
State Certifications, Accreditations, and Licenses

| Agency | Web Site | Number |
|--------------------------|---|---------------|
| Alaska DEC UST | http://dec.alaska.gov/applications/eh/ehllabreports/USTLabs.aspx | UST-040 |
| Arizona DHS | http://www.azdhs.gov/lab/license/env.htm | AZ0339 |
| Arkansas - DEQ | http://www.adeq.state.ar.us/techsvs/labcert.htm | 88-0637 |
| California DHS (ELAP) | http://www.cdph.ca.gov/certlic/labs/Pages/ELAP.aspx | 2795 |
| DOD ELAP | http://www.denix.osd.mil/edqw/Accreditation/AccreditedLabs.cfm | L14-51 |
| Florida DOH | http://www.doh.state.fl.us/lab/EnvLabCert/WaterCert.htm | E87412 |
| Hawaii DOH | Not available | - |
| Idaho DHW | http://www.healthandwelfare.idaho.gov/Health/Labs/CertificationDrinkingWaterLabs/tabid/1833/Default.aspx | - |
| ISO 17025 | http://www.pjllabs.com/ | L14-50 |
| Louisiana DEQ | http://www.deq.louisiana.gov/portal/DIVISIONS/PublicParticipationandPermitSupport/LouisianaLaboratoryAccreditationProgram.aspx | 03016 |
| Maine DHS | Not available | WA01276 |
| Michigan DEQ | http://www.michigan.gov/deq/0,1607,7-135-3307_4131_4156---,00.html | 9949 |
| Minnesota DOH | http://www.health.state.mn.us/accreditation | 053-999-457 |
| Montana DPHHS | http://www.dphhs.mt.gov/publichealth/ | CERT0047 |
| Nevada DEP | http://ndep.nv.gov/bsdwlabservice.htm | WA01276 |
| New Jersey DEP | http://www.nj.gov/dep/oqa/ | WA005 |
| North Carolina DWQ | http://www.dwqlab.org/ | 605 |
| Oklahoma DEQ | http://www.deq.state.ok.us/CSDnew/labcert.htm | 9801 |
| Oregon – DEQ (NELAP) | http://public.health.oregon.gov/LaboratoryServices/EnvironmentalLaboratoryAccreditation/Pages/index.aspx | WA100010 |
| South Carolina DHEC | http://www.scdhec.gov/environment/envserv/ | 61002 |
| Texas CEQ | http://www.tceq.texas.gov/field/qa/env_lab_accreditation.html | T104704427 |
| Washington DOE | http://www.ecy.wa.gov/programs/eap/labs/lab-accreditation.html | C544 |
| Wisconsin DNR | http://dnr.wi.gov/ | 998386840 |
| Wyoming (EPA Region 8) | http://www.epa.gov/region8/water/dwhome/wyomingdi.html | - |
| Kelso Laboratory Website | www.alsglobal.com | NA |

Analyses were performed according to our laboratory's NELAP-approved quality assurance program. A complete listing of specific NELAP-certified analytes, can be found in the certification section at www.ALSGlobal.com or at the accreditation bodies web site.

Please refer to the certification and/or accreditation body's web site if samples are submitted for compliance purposes. The states highlighted above, require the analysis be listed on the state certification if used for compliance purposes and if the method/analyte is offered by that state.



Chain of Custody


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ALS Environmental—Kelso Laboratory
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www.alsglobal.com



For courier service / sample drop off information, contact us26_sales@eurofinsus.com or call us.

K1510500

To: ALS - Kelso

CHAIN OF CUSTODY RECORD

DATE: 09/21/15

PAGE: 1 OF 1

[illegible]

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Cooler Receipt and Preservation Form

Client / Project: Eurohns Calscience, Inc Service Request K15 10500Received: 9/22/15 Opened: 9/22/15 By: KR Unloaded: 9/22/15 By: KE

1. Samples were received via? Mail Fed Ex UPS DHL PDX Courier Hand Delivered
2. Samples were received in: (circle) Cooler Box Envelope Other NA
3. Were custody seals on coolers? NA (Y) N If yes, how many and where? 1 F
- If present, were custody seals intact? (Y) N If present, were they signed and dated? (Y) N

| Raw Cooler Temp | Corrected Cooler Temp | Raw Temp Blank | Corrected Temp Blank | Corr. Factor | Thermometer ID | Cooler/COC ID | Tracking Number | NA | Filed |
|-----------------|-----------------------|----------------|----------------------|--------------|----------------|---------------|-----------------|----|-------|
| -0.5 | -0.2 | — | — | +0.3 | 361 | NA | 774561083017 | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |

4. Packing material: Inserts Baggies Bubble Wrap GelPacks (Wet Ice) Dry Ice Sleeves large plastic bag
5. Were custody papers properly filled out (ink, signed, etc.)? NA (Y) N
6. Did all bottles arrive in good condition (unbroken)? Indicate in the table below. NA (Y) N
7. Were all sample labels complete (i.e analysis, preservation, etc.)? NA (Y) N
8. Did all sample labels and tags agree with custody papers? Indicate major discrepancies in the table on page 2. NA (Y) N
9. Were appropriate bottles/containers and volumes received for the tests indicated? NA (Y) N
10. Were the pH-preserved bottles (see SMO GEN SOP) received at the appropriate pH? Indicate in the table below (NA) Y N
11. Were VOA vials received without headspace? Indicate in the table below. (NA) Y N
12. Was C12/Res negative? (NA) Y N

| Sample ID on Bottle | Sample ID on COC | Identified by: |
|---------------------|------------------|----------------|
| | | |
| | | |
| | | |

| Sample ID | Bottle Count Bottle Type | Out of Temp | Head- space | Broke | pH | Reagent | Volume added | Reagent Lot Number | Initials | Time |
|-----------|-----------------------------|----------------|----------------|-------|----|---------|-----------------|-----------------------|----------|------|
| | | | | | | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |

Notes, Discrepancies, & Resolutions: _____



Steroids and Endocrine Disrupting Compounds


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WORK ORDER NUMBER: 15-06-0463

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: SWCA Environmental Consultants

Client Project Name: EAA 27122

Attention: Philip Pearce
6200 UTSA Blvd.
Suite 102
San Antonio, TX 78249-1618

H. Burley FOR

Approved for release on 06/19/2015 by:
Don Burley
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

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Client Project Name: EAA 27122
Work Order Number: 15-06-0463

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Work Order Narrative

Work Order: 15-06-0463

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Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 06/05/15. They were assigned to Work Order 15-06-0463.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.



Calscience

Sample Summary

| | |
|--|------------------------------------|
| Client: SWCA Environmental Consultants | Work Order: 15-06-0463 |
| 6200 UTSA Blvd., Suite 102 | Project Name: EAA 27122 |
| San Antonio, TX 78249-1618 | PO Number: |
| | Date/Time Received: 06/05/15 09:40 |
| | Number of Containers: 13 |
| Attn: Philip Pearce | |

| Sample Identification | Lab Number | Collection Date and Time | Number of Containers | Matrix |
|-----------------------|--------------|--------------------------|----------------------|---------|
| HCS310 | 15-06-0463-1 | 06/04/15 11:22 | 2 | Solid |
| HCS320 | 15-06-0463-2 | 06/04/15 11:43 | 2 | Solid |
| HCS330 | 15-06-0463-3 | 06/04/15 12:20 | 2 | Solid |
| HCS340 | 15-06-0463-4 | 06/04/15 13:57 | 2 | Solid |
| HCS360 | 15-06-0463-5 | 06/04/15 14:24 | 2 | Solid |
| FDHCS360 | 15-06-0463-6 | 06/04/15 14:24 | 2 | Solid |
| TB07 | 15-06-0463-7 | 06/04/15 00:00 | 1 | Aqueous |

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Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: EPA 300.0
Units: mg/kg

Project: EAA 27122

Page 1 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS310 | 15-06-0463-1-B | 06/04/15 11:22 | Solid | IC 10 | 06/10/15 | 06/11/15 00:54 | 150610L01P |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Fluoride | ND | 1.2 | 1.00 | |
| Chloride | ND | 12 | 1.00 | |
| Bromide | ND | 1.2 | 1.00 | |
| Nitrate (as N) | ND | 1.2 | 1.00 | |
| Sulfate | 56 | 12 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|
| HCS320 | 15-06-0463-2-B | 06/04/15 11:43 | Solid | IC 10 | 06/10/15 | 06/11/15 01:10 | 150610L01P |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Fluoride | 1.6 | 1.4 | 1.00 | |
| Chloride | ND | 14 | 1.00 | |
| Bromide | ND | 1.4 | 1.00 | |
| Nitrate (as N) | 2.4 | 1.4 | 1.00 | |
| Sulfate | 26 | 14 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|
| HCS330 | 15-06-0463-3-B | 06/04/15 12:20 | Solid | IC 10 | 06/10/15 | 06/11/15 01:26 | 150610L01P |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Fluoride | 2.3 | 1.4 | 1.00 | |
| Chloride | ND | 14 | 1.00 | |
| Bromide | ND | 1.4 | 1.00 | |
| Nitrate (as N) | ND | 1.4 | 1.00 | |
| Sulfate | 34 | 14 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: EPA 300.0
Units: mg/kg

Project: EAA 27122

Page 2 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS340 | 15-06-0463-4-B | 06/04/15 13:57 | Solid | IC 10 | 06/10/15 | 06/11/15 01:43 | 150610L01P |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|----------------|--------|-----|------|------------|
| Fluoride | 2.7 | 2.3 | 1.00 | |
| Chloride | 33 | 23 | 1.00 | |
| Bromide | ND | 2.3 | 1.00 | |
| Nitrate (as N) | ND | 2.3 | 1.00 | |
| Sulfate | 63 | 23 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|
| HCS360 | 15-06-0463-5-B | 06/04/15 14:24 | Solid | IC 10 | 06/10/15 | 06/11/15 01:59 | 150610L01P |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|----------------|--------|-----|------|------------|
| Fluoride | 3.1 | 2.0 | 1.00 | |
| Chloride | 22 | 20 | 1.00 | |
| Bromide | ND | 2.0 | 1.00 | |
| Nitrate (as N) | ND | 2.0 | 1.00 | |
| Sulfate | 49 | 20 | 1.00 | |

| | | | | | | | |
|----------|----------------|----------------|-------|-------|----------|----------------|------------|
| FDHCS360 | 15-06-0463-6-B | 06/04/15 14:24 | Solid | IC 10 | 06/10/15 | 06/11/15 02:16 | 150610L01P |
|----------|----------------|----------------|-------|-------|----------|----------------|------------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|----------------|--------|-----|------|------------|
| Fluoride | 3.2 | 2.0 | 1.00 | |
| Chloride | 22 | 20 | 1.00 | |
| Bromide | ND | 2.0 | 1.00 | |
| Nitrate (as N) | ND | 2.0 | 1.00 | |
| Sulfate | 68 | 20 | 1.00 | |

| | | | | | | | |
|--------------|----------------|-----|-------|-------|----------|----------------|------------|
| Method Blank | 099-12-922-584 | N/A | Solid | IC 10 | 06/10/15 | 06/11/15 00:21 | 150610L01P |
|--------------|----------------|-----|-------|-------|----------|----------------|------------|

| Parameter | Result | RL | DF | Qualifiers |
|----------------|--------|-----|------|------------|
| Fluoride | ND | 1.0 | 1.00 | |
| Chloride | ND | 10 | 1.00 | |
| Bromide | ND | 1.0 | 1.00 | |
| Nitrate (as N) | ND | 1.0 | 1.00 | |
| Sulfate | ND | 10 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6010B
Units: mg/kg

Project: EAA 27122

Page 1 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS310 | 15-06-0463-1-AA | 06/04/15 11:22 | Solid | ICP 7300 | 06/10/15 | 06/11/15 16:45 | 150610L04 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|------|-------|------------|
| Magnesium | 1250 | 6.16 | 0.995 | |
| Potassium | 401 | 30.8 | 0.995 | |
| Sodium | 87.7 | 30.8 | 0.995 | |
| Strontium | 127 | 1.85 | 0.995 | |
| Silicon | 647 | 6.16 | 0.995 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|
| HCS310 | 15-06-0463-1-AA | 06/04/15 11:22 | Solid | ICP 7300 | 06/10/15 | 06/17/15 22:46 | 150610L04 |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|------|------|------------|
| Calcium | 249000 | 61.6 | 9.95 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|
| HCS320 | 15-06-0463-2-AA | 06/04/15 11:43 | Solid | ICP 7300 | 06/10/15 | 06/11/15 16:46 | 150610L04 |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|------|------|------------|
| Magnesium | 1900 | 7.16 | 1.02 | |
| Potassium | 456 | 35.8 | 1.02 | |
| Sodium | 353 | 35.8 | 1.02 | |
| Strontium | 235 | 2.15 | 1.02 | |
| Silicon | 803 | 7.16 | 1.02 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|
| HCS320 | 15-06-0463-2-AA | 06/04/15 11:43 | Solid | ICP 7300 | 06/10/15 | 06/17/15 22:48 | 150610L04 |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|------|------|------------|
| Calcium | 328000 | 71.6 | 10.2 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6010B
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS330 | 15-06-0463-3-AA | 06/04/15 12:20 | Solid | ICP 7300 | 06/10/15 | 06/11/15 16:48 | 150610L04 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|------|-------|------------|
| Magnesium | 3640 | 6.96 | 0.966 | |
| Potassium | 2410 | 34.8 | 0.966 | |
| Sodium | 152 | 34.8 | 0.966 | |
| Strontium | 255 | 2.09 | 0.966 | |
| Silicon | 2250 | 6.96 | 0.966 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|
| HCS330 | 15-06-0463-3-AA | 06/04/15 12:20 | Solid | ICP 7300 | 06/10/15 | 06/17/15 22:50 | 150610L04 |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|------|------|------------|
| Calcium | 168000 | 69.6 | 9.66 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|
| HCS340 | 15-06-0463-4-AA | 06/04/15 13:57 | Solid | ICP 7300 | 06/10/15 | 06/11/15 16:49 | 150610L04 |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|------|------|------------|
| Magnesium | 2320 | 11.8 | 1.04 | |
| Potassium | 1180 | 58.8 | 1.04 | |
| Sodium | 736 | 58.8 | 1.04 | |
| Strontium | 447 | 3.53 | 1.04 | |
| Silicon | 1680 | 11.8 | 1.04 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|
| HCS340 | 15-06-0463-4-AA | 06/04/15 13:57 | Solid | ICP 7300 | 06/10/15 | 06/17/15 22:51 | 150610L04 |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|-----|------|------------|
| Calcium | 293000 | 118 | 10.4 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6010B
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS360 | 15-06-0463-5-AA | 06/04/15 14:24 | Solid | ICP 7300 | 06/10/15 | 06/11/15 16:50 | 150610L04 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|------|------|------------|
| Magnesium | 3060 | 10.4 | 1.02 | |
| Potassium | 2300 | 52.2 | 1.02 | |
| Sodium | 207 | 52.2 | 1.02 | |
| Strontium | 199 | 3.13 | 1.02 | |
| Silicon | 1410 | 10.4 | 1.02 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|
| HCS360 | 15-06-0463-5-AA | 06/04/15 14:24 | Solid | ICP 7300 | 06/10/15 | 06/17/15 22:53 | 150610L04 |
|--------|-----------------|----------------|-------|----------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|-----|------|------------|
| Calcium | 146000 | 104 | 10.2 | |

| | | | | | | | |
|----------|-----------------|----------------|-------|----------|----------|----------------|-----------|
| FDHCS360 | 15-06-0463-6-AA | 06/04/15 14:24 | Solid | ICP 7300 | 06/10/15 | 06/11/15 16:56 | 150610L04 |
|----------|-----------------|----------------|-------|----------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|------|------|------------|
| Magnesium | 3560 | 10.5 | 1.03 | |
| Potassium | 2580 | 52.5 | 1.03 | |
| Sodium | 193 | 52.5 | 1.03 | |
| Strontium | 222 | 3.15 | 1.03 | |
| Silicon | 1690 | 10.5 | 1.03 | |

| | | | | | | | |
|----------|-----------------|----------------|-------|----------|----------|----------------|-----------|
| FDHCS360 | 15-06-0463-6-AA | 06/04/15 14:24 | Solid | ICP 7300 | 06/10/15 | 06/17/15 22:55 | 150610L04 |
|----------|-----------------|----------------|-------|----------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|-----|------|------------|
| Calcium | 165000 | 105 | 10.3 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6010B
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 097-01-002-21188 | N/A | Solid | ICP 7300 | 06/10/15 | 06/11/15 16:37 | 150610L04 |

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|------|-------|------------|
| Calcium | ND | 4.88 | 0.976 | |
| Magnesium | ND | 4.88 | 0.976 | |
| Potassium | ND | 24.4 | 0.976 | |
| Sodium | ND | 24.4 | 0.976 | |
| Strontium | ND | 1.46 | 0.976 | |
| Silicon | ND | 4.88 | 0.976 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS310 | 15-06-0463-1-AA | 06/04/15 11:22 | Solid | ICP/MS 04 | 06/08/15 | 06/12/15 00:51 | 150608L01 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Antimony | ND | 2.48 | 1.00 | |
| Arsenic | ND | 1.24 | 1.00 | |
| Barium | 16.0 | 1.24 | 1.00 | |
| Beryllium | ND | 1.24 | 1.00 | |
| Cadmium | ND | 1.24 | 1.00 | |
| Chromium | 3.38 | 2.48 | 1.00 | |
| Copper | 3.65 | 1.24 | 1.00 | |
| Lead | 5.53 | 1.24 | 1.00 | |
| Nickel | 14.5 | 1.24 | 1.00 | |
| Selenium | ND | 1.24 | 1.00 | |
| Silver | ND | 1.24 | 1.00 | |
| Thallium | ND | 1.24 | 1.00 | |
| Zinc | 24.6 | 6.20 | 1.00 | |
| Aluminum | 1390 | 31.0 | 1.00 | |
| Iron | 1710 | 31.0 | 1.00 | |
| Manganese | 31.6 | 3.10 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS320 | 15-06-0463-2-AA | 06/04/15 11:43 | Solid | ICP/MS 04 | 06/08/15 | 06/12/15 00:55 | 150608L01 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Antimony | ND | 2.82 | 1.00 | |
| Arsenic | ND | 1.41 | 1.00 | |
| Barium | 23.5 | 1.41 | 1.00 | |
| Beryllium | ND | 1.41 | 1.00 | |
| Cadmium | ND | 1.41 | 1.00 | |
| Chromium | 6.41 | 2.82 | 1.00 | |
| Copper | 3.55 | 1.41 | 1.00 | |
| Lead | 6.33 | 1.41 | 1.00 | |
| Nickel | 19.1 | 1.41 | 1.00 | |
| Selenium | ND | 1.41 | 1.00 | |
| Silver | ND | 1.41 | 1.00 | |
| Thallium | ND | 1.41 | 1.00 | |
| Zinc | 13.1 | 7.05 | 1.00 | |
| Aluminum | 1580 | 35.3 | 1.00 | |
| Iron | 2620 | 35.3 | 1.00 | |
| Manganese | 47.8 | 3.53 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS330 | 15-06-0463-3-AA | 06/04/15 12:20 | Solid | ICP/MS 04 | 06/08/15 | 06/12/15 01:11 | 150608L01 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Antimony | ND | 2.88 | 1.00 | |
| Arsenic | 4.92 | 1.44 | 1.00 | |
| Barium | 61.0 | 1.44 | 1.00 | |
| Beryllium | ND | 1.44 | 1.00 | |
| Cadmium | ND | 1.44 | 1.00 | |
| Chromium | 15.3 | 2.88 | 1.00 | |
| Copper | 9.74 | 1.44 | 1.00 | |
| Lead | 10.7 | 1.44 | 1.00 | |
| Nickel | 20.5 | 1.44 | 1.00 | |
| Selenium | ND | 1.44 | 1.00 | |
| Silver | ND | 1.44 | 1.00 | |
| Thallium | ND | 1.44 | 1.00 | |
| Zinc | 28.2 | 7.20 | 1.00 | |
| Aluminum | 10300 | 36.0 | 1.00 | |
| Iron | 10300 | 36.0 | 1.00 | |
| Manganese | 195 | 3.60 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS340 | 15-06-0463-4-AA | 06/04/15 13:57 | Solid | ICP/MS 04 | 06/08/15 | 06/12/15 01:15 | 150608L01 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Antimony | ND | 4.51 | 1.00 | |
| Arsenic | 3.14 | 2.26 | 1.00 | |
| Barium | 67.9 | 2.26 | 1.00 | |
| Beryllium | ND | 2.26 | 1.00 | |
| Cadmium | ND | 2.26 | 1.00 | |
| Chromium | 11.7 | 4.51 | 1.00 | |
| Copper | 12.9 | 2.26 | 1.00 | |
| Lead | 13.5 | 2.26 | 1.00 | |
| Nickel | 21.1 | 2.26 | 1.00 | |
| Selenium | ND | 2.26 | 1.00 | |
| Silver | ND | 2.26 | 1.00 | |
| Thallium | ND | 2.26 | 1.00 | |
| Zinc | 85.1 | 11.3 | 1.00 | |
| Aluminum | 4880 | 56.4 | 1.00 | |
| Iron | 9730 | 56.4 | 1.00 | |
| Manganese | 108 | 5.64 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS360 | 15-06-0463-5-AA | 06/04/15 14:24 | Solid | ICP/MS 04 | 06/08/15 | 06/12/15 01:19 | 150608L01 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Antimony | ND | 4.09 | 1.00 | |
| Arsenic | 3.83 | 2.04 | 1.00 | |
| Barium | 68.9 | 2.04 | 1.00 | |
| Beryllium | ND | 2.04 | 1.00 | |
| Cadmium | ND | 2.04 | 1.00 | |
| Chromium | 16.1 | 4.09 | 1.00 | |
| Copper | 14.3 | 2.04 | 1.00 | |
| Lead | 23.9 | 2.04 | 1.00 | |
| Nickel | 17.0 | 2.04 | 1.00 | |
| Selenium | ND | 2.04 | 1.00 | |
| Silver | ND | 2.04 | 1.00 | |
| Thallium | ND | 2.04 | 1.00 | |
| Zinc | 142 | 10.2 | 1.00 | |
| Aluminum | 10800 | 51.1 | 1.00 | |
| Iron | 9750 | 51.1 | 1.00 | |
| Manganese | 216 | 5.11 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| FDHCS360 | 15-06-0463-6-AA | 06/04/15 14:24 | Solid | ICP/MS 04 | 06/08/15 | 06/12/15 01:23 | 150608L01 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Antimony | ND | 4.10 | 1.00 | |
| Arsenic | 4.15 | 2.05 | 1.00 | |
| Barium | 72.9 | 2.05 | 1.00 | |
| Beryllium | ND | 2.05 | 1.00 | |
| Cadmium | ND | 2.05 | 1.00 | |
| Chromium | 18.0 | 4.10 | 1.00 | |
| Copper | 16.0 | 2.05 | 1.00 | |
| Lead | 26.2 | 2.05 | 1.00 | |
| Nickel | 18.4 | 2.05 | 1.00 | |
| Selenium | ND | 2.05 | 1.00 | |
| Silver | ND | 2.05 | 1.00 | |
| Thallium | ND | 2.05 | 1.00 | |
| Zinc | 160 | 10.2 | 1.00 | |
| Aluminum | 11900 | 51.2 | 1.00 | |
| Iron | 10700 | 51.2 | 1.00 | |
| Manganese | 227 | 5.12 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-621-953 | N/A | Solid | ICP/MS 04 | 06/08/15 | 06/09/15 13:26 | 150608L01 |

| Parameter | Result | RL | DF | Qualifiers |
|-----------|--------|------|------|------------|
| Antimony | ND | 2.00 | 1.00 | |
| Arsenic | ND | 1.00 | 1.00 | |
| Barium | ND | 1.00 | 1.00 | |
| Beryllium | ND | 1.00 | 1.00 | |
| Cadmium | ND | 1.00 | 1.00 | |
| Chromium | ND | 2.00 | 1.00 | |
| Copper | ND | 1.00 | 1.00 | |
| Lead | ND | 1.00 | 1.00 | |
| Nickel | ND | 1.00 | 1.00 | |
| Selenium | ND | 1.00 | 1.00 | |
| Silver | ND | 1.00 | 1.00 | |
| Thallium | ND | 1.00 | 1.00 | |
| Zinc | ND | 5.00 | 1.00 | |
| Aluminum | ND | 25.0 | 1.00 | |
| Iron | ND | 25.0 | 1.00 | |
| Manganese | ND | 2.50 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 7471A Total
Method: EPA 7471A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS310 | 15-06-0463-1-AA | 06/04/15 11:22 | Solid | Mercury 05 | 06/16/15 | 06/16/15 14:39 | 150616L01 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Mercury | ND | 0.103 | 1.00 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|------------|----------|----------------|-----------|
| HCS320 | 15-06-0463-2-AA | 06/04/15 11:43 | Solid | Mercury 05 | 06/16/15 | 06/16/15 14:41 | 150616L01 |
|--------|-----------------|----------------|-------|------------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Mercury | ND | 0.114 | 1.00 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|------------|----------|----------------|-----------|
| HCS330 | 15-06-0463-3-AA | 06/04/15 12:20 | Solid | Mercury 05 | 06/16/15 | 06/16/15 14:43 | 150616L01 |
|--------|-----------------|----------------|-------|------------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Mercury | ND | 0.114 | 1.00 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|------------|----------|----------------|-----------|
| HCS340 | 15-06-0463-4-AA | 06/04/15 13:57 | Solid | Mercury 05 | 06/16/15 | 06/16/15 14:45 | 150616L01 |
|--------|-----------------|----------------|-------|------------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Mercury | ND | 0.182 | 1.00 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|------------|----------|----------------|-----------|
| HCS360 | 15-06-0463-5-AA | 06/04/15 14:24 | Solid | Mercury 05 | 06/16/15 | 06/16/15 14:47 | 150616L01 |
|--------|-----------------|----------------|-------|------------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Mercury | ND | 0.173 | 1.00 | |

| | | | | | | | |
|----------|-----------------|----------------|-------|------------|----------|----------------|-----------|
| FDHCS360 | 15-06-0463-6-AA | 06/04/15 14:24 | Solid | Mercury 05 | 06/16/15 | 06/16/15 14:32 | 150616L01 |
|----------|-----------------|----------------|-------|------------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Mercury | ND | 0.171 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 7471A Total
Method: EPA 7471A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-16-272-1352 | N/A | Solid | Mercury 05 | 06/16/15 | 06/16/15 14:27 | 150616L01 |

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Mercury | ND | 0.0833 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

Page 1 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS310 | 15-06-0463-1-B | 06/04/15 11:22 | Solid | GC 44 | 06/13/15 | 06/16/15 18:22 | 150613L05 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|-----------|-------------------|
| Aldrin | ND | 6.3 | 1.00 | |
| Alpha-BHC | ND | 13 | 1.00 | |
| Beta-BHC | ND | 6.3 | 1.00 | |
| Chlordane | ND | 63 | 1.00 | |
| 4,4'-DDD | ND | 6.3 | 1.00 | |
| 4,4'-DDE | ND | 6.3 | 1.00 | |
| 4,4'-DDT | ND | 6.3 | 1.00 | |
| Delta-BHC | ND | 13 | 1.00 | |
| Dieldrin | ND | 6.3 | 1.00 | |
| Endosulfan I | ND | 6.3 | 1.00 | |
| Endosulfan II | ND | 6.3 | 1.00 | |
| Endosulfan Sulfate | ND | 6.3 | 1.00 | |
| Endrin | ND | 6.3 | 1.00 | |
| Endrin Aldehyde | ND | 6.3 | 1.00 | |
| Endrin Ketone | ND | 6.3 | 1.00 | |
| Gamma-BHC | ND | 6.3 | 1.00 | |
| Heptachlor | ND | 6.3 | 1.00 | |
| Heptachlor Epoxide | ND | 13 | 1.00 | |
| Methoxychlor | ND | 6.3 | 1.00 | |
| Toxaphene | ND | 130 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 125 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 128 | 25-145 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS320 | 15-06-0463-2-B | 06/04/15 11:43 | Solid | GC 44 | 06/13/15 | 06/16/15 18:37 | 150613L05 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|-----------|-------------------|
| Aldrin | ND | 7.1 | 1.00 | |
| Alpha-BHC | ND | 14 | 1.00 | |
| Beta-BHC | ND | 7.1 | 1.00 | |
| Chlordane | ND | 71 | 1.00 | |
| 4,4'-DDD | ND | 7.1 | 1.00 | |
| 4,4'-DDE | ND | 7.1 | 1.00 | |
| 4,4'-DDT | ND | 7.1 | 1.00 | |
| Delta-BHC | ND | 14 | 1.00 | |
| Dieldrin | ND | 7.1 | 1.00 | |
| Endosulfan I | ND | 7.1 | 1.00 | |
| Endosulfan II | ND | 7.1 | 1.00 | |
| Endosulfan Sulfate | ND | 7.1 | 1.00 | |
| Endrin | ND | 7.1 | 1.00 | |
| Endrin Aldehyde | ND | 7.1 | 1.00 | |
| Endrin Ketone | ND | 7.1 | 1.00 | |
| Gamma-BHC | ND | 7.1 | 1.00 | |
| Heptachlor | ND | 7.1 | 1.00 | |
| Heptachlor Epoxide | ND | 14 | 1.00 | |
| Methoxychlor | ND | 7.1 | 1.00 | |
| Toxaphene | ND | 140 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 118 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 122 | 25-145 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS330 | 15-06-0463-3-B | 06/04/15 12:20 | Solid | GC 44 | 06/13/15 | 06/16/15 18:51 | 150613L05 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|-----------|-------------------|
| Aldrin | ND | 7.2 | 1.00 | |
| Alpha-BHC | ND | 14 | 1.00 | |
| Beta-BHC | ND | 7.2 | 1.00 | |
| Chlordane | ND | 72 | 1.00 | |
| 4,4'-DDD | ND | 7.2 | 1.00 | |
| 4,4'-DDE | ND | 7.2 | 1.00 | |
| 4,4'-DDT | ND | 7.2 | 1.00 | |
| Delta-BHC | ND | 14 | 1.00 | |
| Dieldrin | ND | 7.2 | 1.00 | |
| Endosulfan I | ND | 7.2 | 1.00 | |
| Endosulfan II | ND | 7.2 | 1.00 | |
| Endosulfan Sulfate | ND | 7.2 | 1.00 | |
| Endrin | ND | 7.2 | 1.00 | |
| Endrin Aldehyde | ND | 7.2 | 1.00 | |
| Endrin Ketone | ND | 7.2 | 1.00 | |
| Gamma-BHC | ND | 7.2 | 1.00 | |
| Heptachlor | ND | 7.2 | 1.00 | |
| Heptachlor Epoxide | ND | 14 | 1.00 | |
| Methoxychlor | ND | 7.2 | 1.00 | |
| Toxaphene | ND | 140 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 127 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 131 | 25-145 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS340 | 15-06-0463-4-B | 06/04/15 13:57 | Solid | GC 44 | 06/13/15 | 06/16/15 19:05 | 150613L05 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|-----------|-------------------|
| Aldrin | ND | 11 | 1.00 | |
| Alpha-BHC | ND | 22 | 1.00 | |
| Beta-BHC | ND | 11 | 1.00 | |
| Chlordane | ND | 110 | 1.00 | |
| 4,4'-DDD | ND | 11 | 1.00 | |
| 4,4'-DDE | ND | 11 | 1.00 | |
| 4,4'-DDT | ND | 11 | 1.00 | |
| Delta-BHC | ND | 22 | 1.00 | |
| Dieldrin | ND | 11 | 1.00 | |
| Endosulfan I | ND | 11 | 1.00 | |
| Endosulfan II | ND | 11 | 1.00 | |
| Endosulfan Sulfate | ND | 11 | 1.00 | |
| Endrin | ND | 11 | 1.00 | |
| Endrin Aldehyde | ND | 11 | 1.00 | |
| Endrin Ketone | ND | 11 | 1.00 | |
| Gamma-BHC | ND | 11 | 1.00 | |
| Heptachlor | ND | 11 | 1.00 | |
| Heptachlor Epoxide | ND | 22 | 1.00 | |
| Methoxychlor | ND | 11 | 1.00 | |
| Toxaphene | ND | 220 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 117 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 121 | 25-145 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS360 | 15-06-0463-5-B | 06/04/15 14:24 | Solid | GC 44 | 06/13/15 | 06/16/15 19:20 | 150613L05 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|-----------|-------------------|
| Aldrin | ND | 10 | 1.00 | |
| Alpha-BHC | ND | 21 | 1.00 | |
| Beta-BHC | ND | 10 | 1.00 | |
| Chlordane | ND | 100 | 1.00 | |
| 4,4'-DDD | ND | 10 | 1.00 | |
| 4,4'-DDE | ND | 10 | 1.00 | |
| 4,4'-DDT | ND | 10 | 1.00 | |
| Delta-BHC | ND | 21 | 1.00 | |
| Dieldrin | ND | 10 | 1.00 | |
| Endosulfan I | ND | 10 | 1.00 | |
| Endosulfan II | ND | 10 | 1.00 | |
| Endosulfan Sulfate | ND | 10 | 1.00 | |
| Endrin | ND | 10 | 1.00 | |
| Endrin Aldehyde | ND | 10 | 1.00 | |
| Endrin Ketone | ND | 10 | 1.00 | |
| Gamma-BHC | ND | 10 | 1.00 | |
| Heptachlor | ND | 10 | 1.00 | |
| Heptachlor Epoxide | ND | 21 | 1.00 | |
| Methoxychlor | ND | 10 | 1.00 | |
| Toxaphene | ND | 210 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 112 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 125 | 25-145 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|--------------|--------------|-----------------|-----------------------|------------------|
| FDHCS360 | 15-06-0463-6-B | 06/04/15 14:24 | Solid | GC 44 | 06/13/15 | 06/16/15 19:34 | 150613L05 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|-----------|-------------------|
| Aldrin | ND | 10 | 1.00 | |
| Alpha-BHC | ND | 20 | 1.00 | |
| Beta-BHC | ND | 10 | 1.00 | |
| Chlordane | ND | 100 | 1.00 | |
| 4,4'-DDD | ND | 10 | 1.00 | |
| 4,4'-DDE | ND | 10 | 1.00 | |
| 4,4'-DDT | ND | 10 | 1.00 | |
| Delta-BHC | ND | 20 | 1.00 | |
| Dieldrin | ND | 10 | 1.00 | |
| Endosulfan I | ND | 10 | 1.00 | |
| Endosulfan II | ND | 10 | 1.00 | |
| Endosulfan Sulfate | ND | 10 | 1.00 | |
| Endrin | ND | 10 | 1.00 | |
| Endrin Aldehyde | ND | 10 | 1.00 | |
| Endrin Ketone | ND | 10 | 1.00 | |
| Gamma-BHC | ND | 10 | 1.00 | |
| Heptachlor | ND | 10 | 1.00 | |
| Heptachlor Epoxide | ND | 20 | 1.00 | |
| Methoxychlor | ND | 10 | 1.00 | |
| Toxaphene | ND | 200 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 102 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 105 | 25-145 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-537-2140 | N/A | Solid | GC 44 | 06/13/15 | 06/16/15 17:39 | 150613L05 |

| Parameter | Result | RL | DF | Qualifiers |
|--------------------|--------|-----|------|------------|
| Aldrin | ND | 5.0 | 1.00 | |
| Alpha-BHC | ND | 10 | 1.00 | |
| Beta-BHC | ND | 5.0 | 1.00 | |
| Chlordane | ND | 50 | 1.00 | |
| 4,4'-DDD | ND | 5.0 | 1.00 | |
| 4,4'-DDE | ND | 5.0 | 1.00 | |
| 4,4'-DDT | ND | 5.0 | 1.00 | |
| Delta-BHC | ND | 10 | 1.00 | |
| Dieldrin | ND | 5.0 | 1.00 | |
| Endosulfan I | ND | 5.0 | 1.00 | |
| Endosulfan II | ND | 5.0 | 1.00 | |
| Endosulfan Sulfate | ND | 5.0 | 1.00 | |
| Endrin | ND | 5.0 | 1.00 | |
| Endrin Aldehyde | ND | 5.0 | 1.00 | |
| Endrin Ketone | ND | 5.0 | 1.00 | |
| Gamma-BHC | ND | 5.0 | 1.00 | |
| Heptachlor | ND | 5.0 | 1.00 | |
| Heptachlor Epoxide | ND | 10 | 1.00 | |
| Methoxychlor | ND | 5.0 | 1.00 | |
| Toxaphene | ND | 100 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 96 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 101 | 25-145 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8082
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS310 | 15-06-0463-1-B | 06/04/15 11:22 | Solid | GC 31 | 06/13/15 | 06/16/15 14:32 | 150613L06 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|--------------|--------|----|------|------------|
| Aroclor-1016 | ND | 63 | 1.00 | |
| Aroclor-1221 | ND | 63 | 1.00 | |
| Aroclor-1232 | ND | 63 | 1.00 | |
| Aroclor-1242 | ND | 63 | 1.00 | |
| Aroclor-1248 | ND | 63 | 1.00 | |
| Aroclor-1254 | ND | 63 | 1.00 | |
| Aroclor-1260 | ND | 63 | 1.00 | |
| Aroclor-1262 | ND | 63 | 1.00 | |
| Aroclor-1268 | ND | 63 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 108 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 113 | 25-145 | |

| HCS320 | 15-06-0463-2-B | 06/04/15 11:43 | Solid | GC 31 | 06/13/15 | 06/16/15 14:51 | 150613L06 |
|--------|----------------|----------------|-------|-------|----------|----------------|-----------|
|--------|----------------|----------------|-------|-------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|--------------|--------|----|------|------------|
| Aroclor-1016 | ND | 71 | 1.00 | |
| Aroclor-1221 | ND | 71 | 1.00 | |
| Aroclor-1232 | ND | 71 | 1.00 | |
| Aroclor-1242 | ND | 71 | 1.00 | |
| Aroclor-1248 | ND | 71 | 1.00 | |
| Aroclor-1254 | ND | 71 | 1.00 | |
| Aroclor-1260 | ND | 71 | 1.00 | |
| Aroclor-1262 | ND | 71 | 1.00 | |
| Aroclor-1268 | ND | 71 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 89 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 103 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8082
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS330 | 15-06-0463-3-B | 06/04/15 12:20 | Solid | GC 31 | 06/13/15 | 06/16/15 15:10 | 150613L06 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|--------------|--------|----|------|------------|
| Aroclor-1016 | ND | 72 | 1.00 | |
| Aroclor-1221 | ND | 72 | 1.00 | |
| Aroclor-1232 | ND | 72 | 1.00 | |
| Aroclor-1242 | ND | 72 | 1.00 | |
| Aroclor-1248 | ND | 72 | 1.00 | |
| Aroclor-1254 | ND | 72 | 1.00 | |
| Aroclor-1260 | ND | 72 | 1.00 | |
| Aroclor-1262 | ND | 72 | 1.00 | |
| Aroclor-1268 | ND | 72 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 86 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 89 | 25-145 | |

| HCS340 | 15-06-0463-4-B | 06/04/15 13:57 | Solid | GC 31 | 06/13/15 | 06/16/15 15:29 | 150613L06 |
|--------|----------------|----------------|-------|-------|----------|----------------|-----------|
|--------|----------------|----------------|-------|-------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|--------------|--------|-----|------|------------|
| Aroclor-1016 | ND | 110 | 1.00 | |
| Aroclor-1221 | ND | 110 | 1.00 | |
| Aroclor-1232 | ND | 110 | 1.00 | |
| Aroclor-1242 | ND | 110 | 1.00 | |
| Aroclor-1248 | ND | 110 | 1.00 | |
| Aroclor-1254 | ND | 110 | 1.00 | |
| Aroclor-1260 | ND | 110 | 1.00 | |
| Aroclor-1262 | ND | 110 | 1.00 | |
| Aroclor-1268 | ND | 110 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 93 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 96 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8082
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS360 | 15-06-0463-5-B | 06/04/15 14:24 | Solid | GC 31 | 06/13/15 | 06/16/15 15:48 | 150613L06 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|--------------|--------|-----|------|------------|
| Aroclor-1016 | ND | 100 | 1.00 | |
| Aroclor-1221 | ND | 100 | 1.00 | |
| Aroclor-1232 | ND | 100 | 1.00 | |
| Aroclor-1242 | ND | 100 | 1.00 | |
| Aroclor-1248 | ND | 100 | 1.00 | |
| Aroclor-1254 | ND | 100 | 1.00 | |
| Aroclor-1260 | ND | 100 | 1.00 | |
| Aroclor-1262 | ND | 100 | 1.00 | |
| Aroclor-1268 | ND | 100 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 98 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 93 | 25-145 | |

| FDHCS360 | 15-06-0463-6-B | 06/04/15 14:24 | Solid | GC 31 | 06/13/15 | 06/16/15 16:07 | 150613L06 |
|----------|----------------|----------------|-------|-------|----------|----------------|-----------|
|----------|----------------|----------------|-------|-------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|--------------|--------|-----|------|------------|
| Aroclor-1016 | ND | 100 | 1.00 | |
| Aroclor-1221 | ND | 100 | 1.00 | |
| Aroclor-1232 | ND | 100 | 1.00 | |
| Aroclor-1242 | ND | 100 | 1.00 | |
| Aroclor-1248 | ND | 100 | 1.00 | |
| Aroclor-1254 | ND | 100 | 1.00 | |
| Aroclor-1260 | ND | 100 | 1.00 | |
| Aroclor-1262 | ND | 100 | 1.00 | |
| Aroclor-1268 | ND | 100 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 98 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 94 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8082
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-535-3269 | N/A | Solid | GC 31 | 06/13/15 | 06/16/15 11:21 | 150613L06 |

| Parameter | Result | RL | DF | Qualifiers |
|--------------|--------|----|------|------------|
| Aroclor-1016 | ND | 50 | 1.00 | |
| Aroclor-1221 | ND | 50 | 1.00 | |
| Aroclor-1232 | ND | 50 | 1.00 | |
| Aroclor-1242 | ND | 50 | 1.00 | |
| Aroclor-1248 | ND | 50 | 1.00 | |
| Aroclor-1254 | ND | 50 | 1.00 | |
| Aroclor-1260 | ND | 50 | 1.00 | |
| Aroclor-1262 | ND | 50 | 1.00 | |
| Aroclor-1268 | ND | 50 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 89 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 106 | 25-145 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

Page 1 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS310 | 15-06-0463-1-B | 06/04/15 11:22 | Solid | GC 26 | 06/08/15 | 06/16/15 11:59 | 150608L12 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Demeton-o/s | ND | 0.62 | 1.00 | |
| Azinphos Methyl | ND | 0.62 | 1.00 | |
| Bolstar | ND | 0.62 | 1.00 | |
| Chlorpyrifos | ND | 0.62 | 1.00 | |
| Coumaphos | ND | 0.62 | 1.00 | |
| Diazinon | ND | 0.62 | 1.00 | |
| Dichlorvos | ND | 0.62 | 1.00 | |
| Disulfoton | ND | 0.62 | 1.00 | |
| Ethoprop | ND | 0.62 | 1.00 | |
| Fensulfothion | ND | 0.62 | 1.00 | |
| Fenthion | ND | 0.62 | 1.00 | |
| Merphos | ND | 0.62 | 1.00 | |
| Methyl Parathion | ND | 0.62 | 1.00 | |
| Mevinphos | ND | 0.62 | 1.00 | |
| Naled | ND | 5.0 | 1.00 | |
| Phorate | ND | 0.62 | 1.00 | |
| Ronnel | ND | 0.62 | 1.00 | |
| Stirophos | ND | 2.5 | 1.00 | |
| Tokuthion | ND | 0.62 | 1.00 | |
| Trichloronate | ND | 0.62 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 94 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS320 | 15-06-0463-2-B | 06/04/15 11:43 | Solid | GC 26 | 06/08/15 | 06/16/15 14:17 | 150608L12 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Demeton-o/s | ND | 0.70 | 1.00 | |
| Azinphos Methyl | ND | 0.70 | 1.00 | |
| Bolstar | ND | 0.70 | 1.00 | |
| Chlorpyrifos | ND | 0.70 | 1.00 | |
| Coumaphos | ND | 0.70 | 1.00 | |
| Diazinon | ND | 0.70 | 1.00 | |
| Dichlorvos | ND | 0.70 | 1.00 | |
| Disulfoton | ND | 0.70 | 1.00 | |
| Ethoprop | ND | 0.70 | 1.00 | |
| Fensulfothion | ND | 0.70 | 1.00 | |
| Fenthion | ND | 0.70 | 1.00 | |
| Merphos | ND | 0.70 | 1.00 | |
| Methyl Parathion | ND | 0.70 | 1.00 | |
| Mevinphos | ND | 0.70 | 1.00 | |
| Naled | ND | 5.6 | 1.00 | |
| Phorate | ND | 0.70 | 1.00 | |
| Ronnel | ND | 0.70 | 1.00 | |
| Stirophos | ND | 2.8 | 1.00 | |
| Tokuthion | ND | 0.70 | 1.00 | |
| Trichloronate | ND | 0.70 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 98 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS330 | 15-06-0463-3-B | 06/04/15 12:20 | Solid | GC 26 | 06/08/15 | 06/16/15 15:02 | 150608L12 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Demeton-o/s | ND | 0.72 | 1.00 | |
| Azinphos Methyl | ND | 0.72 | 1.00 | |
| Bolstar | ND | 0.72 | 1.00 | |
| Chlorpyrifos | ND | 0.72 | 1.00 | |
| Coumaphos | ND | 0.72 | 1.00 | |
| Diazinon | ND | 0.72 | 1.00 | |
| Dichlorvos | ND | 0.72 | 1.00 | |
| Disulfoton | ND | 0.72 | 1.00 | |
| Ethoprop | ND | 0.72 | 1.00 | |
| Fensulfothion | ND | 0.72 | 1.00 | |
| Fenthion | ND | 0.72 | 1.00 | |
| Merphos | ND | 0.72 | 1.00 | |
| Methyl Parathion | ND | 0.72 | 1.00 | |
| Mevinphos | ND | 0.72 | 1.00 | |
| Naled | ND | 5.8 | 1.00 | |
| Phorate | ND | 0.72 | 1.00 | |
| Ronnel | ND | 0.72 | 1.00 | |
| Stirophos | ND | 2.9 | 1.00 | |
| Tokuthion | ND | 0.72 | 1.00 | |
| Trichloronate | ND | 0.72 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 103 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS340 | 15-06-0463-4-B | 06/04/15 13:57 | Solid | GC 26 | 06/08/15 | 06/16/15 15:46 | 150608L12 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Demeton-o/s | ND | 1.1 | 1.00 | |
| Azinphos Methyl | ND | 1.1 | 1.00 | |
| Bolstar | ND | 1.1 | 1.00 | |
| Chlorpyrifos | ND | 1.1 | 1.00 | |
| Coumaphos | ND | 1.1 | 1.00 | |
| Diazinon | ND | 1.1 | 1.00 | |
| Dichlorvos | ND | 1.1 | 1.00 | |
| Disulfoton | ND | 1.1 | 1.00 | |
| Ethoprop | ND | 1.1 | 1.00 | |
| Fensulfothion | ND | 1.1 | 1.00 | |
| Fenthion | ND | 1.1 | 1.00 | |
| Merphos | ND | 1.1 | 1.00 | |
| Methyl Parathion | ND | 1.1 | 1.00 | |
| Mevinphos | ND | 1.1 | 1.00 | |
| Naled | ND | 8.9 | 1.00 | |
| Phorate | ND | 1.1 | 1.00 | |
| Ronnel | ND | 1.1 | 1.00 | |
| Stirophos | ND | 4.5 | 1.00 | |
| Tokuthion | ND | 1.1 | 1.00 | |
| Trichloronate | ND | 1.1 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 94 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS360 | 15-06-0463-5-B | 06/04/15 14:24 | Solid | GC 26 | 06/08/15 | 06/16/15 17:14 | 150608L12 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|-----------|-------------------|
| Demeton-o/s | ND | 1.0 | 1.00 | |
| Azinphos Methyl | ND | 1.0 | 1.00 | |
| Bolstar | ND | 1.0 | 1.00 | |
| Chlorpyrifos | ND | 1.0 | 1.00 | |
| Coumaphos | ND | 1.0 | 1.00 | |
| Diazinon | ND | 1.0 | 1.00 | |
| Dichlorvos | ND | 1.0 | 1.00 | |
| Disulfoton | ND | 1.0 | 1.00 | |
| Ethoprop | ND | 1.0 | 1.00 | |
| Fensulfothion | ND | 1.0 | 1.00 | |
| Fenthion | ND | 1.0 | 1.00 | |
| Merphos | ND | 1.0 | 1.00 | |
| Methyl Parathion | ND | 1.0 | 1.00 | |
| Mevinphos | ND | 1.0 | 1.00 | |
| Naled | ND | 8.2 | 1.00 | |
| Phorate | ND | 1.0 | 1.00 | |
| Ronnel | ND | 1.0 | 1.00 | |
| Stirophos | ND | 4.1 | 1.00 | |
| Tokuthion | ND | 1.0 | 1.00 | |
| Trichloronate | ND | 1.0 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 97 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| FDHCS360 | 15-06-0463-6-B | 06/04/15 14:24 | Solid | GC 26 | 06/08/15 | 06/16/15 19:02 | 150608L12 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|-------------------|
| Demeton-o/s | ND | 1.0 | 1.00 | |
| Azinphos Methyl | ND | 1.0 | 1.00 | |
| Bolstar | ND | 1.0 | 1.00 | |
| Chlorpyrifos | ND | 1.0 | 1.00 | |
| Coumaphos | ND | 1.0 | 1.00 | |
| Diazinon | ND | 1.0 | 1.00 | |
| Dichlorvos | ND | 1.0 | 1.00 | |
| Disulfoton | ND | 1.0 | 1.00 | |
| Ethoprop | ND | 1.0 | 1.00 | |
| Fensulfothion | ND | 1.0 | 1.00 | |
| Fenthion | ND | 1.0 | 1.00 | |
| Merphos | ND | 1.0 | 1.00 | |
| Methyl Parathion | ND | 1.0 | 1.00 | |
| Mevinphos | ND | 1.0 | 1.00 | |
| Naled | ND | 8.2 | 1.00 | |
| Phorate | ND | 1.0 | 1.00 | |
| Ronnel | ND | 1.0 | 1.00 | |
| Stirophos | ND | 4.1 | 1.00 | |
| Tokuthion | ND | 1.0 | 1.00 | |
| Trichloronate | ND | 1.0 | 1.00 | |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | |
| Tributylphosphate | 119 | 30-130 | | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-973-187 | N/A | Solid | GC 26 | 06/08/15 | 06/09/15 11:41 | 150608L12 |

| Parameter | Result | RL | DF | Qualifiers |
|------------------|--------|------|------|------------|
| Demeton-o/s | ND | 0.50 | 1.00 | |
| Azinphos Methyl | ND | 0.50 | 1.00 | |
| Bolstar | ND | 0.50 | 1.00 | |
| Chlorpyrifos | ND | 0.50 | 1.00 | |
| Coumaphos | ND | 0.50 | 1.00 | |
| Diazinon | ND | 0.50 | 1.00 | |
| Dichlorvos | ND | 0.50 | 1.00 | |
| Disulfoton | ND | 0.50 | 1.00 | |
| Ethoprop | ND | 0.50 | 1.00 | |
| Fensulfothion | ND | 0.50 | 1.00 | |
| Fenthion | ND | 0.50 | 1.00 | |
| Merphos | ND | 0.50 | 1.00 | |
| Methyl Parathion | ND | 0.50 | 1.00 | |
| Mevinphos | ND | 0.50 | 1.00 | |
| Naled | ND | 4.0 | 1.00 | |
| Phorate | ND | 0.50 | 1.00 | |
| Ronnel | ND | 0.50 | 1.00 | |
| Stirophos | ND | 2.0 | 1.00 | |
| Tokuthion | ND | 0.50 | 1.00 | |
| Trichloronate | ND | 0.50 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 99 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS310 | 15-06-0463-1-B | 06/04/15 11:22 | Solid | GC 40 | 06/10/15 | 06/16/15 15:14 | 150610L09 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-------------------|--------|-------|------|------------|
| Dalapon | ND | 310 | 1.00 | |
| Dicamba | ND | 12 | 1.00 | |
| MCP | ND | 12000 | 1.00 | |
| MCPA | ND | 12000 | 1.00 | |
| Dichlorprop | ND | 120 | 1.00 | |
| 2,4-D | ND | 120 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 12 | 1.00 | |
| 2,4,5-T | ND | 12 | 1.00 | |
| 2,4-DB | ND | 120 | 1.00 | |
| Dinoseb | ND | 62 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 51 | 30-130 | |

| | | | | | | | |
|--------|----------------|----------------|-------|-------|----------|----------------|-----------|
| HCS320 | 15-06-0463-2-B | 06/04/15 11:43 | Solid | GC 40 | 06/10/15 | 06/16/15 15:41 | 150610L09 |
|--------|----------------|----------------|-------|-------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-------------------|--------|-------|------|------------|
| Dalapon | ND | 350 | 1.00 | |
| Dicamba | ND | 14 | 1.00 | |
| MCP | ND | 14000 | 1.00 | |
| MCPA | ND | 14000 | 1.00 | |
| Dichlorprop | ND | 140 | 1.00 | |
| 2,4-D | ND | 140 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 14 | 1.00 | |
| 2,4,5-T | ND | 14 | 1.00 | |
| 2,4-DB | ND | 140 | 1.00 | |
| Dinoseb | ND | 71 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 58 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS330 | 15-06-0463-3-B | 06/04/15 12:20 | Solid | GC 40 | 06/10/15 | 06/16/15 16:04 | 150610L09 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-------------------|--------|-------|------|------------|
| Dalapon | ND | 360 | 1.00 | |
| Dicamba | ND | 14 | 1.00 | |
| MCP | ND | 14000 | 1.00 | |
| MCPA | ND | 14000 | 1.00 | |
| Dichlorprop | ND | 140 | 1.00 | |
| 2,4-D | ND | 140 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 14 | 1.00 | |
| 2,4,5-T | ND | 14 | 1.00 | |
| 2,4-DB | ND | 140 | 1.00 | |
| Dinoseb | ND | 72 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 64 | 30-130 | |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS340 | 15-06-0463-4-B | 06/04/15 13:57 | Solid | GC 40 | 06/10/15 | 06/16/15 16:27 | 150610L09 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-------------------|--------|-------|------|------------|
| Dalapon | ND | 570 | 1.00 | |
| Dicamba | ND | 23 | 1.00 | |
| MCP | ND | 23000 | 1.00 | |
| MCPA | ND | 23000 | 1.00 | |
| Dichlorprop | ND | 230 | 1.00 | |
| 2,4-D | ND | 230 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 23 | 1.00 | |
| 2,4,5-T | ND | 23 | 1.00 | |
| 2,4-DB | ND | 230 | 1.00 | |
| Dinoseb | ND | 110 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 66 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS360 | 15-06-0463-5-B | 06/04/15 14:24 | Solid | GC 40 | 06/10/15 | 06/16/15 16:51 | 150610L09 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-------------------|--------|-------|------|------------|
| Dalapon | ND | 510 | 1.00 | |
| Dicamba | ND | 21 | 1.00 | |
| MCP | ND | 21000 | 1.00 | |
| MCPA | ND | 21000 | 1.00 | |
| Dichlorprop | ND | 210 | 1.00 | |
| 2,4-D | ND | 210 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 21 | 1.00 | |
| 2,4,5-T | ND | 21 | 1.00 | |
| 2,4-DB | ND | 210 | 1.00 | |
| Dinoseb | ND | 100 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 71 | 30-130 | |

| FDHCS360 | 15-06-0463-6-B | 06/04/15 14:24 | Solid | GC 40 | 06/10/15 | 06/16/15 17:14 | 150610L09 |
|----------|----------------|----------------|-------|-------|----------|----------------|-----------|
|----------|----------------|----------------|-------|-------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-------------------|--------|-------|------|------------|
| Dalapon | ND | 510 | 1.00 | |
| Dicamba | ND | 20 | 1.00 | |
| MCP | ND | 20000 | 1.00 | |
| MCPA | ND | 20000 | 1.00 | |
| Dichlorprop | ND | 200 | 1.00 | |
| 2,4-D | ND | 200 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 20 | 1.00 | |
| 2,4,5-T | ND | 20 | 1.00 | |
| 2,4-DB | ND | 200 | 1.00 | |
| Dinoseb | ND | 100 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 67 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-033-1294 | N/A | Solid | GC 40 | 06/10/15 | 06/11/15 19:20 | 150610L09 |

| Parameter | Result | RL | DF | Qualifiers |
|-------------------|--------|-------|------|------------|
| Dalapon | ND | 250 | 1.00 | |
| Dicamba | ND | 10 | 1.00 | |
| MCPD | ND | 10000 | 1.00 | |
| MCPA | ND | 10000 | 1.00 | |
| Dichlorprop | ND | 100 | 1.00 | |
| 2,4-D | ND | 100 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 10 | 1.00 | |
| 2,4,5-T | ND | 10 | 1.00 | |
| 2,4-DB | ND | 100 | 1.00 | |
| Dinoseb | ND | 50 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 67 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

Page 1 of 21

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS310 | 15-06-0463-1-B | 06/04/15 11:22 | Solid | GC/MS CCC | 06/13/15 | 06/17/15 01:07 | 150613L07 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|------------------------------|--------|------|------|------------|
| Acenaphthene | ND | 0.62 | 1.00 | |
| Acenaphthylene | ND | 0.62 | 1.00 | |
| Aniline | ND | 0.62 | 1.00 | |
| Anthracene | ND | 0.62 | 1.00 | |
| Azobenzene | ND | 0.62 | 1.00 | |
| Benzidine | ND | 12 | 1.00 | |
| Benzo (a) Anthracene | ND | 0.62 | 1.00 | |
| Benzo (a) Pyrene | ND | 0.62 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 0.62 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 0.62 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 0.62 | 1.00 | |
| Benzoic Acid | ND | 3.1 | 1.00 | |
| Benzyl Alcohol | ND | 0.62 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 0.62 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 3.1 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 0.62 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 0.62 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 0.62 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 0.62 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 0.62 | 1.00 | |
| 4-Chloroaniline | ND | 0.62 | 1.00 | |
| 2-Chloronaphthalene | ND | 0.62 | 1.00 | |
| 2-Chlorophenol | ND | 0.62 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 0.62 | 1.00 | |
| Chrysene | ND | 0.62 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 0.62 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 0.62 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 0.62 | 1.00 | |
| Dibenzofuran | ND | 0.62 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 0.62 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 0.62 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 0.62 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 12 | 1.00 | |
| 2,4-Dichlorophenol | ND | 0.62 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Parameter | Result | RL | DF | Qualifiers |
|----------------------------|--------|------|------|------------|
| Diethyl Phthalate | ND | 0.62 | 1.00 | |
| Dimethyl Phthalate | ND | 0.62 | 1.00 | |
| 2,4-Dimethylphenol | ND | 0.62 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 3.1 | 1.00 | |
| 2,4-Dinitrophenol | ND | 3.1 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 0.62 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 0.62 | 1.00 | |
| Fluoranthene | ND | 0.62 | 1.00 | |
| Fluorene | ND | 0.62 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 0.62 | 1.00 | |
| Hexachlorobenzene | ND | 0.62 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 3.1 | 1.00 | |
| Hexachloroethane | ND | 0.62 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 0.62 | 1.00 | |
| Isophorone | ND | 0.62 | 1.00 | |
| 2-Methylnaphthalene | ND | 0.62 | 1.00 | |
| 1-Methylnaphthalene | ND | 0.62 | 1.00 | |
| 2-Methylphenol | ND | 0.62 | 1.00 | |
| 3/4-Methylphenol | ND | 0.62 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 0.62 | 1.00 | |
| N-Nitrosodimethylamine | ND | 0.62 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 0.62 | 1.00 | |
| Naphthalene | ND | 0.62 | 1.00 | |
| 4-Nitroaniline | ND | 0.62 | 1.00 | |
| 3-Nitroaniline | ND | 0.62 | 1.00 | |
| 2-Nitroaniline | ND | 0.62 | 1.00 | |
| Nitrobenzene | ND | 3.1 | 1.00 | |
| 4-Nitrophenol | ND | 0.62 | 1.00 | |
| 2-Nitrophenol | ND | 0.62 | 1.00 | |
| Pentachlorophenol | ND | 3.1 | 1.00 | |
| Phenanthrene | ND | 0.62 | 1.00 | |
| Phenol | ND | 0.62 | 1.00 | |
| Pyrene | ND | 0.62 | 1.00 | |
| Pyridine | ND | 0.62 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 0.62 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 0.62 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 0.62 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 52 | 27-120 | |
| 2-Fluorophenol | 53 | 25-120 | |
| Nitrobenzene-d5 | 43 | 33-123 | |
| p-Terphenyl-d14 | 59 | 27-159 | |
| Phenol-d6 | 56 | 26-122 | |
| 2,4,6-Tribromophenol | 62 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS320 | 15-06-0463-2-B | 06/04/15 11:43 | Solid | GC/MS CCC | 06/13/15 | 06/17/15 01:25 | 150613L07 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|------------------------------|--------|------|------|------------|
| Acenaphthene | ND | 0.72 | 1.00 | |
| Acenaphthylene | ND | 0.72 | 1.00 | |
| Aniline | ND | 0.72 | 1.00 | |
| Anthracene | ND | 0.72 | 1.00 | |
| Azobenzene | ND | 0.72 | 1.00 | |
| Benzidine | ND | 14 | 1.00 | |
| Benzo (a) Anthracene | ND | 0.72 | 1.00 | |
| Benzo (a) Pyrene | ND | 0.72 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 0.72 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 0.72 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 0.72 | 1.00 | |
| Benzoic Acid | ND | 3.6 | 1.00 | |
| Benzyl Alcohol | ND | 0.72 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 0.72 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 3.6 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 0.72 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | 0.94 | 0.72 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 0.72 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 0.72 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 0.72 | 1.00 | |
| 4-Chloroaniline | ND | 0.72 | 1.00 | |
| 2-Chloronaphthalene | ND | 0.72 | 1.00 | |
| 2-Chlorophenol | ND | 0.72 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 0.72 | 1.00 | |
| Chrysene | ND | 0.72 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 0.72 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 0.72 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 0.72 | 1.00 | |
| Dibenzofuran | ND | 0.72 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 0.72 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 0.72 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 0.72 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 14 | 1.00 | |
| 2,4-Dichlorophenol | ND | 0.72 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Parameter | Result | RL | DF | Qualifiers |
|----------------------------|--------|------|------|------------|
| Diethyl Phthalate | ND | 0.72 | 1.00 | |
| Dimethyl Phthalate | ND | 0.72 | 1.00 | |
| 2,4-Dimethylphenol | ND | 0.72 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 3.6 | 1.00 | |
| 2,4-Dinitrophenol | ND | 3.6 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 0.72 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 0.72 | 1.00 | |
| Fluoranthene | ND | 0.72 | 1.00 | |
| Fluorene | ND | 0.72 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 0.72 | 1.00 | |
| Hexachlorobenzene | ND | 0.72 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 3.6 | 1.00 | |
| Hexachloroethane | ND | 0.72 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 0.72 | 1.00 | |
| Isophorone | ND | 0.72 | 1.00 | |
| 2-Methylnaphthalene | ND | 0.72 | 1.00 | |
| 1-Methylnaphthalene | ND | 0.72 | 1.00 | |
| 2-Methylphenol | ND | 0.72 | 1.00 | |
| 3/4-Methylphenol | ND | 0.72 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 0.72 | 1.00 | |
| N-Nitrosodimethylamine | ND | 0.72 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 0.72 | 1.00 | |
| Naphthalene | ND | 0.72 | 1.00 | |
| 4-Nitroaniline | ND | 0.72 | 1.00 | |
| 3-Nitroaniline | ND | 0.72 | 1.00 | |
| 2-Nitroaniline | ND | 0.72 | 1.00 | |
| Nitrobenzene | ND | 3.6 | 1.00 | |
| 4-Nitrophenol | ND | 0.72 | 1.00 | |
| 2-Nitrophenol | ND | 0.72 | 1.00 | |
| Pentachlorophenol | ND | 3.6 | 1.00 | |
| Phenanthrene | ND | 0.72 | 1.00 | |
| Phenol | ND | 0.72 | 1.00 | |
| Pyrene | ND | 0.72 | 1.00 | |
| Pyridine | ND | 0.72 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 0.72 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 0.72 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 0.72 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 53 | 27-120 | |
| 2-Fluorophenol | 51 | 25-120 | |
| Nitrobenzene-d5 | 43 | 33-123 | |
| p-Terphenyl-d14 | 67 | 27-159 | |
| Phenol-d6 | 55 | 26-122 | |
| 2,4,6-Tribromophenol | 70 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS330 | 15-06-0463-3-B | 06/04/15 12:20 | Solid | GC/MS CCC | 06/13/15 | 06/17/15 01:43 | 150613L07 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|------------------------------|--------|------|------|------------|
| Acenaphthene | ND | 0.73 | 1.00 | |
| Acenaphthylene | ND | 0.73 | 1.00 | |
| Aniline | ND | 0.73 | 1.00 | |
| Anthracene | ND | 0.73 | 1.00 | |
| Azobenzene | ND | 0.73 | 1.00 | |
| Benzidine | ND | 15 | 1.00 | |
| Benzo (a) Anthracene | ND | 0.73 | 1.00 | |
| Benzo (a) Pyrene | ND | 0.73 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 0.73 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 0.73 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 0.73 | 1.00 | |
| Benzoic Acid | ND | 3.7 | 1.00 | |
| Benzyl Alcohol | ND | 0.73 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 0.73 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 3.7 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 0.73 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 0.73 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 0.73 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 0.73 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 0.73 | 1.00 | |
| 4-Chloroaniline | ND | 0.73 | 1.00 | |
| 2-Chloronaphthalene | ND | 0.73 | 1.00 | |
| 2-Chlorophenol | ND | 0.73 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 0.73 | 1.00 | |
| Chrysene | ND | 0.73 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 0.73 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 0.73 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 0.73 | 1.00 | |
| Dibenzofuran | ND | 0.73 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 0.73 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 0.73 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 0.73 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 15 | 1.00 | |
| 2,4-Dichlorophenol | ND | 0.73 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|----------------------------|---------------|-----------|-----------|-------------------|
| Diethyl Phthalate | ND | 0.73 | 1.00 | |
| Dimethyl Phthalate | ND | 0.73 | 1.00 | |
| 2,4-Dimethylphenol | ND | 0.73 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 3.7 | 1.00 | |
| 2,4-Dinitrophenol | ND | 3.7 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 0.73 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 0.73 | 1.00 | |
| Fluoranthene | ND | 0.73 | 1.00 | |
| Fluorene | ND | 0.73 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 0.73 | 1.00 | |
| Hexachlorobenzene | ND | 0.73 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 3.7 | 1.00 | |
| Hexachloroethane | ND | 0.73 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 0.73 | 1.00 | |
| Isophorone | ND | 0.73 | 1.00 | |
| 2-Methylnaphthalene | ND | 0.73 | 1.00 | |
| 1-Methylnaphthalene | ND | 0.73 | 1.00 | |
| 2-Methylphenol | ND | 0.73 | 1.00 | |
| 3/4-Methylphenol | ND | 0.73 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 0.73 | 1.00 | |
| N-Nitrosodimethylamine | ND | 0.73 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 0.73 | 1.00 | |
| Naphthalene | ND | 0.73 | 1.00 | |
| 4-Nitroaniline | ND | 0.73 | 1.00 | |
| 3-Nitroaniline | ND | 0.73 | 1.00 | |
| 2-Nitroaniline | ND | 0.73 | 1.00 | |
| Nitrobenzene | ND | 3.7 | 1.00 | |
| 4-Nitrophenol | ND | 0.73 | 1.00 | |
| 2-Nitrophenol | ND | 0.73 | 1.00 | |
| Pentachlorophenol | ND | 3.7 | 1.00 | |
| Phenanthrene | ND | 0.73 | 1.00 | |
| Phenol | ND | 0.73 | 1.00 | |
| Pyrene | ND | 0.73 | 1.00 | |
| Pyridine | ND | 0.73 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 0.73 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 0.73 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 0.73 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 48 | 27-120 | |
| 2-Fluorophenol | 45 | 25-120 | |
| Nitrobenzene-d5 | 44 | 33-123 | |
| p-Terphenyl-d14 | 58 | 27-159 | |
| Phenol-d6 | 48 | 26-122 | |
| 2,4,6-Tribromophenol | 60 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS340 | 15-06-0463-4-B | 06/04/15 13:57 | Solid | GC/MS CCC | 06/13/15 | 06/17/15 02:01 | 150613L07 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|------------------------------|--------|-----|------|------------|
| Acenaphthene | ND | 1.1 | 1.00 | |
| Acenaphthylene | ND | 1.1 | 1.00 | |
| Aniline | ND | 1.1 | 1.00 | |
| Anthracene | ND | 1.1 | 1.00 | |
| Azobenzene | ND | 1.1 | 1.00 | |
| Benzidine | ND | 23 | 1.00 | |
| Benzo (a) Anthracene | ND | 1.1 | 1.00 | |
| Benzo (a) Pyrene | ND | 1.1 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 1.1 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 1.1 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 1.1 | 1.00 | |
| Benzoic Acid | ND | 5.6 | 1.00 | |
| Benzyl Alcohol | ND | 1.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 1.1 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 5.6 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 1.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 1.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 1.1 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 1.1 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 1.1 | 1.00 | |
| 4-Chloroaniline | ND | 1.1 | 1.00 | |
| 2-Chloronaphthalene | ND | 1.1 | 1.00 | |
| 2-Chlorophenol | ND | 1.1 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 1.1 | 1.00 | |
| Chrysene | ND | 1.1 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 1.1 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 1.1 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 1.1 | 1.00 | |
| Dibenzofuran | ND | 1.1 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.1 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.1 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 23 | 1.00 | |
| 2,4-Dichlorophenol | ND | 1.1 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|----------------------------|---------------|-----------|-----------|-------------------|
| Diethyl Phthalate | ND | 1.1 | 1.00 | |
| Dimethyl Phthalate | ND | 1.1 | 1.00 | |
| 2,4-Dimethylphenol | ND | 1.1 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 5.6 | 1.00 | |
| 2,4-Dinitrophenol | ND | 5.6 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 1.1 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 1.1 | 1.00 | |
| Fluoranthene | ND | 1.1 | 1.00 | |
| Fluorene | ND | 1.1 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 1.1 | 1.00 | |
| Hexachlorobenzene | ND | 1.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 5.6 | 1.00 | |
| Hexachloroethane | ND | 1.1 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 1.1 | 1.00 | |
| Isophorone | ND | 1.1 | 1.00 | |
| 2-Methylnaphthalene | ND | 1.1 | 1.00 | |
| 1-Methylnaphthalene | ND | 1.1 | 1.00 | |
| 2-Methylphenol | ND | 1.1 | 1.00 | |
| 3/4-Methylphenol | ND | 1.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 1.1 | 1.00 | |
| N-Nitrosodimethylamine | ND | 1.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 1.1 | 1.00 | |
| Naphthalene | ND | 1.1 | 1.00 | |
| 4-Nitroaniline | ND | 1.1 | 1.00 | |
| 3-Nitroaniline | ND | 1.1 | 1.00 | |
| 2-Nitroaniline | ND | 1.1 | 1.00 | |
| Nitrobenzene | ND | 5.6 | 1.00 | |
| 4-Nitrophenol | ND | 1.1 | 1.00 | |
| 2-Nitrophenol | ND | 1.1 | 1.00 | |
| Pentachlorophenol | ND | 5.6 | 1.00 | |
| Phenanthrene | ND | 1.1 | 1.00 | |
| Phenol | ND | 1.1 | 1.00 | |
| Pyrene | ND | 1.1 | 1.00 | |
| Pyridine | ND | 1.1 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.1 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 1.1 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 1.1 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

Page 12 of 21

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 55 | 27-120 | |
| 2-Fluorophenol | 55 | 25-120 | |
| Nitrobenzene-d5 | 47 | 33-123 | |
| p-Terphenyl-d14 | 62 | 27-159 | |
| Phenol-d6 | 57 | 26-122 | |
| 2,4,6-Tribromophenol | 64 | 18-138 | |



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

Page 13 of 21

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS360 | 15-06-0463-5-B | 06/04/15 14:24 | Solid | GC/MS CCC | 06/13/15 | 06/17/15 02:19 | 150613L07 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|------------------------------|--------|-----|------|------------|
| Acenaphthene | ND | 1.0 | 1.00 | |
| Acenaphthylene | ND | 1.0 | 1.00 | |
| Aniline | ND | 1.0 | 1.00 | |
| Anthracene | ND | 1.0 | 1.00 | |
| Azobenzene | ND | 1.0 | 1.00 | |
| Benzidine | ND | 20 | 1.00 | |
| Benzo (a) Anthracene | ND | 1.0 | 1.00 | |
| Benzo (a) Pyrene | ND | 1.0 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 1.0 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 1.0 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 1.0 | 1.00 | |
| Benzoic Acid | ND | 5.1 | 1.00 | |
| Benzyl Alcohol | ND | 1.0 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 1.0 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 5.1 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 1.0 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 1.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 1.0 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 1.0 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 1.0 | 1.00 | |
| 4-Chloroaniline | ND | 1.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 1.0 | 1.00 | |
| 2-Chlorophenol | ND | 1.0 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 1.0 | 1.00 | |
| Chrysene | ND | 1.0 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 1.0 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 1.0 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 1.0 | 1.00 | |
| Dibenzofuran | ND | 1.0 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 20 | 1.00 | |
| 2,4-Dichlorophenol | ND | 1.0 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|----------------------------|---------------|-----------|-----------|-------------------|
| Diethyl Phthalate | ND | 1.0 | 1.00 | |
| Dimethyl Phthalate | ND | 1.0 | 1.00 | |
| 2,4-Dimethylphenol | ND | 1.0 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 5.1 | 1.00 | |
| 2,4-Dinitrophenol | ND | 5.1 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 1.0 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 1.0 | 1.00 | |
| Fluoranthene | ND | 1.0 | 1.00 | |
| Fluorene | ND | 1.0 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 1.0 | 1.00 | |
| Hexachlorobenzene | ND | 1.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 5.1 | 1.00 | |
| Hexachloroethane | ND | 1.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 1.0 | 1.00 | |
| Isophorone | ND | 1.0 | 1.00 | |
| 2-Methylnaphthalene | ND | 1.0 | 1.00 | |
| 1-Methylnaphthalene | ND | 1.0 | 1.00 | |
| 2-Methylphenol | ND | 1.0 | 1.00 | |
| 3/4-Methylphenol | ND | 1.0 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 1.0 | 1.00 | |
| N-Nitrosodimethylamine | ND | 1.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 1.0 | 1.00 | |
| Naphthalene | ND | 1.0 | 1.00 | |
| 4-Nitroaniline | ND | 1.0 | 1.00 | |
| 3-Nitroaniline | ND | 1.0 | 1.00 | |
| 2-Nitroaniline | ND | 1.0 | 1.00 | |
| Nitrobenzene | ND | 5.1 | 1.00 | |
| 4-Nitrophenol | ND | 1.0 | 1.00 | |
| 2-Nitrophenol | ND | 1.0 | 1.00 | |
| Pentachlorophenol | ND | 5.1 | 1.00 | |
| Phenanthrene | ND | 1.0 | 1.00 | |
| Phenol | ND | 1.0 | 1.00 | |
| Pyrene | ND | 1.0 | 1.00 | |
| Pyridine | ND | 1.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 1.0 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 1.0 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 66 | 27-120 | |
| 2-Fluorophenol | 62 | 25-120 | |
| Nitrobenzene-d5 | 63 | 33-123 | |
| p-Terphenyl-d14 | 70 | 27-159 | |
| Phenol-d6 | 65 | 26-122 | |
| 2,4,6-Tribromophenol | 77 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| FDHCS360 | 15-06-0463-6-B | 06/04/15 14:24 | Solid | GC/MS CCC | 06/13/15 | 06/17/15 02:37 | 150613L07 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|------------------------------|--------|-----|------|------------|
| Acenaphthene | ND | 1.0 | 1.00 | |
| Acenaphthylene | ND | 1.0 | 1.00 | |
| Aniline | ND | 1.0 | 1.00 | |
| Anthracene | ND | 1.0 | 1.00 | |
| Azobenzene | ND | 1.0 | 1.00 | |
| Benzidine | ND | 20 | 1.00 | |
| Benzo (a) Anthracene | ND | 1.0 | 1.00 | |
| Benzo (a) Pyrene | ND | 1.0 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 1.0 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 1.0 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 1.0 | 1.00 | |
| Benzoic Acid | ND | 5.1 | 1.00 | |
| Benzyl Alcohol | ND | 1.0 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 1.0 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 5.1 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 1.0 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 1.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 1.0 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 1.0 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 1.0 | 1.00 | |
| 4-Chloroaniline | ND | 1.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 1.0 | 1.00 | |
| 2-Chlorophenol | ND | 1.0 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 1.0 | 1.00 | |
| Chrysene | ND | 1.0 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 1.0 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 1.0 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 1.0 | 1.00 | |
| Dibenzofuran | ND | 1.0 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 20 | 1.00 | |
| 2,4-Dichlorophenol | ND | 1.0 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|----------------------------|---------------|-----------|-----------|-------------------|
| Diethyl Phthalate | ND | 1.0 | 1.00 | |
| Dimethyl Phthalate | ND | 1.0 | 1.00 | |
| 2,4-Dimethylphenol | ND | 1.0 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 5.1 | 1.00 | |
| 2,4-Dinitrophenol | ND | 5.1 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 1.0 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 1.0 | 1.00 | |
| Fluoranthene | ND | 1.0 | 1.00 | |
| Fluorene | ND | 1.0 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 1.0 | 1.00 | |
| Hexachlorobenzene | ND | 1.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 5.1 | 1.00 | |
| Hexachloroethane | ND | 1.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 1.0 | 1.00 | |
| Isophorone | ND | 1.0 | 1.00 | |
| 2-Methylnaphthalene | ND | 1.0 | 1.00 | |
| 1-Methylnaphthalene | ND | 1.0 | 1.00 | |
| 2-Methylphenol | ND | 1.0 | 1.00 | |
| 3/4-Methylphenol | ND | 1.0 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 1.0 | 1.00 | |
| N-Nitrosodimethylamine | ND | 1.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 1.0 | 1.00 | |
| Naphthalene | ND | 1.0 | 1.00 | |
| 4-Nitroaniline | ND | 1.0 | 1.00 | |
| 3-Nitroaniline | ND | 1.0 | 1.00 | |
| 2-Nitroaniline | ND | 1.0 | 1.00 | |
| Nitrobenzene | ND | 5.1 | 1.00 | |
| 4-Nitrophenol | ND | 1.0 | 1.00 | |
| 2-Nitrophenol | ND | 1.0 | 1.00 | |
| Pentachlorophenol | ND | 5.1 | 1.00 | |
| Phenanthrene | ND | 1.0 | 1.00 | |
| Phenol | ND | 1.0 | 1.00 | |
| Pyrene | ND | 1.0 | 1.00 | |
| Pyridine | ND | 1.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 1.0 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 1.0 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 64 | 27-120 | |
| 2-Fluorophenol | 62 | 25-120 | |
| Nitrobenzene-d5 | 60 | 33-123 | |
| p-Terphenyl-d14 | 68 | 27-159 | |
| Phenol-d6 | 65 | 26-122 | |
| 2,4,6-Tribromophenol | 74 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-549-3310 | N/A | Solid | GC/MS SS | 06/13/15 | 06/15/15 21:16 | 150613L07 |

| Parameter | Result | RL | DF | Qualifiers |
|------------------------------|--------|------|------|------------|
| Acenaphthene | ND | 0.50 | 1.00 | |
| Acenaphthylene | ND | 0.50 | 1.00 | |
| Aniline | ND | 0.50 | 1.00 | |
| Anthracene | ND | 0.50 | 1.00 | |
| Azobenzene | ND | 0.50 | 1.00 | |
| Benzidine | ND | 10 | 1.00 | |
| Benzo (a) Anthracene | ND | 0.50 | 1.00 | |
| Benzo (a) Pyrene | ND | 0.50 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 0.50 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 0.50 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 0.50 | 1.00 | |
| Benzoic Acid | ND | 2.5 | 1.00 | |
| Benzyl Alcohol | ND | 0.50 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 0.50 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 0.50 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 0.50 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 0.50 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 0.50 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 0.50 | 1.00 | |
| 4-Chloroaniline | ND | 0.50 | 1.00 | |
| 2-Chloronaphthalene | ND | 0.50 | 1.00 | |
| 2-Chlorophenol | ND | 0.50 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 0.50 | 1.00 | |
| Chrysene | ND | 0.50 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 0.50 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 0.50 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 0.50 | 1.00 | |
| Dibenzofuran | ND | 0.50 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 0.50 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 0.50 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 0.50 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 10 | 1.00 | |
| 2,4-Dichlorophenol | ND | 0.50 | 1.00 | |
| Diethyl Phthalate | ND | 0.50 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|----------------------------|---------------|-----------|-----------|-------------------|
| Dimethyl Phthalate | ND | 0.50 | 1.00 | |
| 2,4-Dimethylphenol | ND | 0.50 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 2.5 | 1.00 | |
| 2,4-Dinitrophenol | ND | 2.5 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 0.50 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 0.50 | 1.00 | |
| Fluoranthene | ND | 0.50 | 1.00 | |
| Fluorene | ND | 0.50 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 0.50 | 1.00 | |
| Hexachlorobenzene | ND | 0.50 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 2.5 | 1.00 | |
| Hexachloroethane | ND | 0.50 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 0.50 | 1.00 | |
| Isophorone | ND | 0.50 | 1.00 | |
| 2-Methylnaphthalene | ND | 0.50 | 1.00 | |
| 1-Methylnaphthalene | ND | 0.50 | 1.00 | |
| 2-Methylphenol | ND | 0.50 | 1.00 | |
| 3/4-Methylphenol | ND | 0.50 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 0.50 | 1.00 | |
| N-Nitrosodimethylamine | ND | 0.50 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 0.50 | 1.00 | |
| Naphthalene | ND | 0.50 | 1.00 | |
| 4-Nitroaniline | ND | 0.50 | 1.00 | |
| 3-Nitroaniline | ND | 0.50 | 1.00 | |
| 2-Nitroaniline | ND | 0.50 | 1.00 | |
| Nitrobenzene | ND | 2.5 | 1.00 | |
| 4-Nitrophenol | ND | 0.50 | 1.00 | |
| 2-Nitrophenol | ND | 0.50 | 1.00 | |
| Pentachlorophenol | ND | 2.5 | 1.00 | |
| Phenanthrene | ND | 0.50 | 1.00 | |
| Phenol | ND | 0.50 | 1.00 | |
| Pyrene | ND | 0.50 | 1.00 | |
| Pyridine | ND | 0.50 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 0.50 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 0.50 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 0.50 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 77 | 27-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorophenol | 79 | 25-120 | |
| Nitrobenzene-d5 | 66 | 33-123 | |
| p-Terphenyl-d14 | 76 | 27-159 | |
| Phenol-d6 | 79 | 26-122 | |
| 2,4,6-Tribromophenol | 90 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 1 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|------------------|-----------------|-----------------------|-------------------|
| TB07 | 15-06-0463-7-A | 06/04/15 00:00 | Aqueous | GC/MS V V | 06/13/15 | 06/13/15 20:26 | 150613L025 |

| Parameter | Result | RL | DF | Qualifiers |
|-----------------------------|--------|------|------|------------|
| Acetone | ND | 20 | 1.00 | |
| Benzene | ND | 0.50 | 1.00 | |
| Bromobenzene | ND | 1.0 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 1.00 | |
| Bromoform | ND | 1.0 | 1.00 | |
| Bromomethane | ND | 10 | 1.00 | |
| 2-Butanone | ND | 10 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 1.00 | |
| Carbon Disulfide | ND | 10 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 1.00 | |
| Chloroethane | ND | 5.0 | 1.00 | |
| Chloroform | ND | 1.0 | 1.00 | |
| Chloromethane | ND | 10 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 1.00 | |
| Dibromomethane | ND | 1.0 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 1.00 | |
| 2,2-Dichloropropane | ND | 1.0 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 2 of 4

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|-----------|-------------------|
| 1,1-Dichloropropene | ND | 1.0 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 1.00 | |
| 2-Hexanone | ND | 10 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 1.00 | |
| Methylene Chloride | ND | 10 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 1.00 | |
| Naphthalene | ND | 10 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 1.00 | |
| Styrene | ND | 1.0 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 1.00 | |
| Toluene | ND | 1.0 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 1.00 | |
| Trichloroethene | ND | 1.0 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 1.00 | |
| Vinyl Acetate | ND | 10 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 1.00 | |
| o-Xylene | ND | 1.0 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 100 | 80-120 | |
| Dibromofluoromethane | 98 | 78-126 | |
| 1,2-Dichloroethane-d4 | 95 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 3 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-17397 | N/A | Aqueous | GC/MS V V | 06/13/15 | 06/13/15 11:37 | 150613L025 |

| Parameter | Result | RL | DF | Qualifiers |
|-----------------------------|--------|------|------|------------|
| Acetone | ND | 20 | 1.00 | |
| Benzene | ND | 0.50 | 1.00 | |
| Bromobenzene | ND | 1.0 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 1.00 | |
| Bromoform | ND | 1.0 | 1.00 | |
| Bromomethane | ND | 10 | 1.00 | |
| 2-Butanone | ND | 10 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 1.00 | |
| Carbon Disulfide | ND | 10 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 1.00 | |
| Chloroethane | ND | 5.0 | 1.00 | |
| Chloroform | ND | 1.0 | 1.00 | |
| Chloromethane | ND | 10 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 1.00 | |
| Dibromomethane | ND | 1.0 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 1.00 | |
| 2,2-Dichloropropane | ND | 1.0 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 4 of 4

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|-----------|-------------------|
| 1,1-Dichloropropene | ND | 1.0 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 1.00 | |
| 2-Hexanone | ND | 10 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 1.00 | |
| Methylene Chloride | ND | 10 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 1.00 | |
| Naphthalene | ND | 10 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 1.00 | |
| Styrene | ND | 1.0 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 1.00 | |
| Toluene | ND | 1.0 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 1.00 | |
| Trichloroethene | ND | 1.0 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 1.00 | |
| Vinyl Acetate | ND | 10 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 1.00 | |
| o-Xylene | ND | 1.0 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 99 | 80-120 | |
| Dibromofluoromethane | 101 | 78-126 | |
| 1,2-Dichloroethane-d4 | 95 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

Page 1 of 14

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS310 | 15-06-0463-1-B | 06/04/15 11:22 | Solid | GC/MS W | 06/07/15 | 06/08/15 04:06 | 150607L024 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|-----------|-------------------|
| Acetone | ND | 160 | 1.00 | |
| Benzene | ND | 6.3 | 1.00 | |
| Bromobenzene | ND | 6.3 | 1.00 | |
| Bromochloromethane | ND | 6.3 | 1.00 | |
| Bromodichloromethane | ND | 6.3 | 1.00 | |
| Bromoform | ND | 6.3 | 1.00 | |
| Bromomethane | ND | 32 | 1.00 | |
| 2-Butanone | ND | 63 | 1.00 | |
| n-Butylbenzene | ND | 6.3 | 1.00 | |
| sec-Butylbenzene | ND | 6.3 | 1.00 | |
| tert-Butylbenzene | ND | 6.3 | 1.00 | |
| Carbon Disulfide | ND | 63 | 1.00 | |
| Carbon Tetrachloride | ND | 6.3 | 1.00 | |
| Chlorobenzene | ND | 6.3 | 1.00 | |
| Chloroethane | ND | 6.3 | 1.00 | |
| Chloroform | ND | 6.3 | 1.00 | |
| Chloromethane | ND | 32 | 1.00 | |
| 2-Chlorotoluene | ND | 6.3 | 1.00 | |
| 4-Chlorotoluene | ND | 6.3 | 1.00 | |
| Dibromochloromethane | ND | 6.3 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 13 | 1.00 | |
| 1,2-Dibromoethane | ND | 6.3 | 1.00 | |
| Dibromomethane | ND | 6.3 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 6.3 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 6.3 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 6.3 | 1.00 | |
| Dichlorodifluoromethane | ND | 6.3 | 1.00 | |
| 1,1-Dichloroethane | ND | 6.3 | 1.00 | |
| 1,2-Dichloroethane | ND | 6.3 | 1.00 | |
| 1,1-Dichloroethene | ND | 6.3 | 1.00 | |
| c-1,2-Dichloroethene | ND | 6.3 | 1.00 | |
| t-1,2-Dichloroethene | ND | 6.3 | 1.00 | |
| 1,2-Dichloropropane | ND | 6.3 | 1.00 | |
| 1,3-Dichloropropane | ND | 6.3 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

Page 2 of 14

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|-----------------|-----------------------|-------------------|-------------------|
| 2,2-Dichloropropane | ND | 6.3 | 1.00 | |
| 1,1-Dichloropropene | ND | 6.3 | 1.00 | |
| c-1,3-Dichloropropene | ND | 6.3 | 1.00 | |
| t-1,3-Dichloropropene | ND | 6.3 | 1.00 | |
| Ethylbenzene | ND | 6.3 | 1.00 | |
| 2-Hexanone | ND | 63 | 1.00 | |
| Isopropylbenzene | ND | 6.3 | 1.00 | |
| p-Isopropyltoluene | ND | 6.3 | 1.00 | |
| Methylene Chloride | ND | 63 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 63 | 1.00 | |
| Naphthalene | ND | 63 | 1.00 | |
| n-Propylbenzene | ND | 6.3 | 1.00 | |
| Styrene | ND | 6.3 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 6.3 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 6.3 | 1.00 | |
| Tetrachloroethene | ND | 6.3 | 1.00 | |
| Toluene | ND | 6.3 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 13 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 6.3 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 6.3 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 6.3 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 63 | 1.00 | |
| Trichloroethene | ND | 6.3 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 6.3 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 6.3 | 1.00 | |
| Trichlorofluoromethane | ND | 63 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 6.3 | 1.00 | |
| Vinyl Acetate | ND | 63 | 1.00 | |
| Vinyl Chloride | ND | 6.3 | 1.00 | |
| p/m-Xylene | ND | 6.3 | 1.00 | |
| o-Xylene | ND | 6.3 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 6.3 | 1.00 | |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | |
| 1,4-Bromofluorobenzene | 99 | 60-132 | | |
| Dibromofluoromethane | 107 | 63-141 | | |
| 1,2-Dichloroethane-d4 | 115 | 62-146 | | |
| Toluene-d8 | 99 | 80-120 | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

Page 3 of 14

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS320 | 15-06-0463-2-B | 06/04/15 11:43 | Solid | GC/MS W | 06/07/15 | 06/08/15 04:35 | 150607L024 |

Comment(s): - Results are reported on a dry weight basis.

| Parameter | Result | RL | DF | Qualifiers |
|-----------------------------|--------|-----|------|------------|
| Acetone | ND | 180 | 1.00 | |
| Benzene | ND | 7.2 | 1.00 | |
| Bromobenzene | ND | 7.2 | 1.00 | |
| Bromochloromethane | ND | 7.2 | 1.00 | |
| Bromodichloromethane | ND | 7.2 | 1.00 | |
| Bromoform | ND | 7.2 | 1.00 | |
| Bromomethane | ND | 36 | 1.00 | |
| 2-Butanone | ND | 72 | 1.00 | |
| n-Butylbenzene | ND | 7.2 | 1.00 | |
| sec-Butylbenzene | ND | 7.2 | 1.00 | |
| tert-Butylbenzene | ND | 7.2 | 1.00 | |
| Carbon Disulfide | ND | 72 | 1.00 | |
| Carbon Tetrachloride | ND | 7.2 | 1.00 | |
| Chlorobenzene | ND | 7.2 | 1.00 | |
| Chloroethane | ND | 7.2 | 1.00 | |
| Chloroform | ND | 7.2 | 1.00 | |
| Chloromethane | ND | 36 | 1.00 | |
| 2-Chlorotoluene | ND | 7.2 | 1.00 | |
| 4-Chlorotoluene | ND | 7.2 | 1.00 | |
| Dibromochloromethane | ND | 7.2 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 14 | 1.00 | |
| 1,2-Dibromoethane | ND | 7.2 | 1.00 | |
| Dibromomethane | ND | 7.2 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 7.2 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 7.2 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 7.2 | 1.00 | |
| Dichlorodifluoromethane | ND | 7.2 | 1.00 | |
| 1,1-Dichloroethane | ND | 7.2 | 1.00 | |
| 1,2-Dichloroethane | ND | 7.2 | 1.00 | |
| 1,1-Dichloroethene | ND | 7.2 | 1.00 | |
| c-1,2-Dichloroethene | ND | 7.2 | 1.00 | |
| t-1,2-Dichloroethene | ND | 7.2 | 1.00 | |
| 1,2-Dichloropropane | ND | 7.2 | 1.00 | |
| 1,3-Dichloropropane | ND | 7.2 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

Page 4 of 14

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 7.2 | 1.00 | |
| 1,1-Dichloropropene | ND | 7.2 | 1.00 | |
| c-1,3-Dichloropropene | ND | 7.2 | 1.00 | |
| t-1,3-Dichloropropene | ND | 7.2 | 1.00 | |
| Ethylbenzene | ND | 7.2 | 1.00 | |
| 2-Hexanone | ND | 72 | 1.00 | |
| Isopropylbenzene | ND | 7.2 | 1.00 | |
| p-Isopropyltoluene | ND | 7.2 | 1.00 | |
| Methylene Chloride | ND | 72 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 72 | 1.00 | |
| Naphthalene | ND | 72 | 1.00 | |
| n-Propylbenzene | ND | 7.2 | 1.00 | |
| Styrene | ND | 7.2 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 7.2 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 7.2 | 1.00 | |
| Tetrachloroethene | ND | 7.2 | 1.00 | |
| Toluene | ND | 7.2 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 14 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 7.2 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 7.2 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 7.2 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 72 | 1.00 | |
| Trichloroethene | ND | 7.2 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 7.2 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 7.2 | 1.00 | |
| Trichlorofluoromethane | ND | 72 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 7.2 | 1.00 | |
| Vinyl Acetate | ND | 72 | 1.00 | |
| Vinyl Chloride | ND | 7.2 | 1.00 | |
| p/m-Xylene | ND | 7.2 | 1.00 | |
| o-Xylene | ND | 7.2 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 7.2 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 99 | 60-132 | |
| Dibromofluoromethane | 112 | 63-141 | |
| 1,2-Dichloroethane-d4 | 120 | 62-146 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS330 | 15-06-0463-3-B | 06/04/15 12:20 | Solid | GC/MS W | 06/07/15 | 06/08/15 05:05 | 150607L024 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|-----------|-------------------|
| Acetone | ND | 180 | 1.00 | |
| Benzene | ND | 7.2 | 1.00 | |
| Bromobenzene | ND | 7.2 | 1.00 | |
| Bromochloromethane | ND | 7.2 | 1.00 | |
| Bromodichloromethane | ND | 7.2 | 1.00 | |
| Bromoform | ND | 7.2 | 1.00 | |
| Bromomethane | ND | 36 | 1.00 | |
| 2-Butanone | ND | 72 | 1.00 | |
| n-Butylbenzene | ND | 7.2 | 1.00 | |
| sec-Butylbenzene | ND | 7.2 | 1.00 | |
| tert-Butylbenzene | ND | 7.2 | 1.00 | |
| Carbon Disulfide | ND | 72 | 1.00 | |
| Carbon Tetrachloride | ND | 7.2 | 1.00 | |
| Chlorobenzene | ND | 7.2 | 1.00 | |
| Chloroethane | ND | 7.2 | 1.00 | |
| Chloroform | ND | 7.2 | 1.00 | |
| Chloromethane | ND | 36 | 1.00 | |
| 2-Chlorotoluene | ND | 7.2 | 1.00 | |
| 4-Chlorotoluene | ND | 7.2 | 1.00 | |
| Dibromochloromethane | ND | 7.2 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 14 | 1.00 | |
| 1,2-Dibromoethane | ND | 7.2 | 1.00 | |
| Dibromomethane | ND | 7.2 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 7.2 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 7.2 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 7.2 | 1.00 | |
| Dichlorodifluoromethane | ND | 7.2 | 1.00 | |
| 1,1-Dichloroethane | ND | 7.2 | 1.00 | |
| 1,2-Dichloroethane | ND | 7.2 | 1.00 | |
| 1,1-Dichloroethene | ND | 7.2 | 1.00 | |
| c-1,2-Dichloroethene | ND | 7.2 | 1.00 | |
| t-1,2-Dichloroethene | ND | 7.2 | 1.00 | |
| 1,2-Dichloropropane | ND | 7.2 | 1.00 | |
| 1,3-Dichloropropane | ND | 7.2 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

Page 6 of 14

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 7.2 | 1.00 | |
| 1,1-Dichloropropene | ND | 7.2 | 1.00 | |
| c-1,3-Dichloropropene | ND | 7.2 | 1.00 | |
| t-1,3-Dichloropropene | ND | 7.2 | 1.00 | |
| Ethylbenzene | ND | 7.2 | 1.00 | |
| 2-Hexanone | ND | 72 | 1.00 | |
| Isopropylbenzene | ND | 7.2 | 1.00 | |
| p-Isopropyltoluene | ND | 7.2 | 1.00 | |
| Methylene Chloride | ND | 72 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 72 | 1.00 | |
| Naphthalene | ND | 72 | 1.00 | |
| n-Propylbenzene | ND | 7.2 | 1.00 | |
| Styrene | ND | 7.2 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 7.2 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 7.2 | 1.00 | |
| Tetrachloroethene | ND | 7.2 | 1.00 | |
| Toluene | ND | 7.2 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 14 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 7.2 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 7.2 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 7.2 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 72 | 1.00 | |
| Trichloroethene | ND | 7.2 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 7.2 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 7.2 | 1.00 | |
| Trichlorofluoromethane | ND | 72 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 7.2 | 1.00 | |
| Vinyl Acetate | ND | 72 | 1.00 | |
| Vinyl Chloride | ND | 7.2 | 1.00 | |
| p/m-Xylene | ND | 7.2 | 1.00 | |
| o-Xylene | ND | 7.2 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 7.2 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 96 | 60-132 | |
| Dibromofluoromethane | 106 | 63-141 | |
| 1,2-Dichloroethane-d4 | 118 | 62-146 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS340 | 15-06-0463-4-B | 06/04/15 13:57 | Solid | GC/MS W | 06/07/15 | 06/08/15 05:34 | 150607L024 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|-----------|-------------------|
| Acetone | ND | 280 | 1.00 | |
| Benzene | ND | 11 | 1.00 | |
| Bromobenzene | ND | 11 | 1.00 | |
| Bromochloromethane | ND | 11 | 1.00 | |
| Bromodichloromethane | ND | 11 | 1.00 | |
| Bromoform | ND | 11 | 1.00 | |
| Bromomethane | ND | 57 | 1.00 | |
| 2-Butanone | ND | 110 | 1.00 | |
| n-Butylbenzene | ND | 11 | 1.00 | |
| sec-Butylbenzene | ND | 11 | 1.00 | |
| tert-Butylbenzene | ND | 11 | 1.00 | |
| Carbon Disulfide | ND | 110 | 1.00 | |
| Carbon Tetrachloride | ND | 11 | 1.00 | |
| Chlorobenzene | ND | 11 | 1.00 | |
| Chloroethane | ND | 11 | 1.00 | |
| Chloroform | ND | 11 | 1.00 | |
| Chloromethane | ND | 57 | 1.00 | |
| 2-Chlorotoluene | ND | 11 | 1.00 | |
| 4-Chlorotoluene | ND | 11 | 1.00 | |
| Dibromochloromethane | ND | 11 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 23 | 1.00 | |
| 1,2-Dibromoethane | ND | 11 | 1.00 | |
| Dibromomethane | ND | 11 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 11 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 11 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 11 | 1.00 | |
| Dichlorodifluoromethane | ND | 11 | 1.00 | |
| 1,1-Dichloroethane | ND | 11 | 1.00 | |
| 1,2-Dichloroethane | ND | 11 | 1.00 | |
| 1,1-Dichloroethene | ND | 11 | 1.00 | |
| c-1,2-Dichloroethene | ND | 11 | 1.00 | |
| t-1,2-Dichloroethene | ND | 11 | 1.00 | |
| 1,2-Dichloropropane | ND | 11 | 1.00 | |
| 1,3-Dichloropropane | ND | 11 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

Page 8 of 14

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 11 | 1.00 | |
| 1,1-Dichloropropene | ND | 11 | 1.00 | |
| c-1,3-Dichloropropene | ND | 11 | 1.00 | |
| t-1,3-Dichloropropene | ND | 11 | 1.00 | |
| Ethylbenzene | ND | 11 | 1.00 | |
| 2-Hexanone | ND | 110 | 1.00 | |
| Isopropylbenzene | ND | 11 | 1.00 | |
| p-Isopropyltoluene | ND | 11 | 1.00 | |
| Methylene Chloride | ND | 110 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 110 | 1.00 | |
| Naphthalene | ND | 110 | 1.00 | |
| n-Propylbenzene | ND | 11 | 1.00 | |
| Styrene | ND | 11 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 11 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 11 | 1.00 | |
| Tetrachloroethene | ND | 11 | 1.00 | |
| Toluene | ND | 11 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 23 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 11 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 11 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 11 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 110 | 1.00 | |
| Trichloroethene | ND | 11 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 11 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 11 | 1.00 | |
| Trichlorofluoromethane | ND | 110 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 11 | 1.00 | |
| Vinyl Acetate | ND | 110 | 1.00 | |
| Vinyl Chloride | ND | 11 | 1.00 | |
| p/m-Xylene | ND | 11 | 1.00 | |
| o-Xylene | ND | 11 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 11 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 97 | 60-132 | |
| Dibromofluoromethane | 110 | 63-141 | |
| 1,2-Dichloroethane-d4 | 121 | 62-146 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HCS360 | 15-06-0463-5-B | 06/04/15 14:24 | Solid | GC/MS W | 06/07/15 | 06/08/15 06:03 | 150607L024 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|-----------|-------------------|
| Acetone | ND | 260 | 1.00 | |
| Benzene | ND | 10 | 1.00 | |
| Bromobenzene | ND | 10 | 1.00 | |
| Bromochloromethane | ND | 10 | 1.00 | |
| Bromodichloromethane | ND | 10 | 1.00 | |
| Bromoform | ND | 10 | 1.00 | |
| Bromomethane | ND | 52 | 1.00 | |
| 2-Butanone | ND | 100 | 1.00 | |
| n-Butylbenzene | ND | 10 | 1.00 | |
| sec-Butylbenzene | ND | 10 | 1.00 | |
| tert-Butylbenzene | ND | 10 | 1.00 | |
| Carbon Disulfide | ND | 100 | 1.00 | |
| Carbon Tetrachloride | ND | 10 | 1.00 | |
| Chlorobenzene | ND | 10 | 1.00 | |
| Chloroethane | ND | 10 | 1.00 | |
| Chloroform | ND | 10 | 1.00 | |
| Chloromethane | ND | 52 | 1.00 | |
| 2-Chlorotoluene | ND | 10 | 1.00 | |
| 4-Chlorotoluene | ND | 10 | 1.00 | |
| Dibromochloromethane | ND | 10 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 21 | 1.00 | |
| 1,2-Dibromoethane | ND | 10 | 1.00 | |
| Dibromomethane | ND | 10 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 1.00 | |
| Dichlorodifluoromethane | ND | 10 | 1.00 | |
| 1,1-Dichloroethane | ND | 10 | 1.00 | |
| 1,2-Dichloroethane | ND | 10 | 1.00 | |
| 1,1-Dichloroethene | ND | 10 | 1.00 | |
| c-1,2-Dichloroethene | ND | 10 | 1.00 | |
| t-1,2-Dichloroethene | ND | 10 | 1.00 | |
| 1,2-Dichloropropane | ND | 10 | 1.00 | |
| 1,3-Dichloropropane | ND | 10 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

Page 10 of 14

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 10 | 1.00 | |
| 1,1-Dichloropropene | ND | 10 | 1.00 | |
| c-1,3-Dichloropropene | ND | 10 | 1.00 | |
| t-1,3-Dichloropropene | ND | 10 | 1.00 | |
| Ethylbenzene | ND | 10 | 1.00 | |
| 2-Hexanone | ND | 100 | 1.00 | |
| Isopropylbenzene | ND | 10 | 1.00 | |
| p-Isopropyltoluene | ND | 10 | 1.00 | |
| Methylene Chloride | ND | 100 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 100 | 1.00 | |
| Naphthalene | ND | 100 | 1.00 | |
| n-Propylbenzene | ND | 10 | 1.00 | |
| Styrene | ND | 10 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 10 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 10 | 1.00 | |
| Tetrachloroethene | ND | 10 | 1.00 | |
| Toluene | ND | 10 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 21 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 10 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 10 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 100 | 1.00 | |
| Trichloroethene | ND | 10 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 10 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 10 | 1.00 | |
| Trichlorofluoromethane | ND | 100 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 10 | 1.00 | |
| Vinyl Acetate | ND | 100 | 1.00 | |
| Vinyl Chloride | ND | 10 | 1.00 | |
| p/m-Xylene | ND | 10 | 1.00 | |
| o-Xylene | ND | 10 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 10 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 60-132 | |
| Dibromofluoromethane | 112 | 63-141 | |
| 1,2-Dichloroethane-d4 | 124 | 62-146 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

Page 11 of 14

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|--------------|----------------|-----------------|---------------------------|-------------------|
| FDHCS360 | 15-06-0463-6-B | 06/04/15 14:24 | Solid | GC/MS W | 06/07/15 | 06/08/15 06:33 | 150607L024 |

Comment(s): - Results are reported on a dry weight basis.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|-----------|-------------------|
| Acetone | ND | 260 | 1.00 | |
| Benzene | ND | 10 | 1.00 | |
| Bromobenzene | ND | 10 | 1.00 | |
| Bromochloromethane | ND | 10 | 1.00 | |
| Bromodichloromethane | ND | 10 | 1.00 | |
| Bromoform | ND | 10 | 1.00 | |
| Bromomethane | ND | 51 | 1.00 | |
| 2-Butanone | ND | 100 | 1.00 | |
| n-Butylbenzene | ND | 10 | 1.00 | |
| sec-Butylbenzene | ND | 10 | 1.00 | |
| tert-Butylbenzene | ND | 10 | 1.00 | |
| Carbon Disulfide | ND | 100 | 1.00 | |
| Carbon Tetrachloride | ND | 10 | 1.00 | |
| Chlorobenzene | ND | 10 | 1.00 | |
| Chloroethane | ND | 10 | 1.00 | |
| Chloroform | ND | 10 | 1.00 | |
| Chloromethane | ND | 51 | 1.00 | |
| 2-Chlorotoluene | ND | 10 | 1.00 | |
| 4-Chlorotoluene | ND | 10 | 1.00 | |
| Dibromochloromethane | ND | 10 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 21 | 1.00 | |
| 1,2-Dibromoethane | ND | 10 | 1.00 | |
| Dibromomethane | ND | 10 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 1.00 | |
| Dichlorodifluoromethane | ND | 10 | 1.00 | |
| 1,1-Dichloroethane | ND | 10 | 1.00 | |
| 1,2-Dichloroethane | ND | 10 | 1.00 | |
| 1,1-Dichloroethene | ND | 10 | 1.00 | |
| c-1,2-Dichloroethene | ND | 10 | 1.00 | |
| t-1,2-Dichloroethene | ND | 10 | 1.00 | |
| 1,2-Dichloropropane | ND | 10 | 1.00 | |
| 1,3-Dichloropropane | ND | 10 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

Page 12 of 14

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 10 | 1.00 | |
| 1,1-Dichloropropene | ND | 10 | 1.00 | |
| c-1,3-Dichloropropene | ND | 10 | 1.00 | |
| t-1,3-Dichloropropene | ND | 10 | 1.00 | |
| Ethylbenzene | ND | 10 | 1.00 | |
| 2-Hexanone | ND | 100 | 1.00 | |
| Isopropylbenzene | ND | 10 | 1.00 | |
| p-Isopropyltoluene | ND | 10 | 1.00 | |
| Methylene Chloride | ND | 100 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 100 | 1.00 | |
| Naphthalene | ND | 100 | 1.00 | |
| n-Propylbenzene | ND | 10 | 1.00 | |
| Styrene | ND | 10 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 10 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 10 | 1.00 | |
| Tetrachloroethene | ND | 10 | 1.00 | |
| Toluene | ND | 10 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 21 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 10 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 10 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 100 | 1.00 | |
| Trichloroethene | ND | 10 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 10 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 10 | 1.00 | |
| Trichlorofluoromethane | ND | 100 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 10 | 1.00 | |
| Vinyl Acetate | ND | 100 | 1.00 | |
| Vinyl Chloride | ND | 10 | 1.00 | |
| p/m-Xylene | ND | 10 | 1.00 | |
| o-Xylene | ND | 10 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 10 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 60-132 | |
| Dibromofluoromethane | 109 | 63-141 | |
| 1,2-Dichloroethane-d4 | 119 | 62-146 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-796-9798 | N/A | Solid | GC/MS W | 06/07/15 | 06/08/15 00:42 | 150607L024 |

| Parameter | Result | RL | DF | Qualifiers |
|-----------------------------|--------|-----|------|------------|
| Acetone | ND | 120 | 1.00 | |
| Benzene | ND | 5.0 | 1.00 | |
| Bromobenzene | ND | 5.0 | 1.00 | |
| Bromochloromethane | ND | 5.0 | 1.00 | |
| Bromodichloromethane | ND | 5.0 | 1.00 | |
| Bromoform | ND | 5.0 | 1.00 | |
| Bromomethane | ND | 25 | 1.00 | |
| 2-Butanone | ND | 50 | 1.00 | |
| n-Butylbenzene | ND | 5.0 | 1.00 | |
| sec-Butylbenzene | ND | 5.0 | 1.00 | |
| tert-Butylbenzene | ND | 5.0 | 1.00 | |
| Carbon Disulfide | ND | 50 | 1.00 | |
| Carbon Tetrachloride | ND | 5.0 | 1.00 | |
| Chlorobenzene | ND | 5.0 | 1.00 | |
| Chloroethane | ND | 5.0 | 1.00 | |
| Chloroform | ND | 5.0 | 1.00 | |
| Chloromethane | ND | 25 | 1.00 | |
| 2-Chlorotoluene | ND | 5.0 | 1.00 | |
| 4-Chlorotoluene | ND | 5.0 | 1.00 | |
| Dibromochloromethane | ND | 5.0 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 10 | 1.00 | |
| 1,2-Dibromoethane | ND | 5.0 | 1.00 | |
| Dibromomethane | ND | 5.0 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 5.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 5.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 5.0 | 1.00 | |
| Dichlorodifluoromethane | ND | 5.0 | 1.00 | |
| 1,1-Dichloroethane | ND | 5.0 | 1.00 | |
| 1,2-Dichloroethane | ND | 5.0 | 1.00 | |
| 1,1-Dichloroethene | ND | 5.0 | 1.00 | |
| c-1,2-Dichloroethene | ND | 5.0 | 1.00 | |
| t-1,2-Dichloroethene | ND | 5.0 | 1.00 | |
| 1,2-Dichloropropane | ND | 5.0 | 1.00 | |
| 1,3-Dichloropropane | ND | 5.0 | 1.00 | |
| 2,2-Dichloropropane | ND | 5.0 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

Page 14 of 14

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|-----------|-------------------|
| 1,1-Dichloropropene | ND | 5.0 | 1.00 | |
| c-1,3-Dichloropropene | ND | 5.0 | 1.00 | |
| t-1,3-Dichloropropene | ND | 5.0 | 1.00 | |
| Ethylbenzene | ND | 5.0 | 1.00 | |
| 2-Hexanone | ND | 50 | 1.00 | |
| Isopropylbenzene | ND | 5.0 | 1.00 | |
| p-Isopropyltoluene | ND | 5.0 | 1.00 | |
| Methylene Chloride | ND | 50 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 50 | 1.00 | |
| Naphthalene | ND | 50 | 1.00 | |
| n-Propylbenzene | ND | 5.0 | 1.00 | |
| Styrene | ND | 5.0 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 5.0 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 5.0 | 1.00 | |
| Tetrachloroethene | ND | 5.0 | 1.00 | |
| Toluene | ND | 5.0 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 10 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 5.0 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 5.0 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 5.0 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 50 | 1.00 | |
| Trichloroethene | ND | 5.0 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 5.0 | 1.00 | |
| Trichlorofluoromethane | ND | 50 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 5.0 | 1.00 | |
| Vinyl Acetate | ND | 50 | 1.00 | |
| Vinyl Chloride | ND | 5.0 | 1.00 | |
| p/m-Xylene | ND | 5.0 | 1.00 | |
| o-Xylene | ND | 5.0 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 5.0 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 60-132 | |
| Dibromofluoromethane | 105 | 63-141 | |
| 1,2-Dichloroethane-d4 | 114 | 62-146 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

06/05/15

Work Order:

15-06-0463

Page 1 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|--------------|
| HCS310 | 15-06-0463-1 | 06/04/15 11:22 | Solid |

Comment(s): (9) - Results are reported on a dry weight basis.

| Parameter | Results | RL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|---|---------|-------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 19.3 | 0.100 | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 7.82 | 0.01 | 1.00 | | pH units | 06/05/15 | 06/05/15 | EPA 9045D |
| Carbon, Total Organic (9) | 19000 | 620 | 1.00 | | mg/kg | 06/17/15 | 06/17/15 | EPA 9060A |
| Alkalinity, Total (as CaCO ₃) | 440 | 5.0 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO ₃) | 440 | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO ₃) | ND | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved | 4800 | 10.0 | 1.00 | | mg/kg | 06/12/15 | 06/12/15 | SM 2540 C (M) |
| Phosphorus, Total (9) | 2.8 | 0.62 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

| | | | |
|---------------|---------------------|-----------------------|--------------|
| HCS320 | 15-06-0463-2 | 06/04/15 11:43 | Solid |
|---------------|---------------------|-----------------------|--------------|

Comment(s): (9) - Results are reported on a dry weight basis.

| Parameter | Results | RL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|---|---------|-------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 29.1 | 0.100 | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 8.12 | 0.01 | 1.00 | | pH units | 06/05/15 | 06/05/15 | EPA 9045D |
| Carbon, Total Organic (9) | 15000 | 710 | 1.00 | | mg/kg | 06/17/15 | 06/17/15 | EPA 9060A |
| Alkalinity, Total (as CaCO ₃) | 340 | 5.0 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO ₃) | 340 | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO ₃) | ND | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved | 5930 | 10.0 | 1.00 | | mg/kg | 06/12/15 | 06/12/15 | SM 2540 C (M) |
| Phosphorus, Total (9) | 4.3 | 0.71 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

| | | | |
|---------------|---------------------|-----------------------|--------------|
| HCS330 | 15-06-0463-3 | 06/04/15 12:20 | Solid |
|---------------|---------------------|-----------------------|--------------|

Comment(s): (9) - Results are reported on a dry weight basis.

| Parameter | Results | RL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|---|---------|-------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 30.6 | 0.100 | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 7.68 | 0.01 | 1.00 | | pH units | 06/05/15 | 06/05/15 | EPA 9045D |
| Carbon, Total Organic (9) | 15000 | 720 | 1.00 | | mg/kg | 06/17/15 | 06/17/15 | EPA 9060A |
| Alkalinity, Total (as CaCO ₃) | 670 | 5.0 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO ₃) | 670 | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO ₃) | ND | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved | 17900 | 100 | 1.00 | | mg/kg | 06/12/15 | 06/12/15 | SM 2540 C (M) |
| Phosphorus, Total (9) | 2.5 | 0.72 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

06/05/15

Work Order:

15-06-0463

Page 2 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|--------------|
| HCS340 | 15-06-0463-4 | 06/04/15 13:57 | Solid |

Comment(s): (9) - Results are reported on a dry weight basis.

| Parameter | Results | RL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|---|---------|-------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 55.7 | 0.100 | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 7.58 | 0.01 | 1.00 | | pH units | 06/05/15 | 06/05/15 | EPA 9045D |
| Carbon, Total Organic (9) | 51000 | 1100 | 1.00 | | mg/kg | 06/17/15 | 06/17/15 | EPA 9060A |
| Alkalinity, Total (as CaCO ₃) | 580 | 5.0 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO ₃) | 580 | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO ₃) | ND | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved | 4160 | 10.0 | 1.00 | | mg/kg | 06/12/15 | 06/12/15 | SM 2540 C (M) |
| Phosphorus, Total (9) | 4.8 | 2.3 | 2.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

| | | | |
|---------------|---------------------|-----------------------|--------------|
| HCS360 | 15-06-0463-5 | 06/04/15 14:24 | Solid |
|---------------|---------------------|-----------------------|--------------|

Comment(s): (9) - Results are reported on a dry weight basis.

| Parameter | Results | RL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|---|---------|-------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 51.1 | 0.100 | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 7.35 | 0.01 | 1.00 | | pH units | 06/05/15 | 06/05/15 | EPA 9045D |
| Carbon, Total Organic (9) | 53000 | 1000 | 1.00 | | mg/kg | 06/17/15 | 06/17/15 | EPA 9060A |
| Alkalinity, Total (as CaCO ₃) | 580 | 5.0 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO ₃) | 580 | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO ₃) | ND | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved | 3460 | 10.0 | 1.00 | | mg/kg | 06/12/15 | 06/12/15 | SM 2540 C (M) |
| Phosphorus, Total (9) | 5.6 | 1.0 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

| | | | |
|-----------------|---------------------|-----------------------|--------------|
| FDHCS360 | 15-06-0463-6 | 06/04/15 14:24 | Solid |
|-----------------|---------------------|-----------------------|--------------|

Comment(s): (9) - Results are reported on a dry weight basis.

| Parameter | Results | RL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|---|---------|-------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 51.2 | 0.100 | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 7.41 | 0.01 | 1.00 | | pH units | 06/05/15 | 06/05/15 | EPA 9045D |
| Carbon, Total Organic (9) | 50000 | 1000 | 1.00 | | mg/kg | 06/17/15 | 06/17/15 | EPA 9060A |
| Alkalinity, Total (as CaCO ₃) | 590 | 5.0 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO ₃) | 590 | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO ₃) | ND | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved | 4320 | 10.0 | 1.00 | | mg/kg | 06/12/15 | 06/12/15 | SM 2540 C (M) |
| Phosphorus, Total (9) | 5.8 | 1.0 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618
Project: EAA 27122

Date Received: 06/05/15
Work Order: 15-06-0463

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| Client Sample Number | | Lab Sample Number | | | Date/Time Collected | | Matrix | |
|---|----------------|-------------------|-----------|-------------------|---------------------|----------------------|----------------------|-------------------|
| Method Blank | | N/A | | | N/A | | Solid | |
| <u>Parameter</u> | <u>Results</u> | <u>RL</u> | <u>DF</u> | <u>Qualifiers</u> | <u>Units</u> | <u>Date Prepared</u> | <u>Date Analyzed</u> | <u>Method</u> |
| Moisture | ND | 0.100 | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| Carbon, Total Organic | ND | 500 | 1.00 | | mg/kg | 06/17/15 | 06/17/15 | EPA 9060A |
| Alkalinity, Total (as CaCO ₃) | ND | 5.0 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO ₃) | ND | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO ₃) | ND | 5.0 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved | ND | 1.0 | 1.00 | | mg/kg | 06/12/15 | 06/12/15 | SM 2540 C (M) |
| Phosphorus, Total | ND | 0.50 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

Page 1 of 14

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-06-0761-7 | Sample | Solid | IC 10 | 06/10/15 | 06/11/15 04:43 | 150610S01P | | | | |
| 15-06-0761-7 | Matrix Spike | Solid | IC 10 | 06/10/15 | 06/11/15 05:49 | 150610S01P | | | | |
| 15-06-0761-7 | Matrix Spike Duplicate | Solid | IC 10 | 06/10/15 | 06/11/15 06:05 | 150610S01P | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Fluoride | 3.727 | 25.00 | 23.22 | 78 | 22.05 | 73 | 80-120 | 5 | 0-20 | 3 |
| Chloride | 148.7 | 500.0 | 687.6 | 108 | 679.1 | 106 | 80-120 | 1 | 0-20 | |
| Bromide | ND | 50.00 | 51.43 | 103 | 51.11 | 102 | 80-120 | 1 | 0-20 | |
| Nitrate (as N) | 3.879 | 50.00 | 56.64 | 106 | 56.02 | 104 | 80-120 | 1 | 0-20 | |
| Sulfate | 96.78 | 500.0 | 623.2 | 105 | 596.3 | 100 | 80-120 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: EPA 9060A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|--------|------------|---------------|----------------|---------------------|
| 15-06-0792-1 | Sample | Solid | TOC 5 | 06/17/15 | 06/17/15 13:24 | F0617TOCS1 |
| 15-06-0792-1 | Matrix Spike | Solid | TOC 5 | 06/17/15 | 06/17/15 13:24 | F0617TOCS1 |
| 15-06-0792-1 | Matrix Spike Duplicate | Solid | TOC 5 | 06/17/15 | 06/17/15 13:24 | F0617TOCS1 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Carbon, Total Organic | 600.0 | 30000 | 23400 | 76 | 26500 | 86 | 75-125 | 12 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: SM 4500 P B/E (M)

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|--------|------------|---------------|----------------|---------------------|
| 15-06-0567-2 | Sample | Solid | UV 7 | 06/10/15 | 06/11/15 16:26 | F0611TPS2 |
| 15-06-0567-2 | Matrix Spike | Solid | UV 7 | 06/10/15 | 06/11/15 16:26 | F0611TPS2 |
| 15-06-0567-2 | Matrix Spike Duplicate | Solid | UV 7 | 06/10/15 | 06/11/15 16:26 | F0611TPS2 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | 39.46 | 40.00 | 80.30 | 102 | 80.70 | 103 | 70-130 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|--------|------------|---------------|----------------|---------------------|
| 15-06-0761-7 | Sample | Solid | ICP 7300 | 06/10/15 | 06/11/15 16:44 | 150610S04 |
| 15-06-0761-7 | Matrix Spike | Solid | ICP 7300 | 06/10/15 | 06/11/15 16:41 | 150610S04 |
| 15-06-0761-7 | Matrix Spike Duplicate | Solid | ICP 7300 | 06/10/15 | 06/11/15 16:42 | 150610S04 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Calcium | 3794 | 25.00 | 4184 | 4X | 4125 | 4X | 75-125 | 4X | 0-20 | Q |
| Magnesium | 4342 | 25.00 | 4604 | 4X | 4452 | 4X | 75-125 | 4X | 0-20 | Q |
| Potassium | 3381 | 250.0 | 3940 | 4X | 3785 | 4X | 75-125 | 4X | 0-20 | Q |
| Sodium | 192.4 | 250.0 | 460.1 | 107 | 453.1 | 104 | 75-125 | 2 | 0-20 | |
| Strontium | 35.22 | 25.00 | 64.81 | 118 | 63.58 | 113 | 75-125 | 2 | 0-20 | |
| Silicon | 563.3 | 25.00 | 1145 | 4X | 1253 | 4X | 75-125 | 4X | 0-20 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-06-0498-3 | Sample | Solid | ICP/MS 04 | 06/08/15 | 06/09/15 13:52 | 150608S01 | | | | |
| 15-06-0498-3 | Matrix Spike | Solid | ICP/MS 04 | 06/08/15 | 06/09/15 13:37 | 150608S01 | | | | |
| 15-06-0498-3 | Matrix Spike Duplicate | Solid | ICP/MS 04 | 06/08/15 | 06/09/15 13:41 | 150608S01 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 25.00 | 7.402 | 30 | 7.364 | 29 | 1-97 | 1 | 0-39 | |
| Arsenic | 7.643 | 25.00 | 31.89 | 97 | 32.62 | 100 | 72-132 | 2 | 0-13 | |
| Barium | 199.5 | 25.00 | 253.0 | 4X | 251.6 | 4X | 50-152 | 4X | 0-41 | Q |
| Beryllium | ND | 25.00 | 27.99 | 112 | 28.49 | 114 | 61-121 | 2 | 0-13 | |
| Cadmium | 3.397 | 25.00 | 30.51 | 108 | 30.54 | 109 | 85-121 | 0 | 0-12 | |
| Chromium | 26.68 | 25.00 | 56.30 | 118 | 56.01 | 117 | 20-182 | 1 | 0-15 | |
| Copper | 29.47 | 25.00 | 56.49 | 108 | 56.44 | 108 | 25-157 | 0 | 0-22 | |
| Lead | 8.765 | 25.00 | 34.60 | 103 | 34.11 | 101 | 62-134 | 1 | 0-23 | |
| Nickel | 48.32 | 25.00 | 75.40 | 108 | 75.29 | 108 | 46-154 | 0 | 0-15 | |
| Selenium | ND | 25.00 | 24.77 | 99 | 25.16 | 101 | 54-132 | 2 | 0-14 | |
| Silver | ND | 12.50 | 13.70 | 110 | 13.80 | 110 | 78-126 | 1 | 0-15 | |
| Thallium | ND | 25.00 | 23.93 | 96 | 23.94 | 96 | 79-115 | 0 | 0-11 | |
| Zinc | 92.30 | 25.00 | 123.9 | 127 | 125.9 | 134 | 23-173 | 2 | 0-18 | |
| Aluminum | 11780 | 25.00 | 14080 | 4X | 13640 | 4X | 80-120 | 4X | 0-20 | Q |
| Iron | 25280 | 25.00 | 26930 | 4X | 26960 | 4X | 80-120 | 4X | 0-20 | Q |
| Manganese | 332.0 | 25.00 | 368.5 | 4X | 366.9 | 4X | 80-120 | 4X | 0-20 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 7471A Total
Method: EPA 7471A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| FDHCS360 | Sample | Solid | Mercury 05 | 06/16/15 | 06/16/15 14:32 | 150616S01 | | | | |
| FDHCS360 | Matrix Spike | Solid | Mercury 05 | 06/16/15 | 06/16/15 14:34 | 150616S01 | | | | |
| FDHCS360 | Matrix Spike Duplicate | Solid | Mercury 05 | 06/16/15 | 06/16/15 14:36 | 150616S01 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Mercury | ND | 0.8350 | 0.8818 | 106 | 0.8495 | 102 | 71-137 | 4 | 0-14 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-06-0976-4 | Sample | Solid | GC 44 | 06/13/15 | 06/16/15 19:48 | 150613S05 | | | | |
| 15-06-0976-4 | Matrix Spike | Solid | GC 44 | 06/13/15 | 06/16/15 17:54 | 150613S05 | | | | |
| 15-06-0976-4 | Matrix Spike Duplicate | Solid | GC 44 | 06/13/15 | 06/16/15 18:08 | 150613S05 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Aldrin | ND | 25.00 | 19.75 | 79 | 19.77 | 79 | 50-135 | 0 | 0-25 | |
| Alpha-BHC | ND | 25.00 | 20.42 | 82 | 20.49 | 82 | 50-135 | 0 | 0-25 | |
| Beta-BHC | ND | 25.00 | 23.08 | 92 | 22.51 | 90 | 50-135 | 3 | 0-25 | |
| 4,4'-DDD | ND | 25.00 | 24.93 | 100 | 24.93 | 100 | 50-135 | 0 | 0-25 | |
| 4,4'-DDE | ND | 25.00 | 25.33 | 101 | 24.72 | 99 | 50-135 | 2 | 0-25 | |
| 4,4'-DDT | ND | 25.00 | 25.25 | 101 | 25.37 | 101 | 50-135 | 0 | 0-25 | |
| Delta-BHC | ND | 25.00 | 24.11 | 96 | 22.99 | 92 | 50-135 | 5 | 0-25 | |
| Dieldrin | ND | 25.00 | 22.74 | 91 | 22.61 | 90 | 50-135 | 1 | 0-25 | |
| Endosulfan I | ND | 25.00 | 20.87 | 83 | 20.52 | 82 | 50-135 | 2 | 0-25 | |
| Endosulfan II | ND | 25.00 | 23.51 | 94 | 23.35 | 93 | 50-135 | 1 | 0-25 | |
| Endosulfan Sulfate | ND | 25.00 | 24.35 | 97 | 23.41 | 94 | 50-135 | 4 | 0-25 | |
| Endrin | ND | 25.00 | 25.88 | 104 | 25.79 | 103 | 50-135 | 0 | 0-25 | |
| Endrin Aldehyde | ND | 25.00 | 19.93 | 80 | 20.21 | 81 | 50-135 | 1 | 0-25 | |
| Gamma-BHC | ND | 25.00 | 22.06 | 88 | 21.76 | 87 | 50-135 | 1 | 0-25 | |
| Heptachlor | ND | 25.00 | 22.81 | 91 | 21.42 | 86 | 50-135 | 6 | 0-25 | |
| Heptachlor Epoxide | ND | 25.00 | 21.42 | 86 | 20.83 | 83 | 50-135 | 3 | 0-25 | |
| Methoxychlor | ND | 25.00 | 27.39 | 110 | 27.84 | 111 | 50-135 | 2 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|--------|------------|---------------|----------------|---------------------|
| 15-06-0976-4 | Sample | Solid | GC 31 | 06/13/15 | 06/16/15 14:13 | 150613S06 |
| 15-06-0976-4 | Matrix Spike | Solid | GC 31 | 06/13/15 | 06/16/15 12:22 | 150613S06 |
| 15-06-0976-4 | Matrix Spike Duplicate | Solid | GC 31 | 06/13/15 | 06/16/15 12:41 | 150613S06 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|--------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Aroclor-1016 | ND | 100.0 | 88.10 | 88 | 85.38 | 85 | 50-135 | 3 | 0-20 | |
| Aroclor-1260 | ND | 100.0 | 98.71 | 99 | 95.09 | 95 | 50-135 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-06-0559-2 | Sample | Solid | GC 26 | 06/08/15 | 06/09/15 15:22 | 150608S12 | | | | |
| 15-06-0559-2 | Matrix Spike | Solid | GC 26 | 06/08/15 | 06/09/15 13:10 | 150608S12 | | | | |
| 15-06-0559-2 | Matrix Spike Duplicate | Solid | GC 26 | 06/08/15 | 06/09/15 13:54 | 150608S12 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | ND | 4.000 | 4.590 | 115 | 4.846 | 121 | 30-130 | 5 | 0-30 | |
| Bolstar | ND | 4.000 | 4.650 | 116 | 4.736 | 118 | 30-130 | 2 | 0-30 | |
| Chlorpyrifos | ND | 4.000 | 4.557 | 114 | 4.797 | 120 | 30-130 | 5 | 0-30 | |
| Coumaphos | ND | 4.000 | 4.506 | 113 | 4.732 | 118 | 30-130 | 5 | 0-30 | |
| Diazinon | ND | 4.000 | 3.212 | 80 | 3.627 | 91 | 30-130 | 12 | 0-30 | |
| Disulfoton | ND | 4.000 | 4.549 | 114 | 4.551 | 114 | 30-130 | 0 | 0-30 | |
| Ethoprop | ND | 4.000 | 3.455 | 86 | 3.862 | 97 | 30-130 | 11 | 0-30 | |
| Fensulfothion | ND | 4.000 | 2.024 | 51 | 2.574 | 64 | 30-130 | 24 | 0-30 | |
| Fenthion | ND | 4.000 | 4.671 | 117 | 4.800 | 120 | 30-130 | 3 | 0-30 | |
| Merphos | ND | 4.000 | 4.906 | 123 | 4.949 | 124 | 30-130 | 1 | 0-30 | |
| Methyl Parathion | ND | 4.000 | 4.468 | 112 | 4.479 | 112 | 30-130 | 0 | 0-30 | |
| Phorate | ND | 4.000 | 4.502 | 113 | 4.450 | 111 | 30-130 | 1 | 0-30 | |
| Ronnel | ND | 4.000 | 4.413 | 110 | 4.453 | 111 | 30-130 | 1 | 0-30 | |
| Stirophos | ND | 4.000 | 2.909 | 73 | 3.357 | 84 | 30-130 | 14 | 0-30 | |
| Tokuthion | ND | 4.000 | 4.687 | 117 | 4.860 | 122 | 30-130 | 4 | 0-30 | |
| Trichloronate | ND | 4.000 | 4.787 | 120 | 4.899 | 122 | 30-130 | 2 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|--------|------------|---------------|----------------|---------------------|
| 15-06-0559-4 | Sample | Solid | GC 40 | 06/10/15 | 06/11/15 21:39 | 150610S09 |
| 15-06-0559-4 | Matrix Spike | Solid | GC 40 | 06/10/15 | 06/11/15 19:44 | 150610S09 |
| 15-06-0559-4 | Matrix Spike Duplicate | Solid | GC 40 | 06/10/15 | 06/11/15 20:07 | 150610S09 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| 2,4-D | ND | 400.0 | 259.5 | 65 | 263.7 | 66 | 30-130 | 2 | 0-30 | |
| 2,4,5-T | ND | 40.00 | 26.20 | 66 | 26.50 | 66 | 30-130 | 1 | 0-30 | |
| 2,4-DB | ND | 400.0 | 308.8 | 77 | 284.7 | 71 | 30-130 | 8 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-06-0929-6 | Sample | Solid | GC/MS SS | 06/13/15 | 06/15/15 22:33 | 150613S07 | | | | |
| 15-06-0929-6 | Matrix Spike | Solid | GC/MS SS | 06/13/15 | 06/15/15 21:54 | 150613S07 | | | | |
| 15-06-0929-6 | Matrix Spike Duplicate | Solid | GC/MS SS | 06/13/15 | 06/15/15 22:13 | 150613S07 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | ND | 10.00 | 6.707 | 67 | 6.505 | 65 | 34-148 | 3 | 0-20 | |
| Acenaphthylene | ND | 10.00 | 6.711 | 67 | 6.521 | 65 | 53-120 | 3 | 0-20 | |
| Butyl Benzyl Phthalate | ND | 10.00 | 6.733 | 67 | 6.630 | 66 | 15-189 | 2 | 0-20 | |
| 4-Chloro-3-Methylphenol | ND | 10.00 | 6.086 | 61 | 6.019 | 60 | 32-120 | 1 | 0-20 | |
| 2-Chlorophenol | ND | 10.00 | 5.673 | 57 | 5.539 | 55 | 53-120 | 2 | 0-20 | |
| 1,4-Dichlorobenzene | ND | 10.00 | 4.225 | 42 | 3.945 | 39 | 43-120 | 7 | 0-26 | 3 |
| Dimethyl Phthalate | ND | 10.00 | 6.540 | 65 | 6.478 | 65 | 44-122 | 1 | 0-20 | |
| 2,4-Dinitrotoluene | ND | 10.00 | 7.530 | 75 | 7.582 | 76 | 28-120 | 1 | 0-20 | |
| Fluorene | ND | 10.00 | 7.081 | 71 | 7.049 | 70 | 12-186 | 0 | 0-20 | |
| N-Nitroso-di-n-propylamine | ND | 10.00 | 4.716 | 47 | 4.522 | 45 | 38-140 | 4 | 0-20 | |
| Naphthalene | ND | 10.00 | 5.303 | 53 | 4.971 | 50 | 20-140 | 6 | 0-20 | |
| 4-Nitrophenol | ND | 10.00 | 5.622 | 56 | 5.872 | 59 | 14-128 | 4 | 0-59 | |
| Pentachlorophenol | ND | 10.00 | 6.067 | 61 | 6.228 | 62 | 10-124 | 3 | 0-20 | |
| Phenol | ND | 10.00 | 5.633 | 56 | 5.517 | 55 | 22-124 | 2 | 0-20 | |
| Pyrene | ND | 10.00 | 6.422 | 64 | 6.249 | 62 | 31-169 | 3 | 0-20 | |
| 1,2,4-Trichlorobenzene | ND | 10.00 | 5.249 | 52 | 4.815 | 48 | 56-120 | 9 | 0-20 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-06-0712-1 | Sample | Aqueous | GC/MS V V | 06/13/15 | 06/13/15 17:14 | 150613S007 |
| 15-06-0712-1 | Matrix Spike | Aqueous | GC/MS V V | 06/13/15 | 06/13/15 10:14 | 150613S007 |
| 15-06-0712-1 | Matrix Spike Duplicate | Aqueous | GC/MS V V | 06/13/15 | 06/13/15 10:42 | 150613S007 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Acetone | ND | 50.00 | 46.15 | 92 | 47.23 | 94 | 51-171 | 2 | 0-20 | |
| Benzene | ND | 50.00 | 60.32 | 121 | 54.65 | 109 | 80-120 | 10 | 0-20 | 3 |
| Bromobenzene | ND | 50.00 | 65.62 | 131 | 60.39 | 121 | 54-150 | 8 | 0-20 | |
| Bromochloromethane | ND | 50.00 | 50.17 | 100 | 45.86 | 92 | 77-125 | 9 | 0-20 | |
| Bromodichloromethane | ND | 50.00 | 65.43 | 131 | 60.26 | 121 | 78-126 | 8 | 0-20 | 3 |
| Bromoform | ND | 50.00 | 59.63 | 119 | 57.20 | 114 | 41-155 | 4 | 0-20 | |
| Bromomethane | ND | 50.00 | 70.80 | 142 | 64.07 | 128 | 30-152 | 10 | 0-20 | |
| 2-Butanone | ND | 50.00 | 47.59 | 95 | 49.76 | 100 | 52-160 | 4 | 0-20 | |
| n-Butylbenzene | ND | 50.00 | 59.42 | 119 | 53.53 | 107 | 50-164 | 10 | 0-20 | |
| sec-Butylbenzene | ND | 50.00 | 58.33 | 117 | 52.26 | 105 | 49-157 | 11 | 0-20 | |
| tert-Butylbenzene | ND | 50.00 | 61.28 | 123 | 54.78 | 110 | 48-156 | 11 | 0-20 | |
| Carbon Disulfide | ND | 50.00 | 70.54 | 141 | 62.65 | 125 | 69-123 | 12 | 0-20 | 3 |
| Carbon Tetrachloride | ND | 50.00 | 60.98 | 122 | 53.99 | 108 | 62-140 | 12 | 0-20 | |
| Chlorobenzene | ND | 50.00 | 62.35 | 125 | 56.67 | 113 | 52-148 | 10 | 0-20 | |
| Chloroethane | ND | 50.00 | 52.19 | 104 | 46.08 | 92 | 66-132 | 12 | 0-20 | |
| Chloroform | ND | 50.00 | 62.85 | 126 | 57.63 | 115 | 80-122 | 9 | 0-20 | 3 |
| Chloromethane | ND | 50.00 | 65.49 | 131 | 56.98 | 114 | 45-147 | 14 | 0-20 | |
| 2-Chlorotoluene | ND | 50.00 | 61.84 | 124 | 56.68 | 113 | 51-153 | 9 | 0-20 | |
| 4-Chlorotoluene | ND | 50.00 | 58.24 | 116 | 53.32 | 107 | 49-151 | 9 | 0-20 | |
| Dibromochloromethane | ND | 50.00 | 66.08 | 132 | 62.12 | 124 | 48-150 | 6 | 0-20 | |
| 1,2-Dibromo-3-Chloropropane | ND | 50.00 | 52.16 | 104 | 50.64 | 101 | 46-142 | 3 | 0-20 | |
| 1,2-Dibromoethane | ND | 50.00 | 63.97 | 128 | 60.94 | 122 | 51-147 | 5 | 0-20 | |
| Dibromomethane | ND | 50.00 | 60.93 | 122 | 57.97 | 116 | 80-123 | 5 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 58.79 | 118 | 55.38 | 111 | 51-147 | 6 | 0-20 | |
| 1,3-Dichlorobenzene | ND | 50.00 | 59.71 | 119 | 54.64 | 109 | 49-151 | 9 | 0-20 | |
| 1,4-Dichlorobenzene | ND | 50.00 | 60.36 | 121 | 55.49 | 111 | 51-147 | 8 | 0-20 | |
| Dichlorodifluoromethane | ND | 50.00 | 58.89 | 118 | 51.88 | 104 | 30-170 | 13 | 0-20 | |
| 1,1-Dichloroethane | ND | 50.00 | 58.06 | 116 | 53.13 | 106 | 67-127 | 9 | 0-20 | |
| 1,2-Dichloroethane | ND | 50.00 | 61.85 | 124 | 58.79 | 118 | 73-133 | 5 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 68.97 | 138 | 61.73 | 123 | 68-128 | 11 | 0-20 | 3 |
| c-1,2-Dichloroethene | ND | 50.00 | 64.22 | 128 | 58.47 | 117 | 77-125 | 9 | 0-20 | 3 |
| t-1,2-Dichloroethene | ND | 50.00 | 65.81 | 132 | 59.15 | 118 | 71-131 | 11 | 0-20 | 3 |
| 1,2-Dichloropropane | ND | 50.00 | 61.70 | 123 | 56.77 | 114 | 80-120 | 8 | 0-20 | 3 |
| 1,3-Dichloropropane | ND | 50.00 | 64.05 | 128 | 60.98 | 122 | 50-146 | 5 | 0-20 | |
| 2,2-Dichloropropane | ND | 50.00 | 69.48 | 139 | 61.93 | 124 | 30-170 | 11 | 0-20 | |

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|---------------------------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| 1,1-Dichloropropene | ND | 50.00 | 58.82 | 118 | 52.21 | 104 | 75-129 | 12 | 0-20 | |
| c-1,3-Dichloropropene | ND | 50.00 | 66.55 | 133 | 61.44 | 123 | 80-124 | 8 | 0-20 | 3 |
| t-1,3-Dichloropropene | ND | 50.00 | 64.61 | 129 | 60.32 | 121 | 47-143 | 7 | 0-20 | |
| Ethylbenzene | ND | 50.00 | 65.16 | 130 | 58.67 | 117 | 54-150 | 10 | 0-20 | |
| 2-Hexanone | ND | 50.00 | 53.27 | 107 | 53.33 | 107 | 44-152 | 0 | 0-20 | |
| Isopropylbenzene | ND | 50.00 | 62.91 | 126 | 56.56 | 113 | 52-154 | 11 | 0-20 | |
| p-Isopropyltoluene | ND | 50.00 | 61.89 | 124 | 55.68 | 111 | 49-151 | 11 | 0-20 | |
| Methylene Chloride | ND | 50.00 | 66.23 | 132 | 61.69 | 123 | 73-127 | 7 | 0-20 | 3 |
| 4-Methyl-2-Pentanone | ND | 50.00 | 53.25 | 107 | 55.82 | 112 | 70-124 | 5 | 0-20 | |
| Naphthalene | ND | 50.00 | 50.20 | 100 | 49.55 | 99 | 39-153 | 1 | 0-20 | |
| n-Propylbenzene | ND | 50.00 | 60.95 | 122 | 54.84 | 110 | 49-157 | 11 | 0-20 | |
| Styrene | ND | 50.00 | 64.25 | 128 | 59.29 | 119 | 54-150 | 8 | 0-20 | |
| 1,1,1,2-Tetrachloroethane | ND | 50.00 | 70.23 | 140 | 65.63 | 131 | 50-152 | 7 | 0-20 | |
| 1,1,2,2-Tetrachloroethane | ND | 50.00 | 59.40 | 119 | 57.57 | 115 | 44-146 | 3 | 0-20 | |
| Tetrachloroethene | ND | 50.00 | 60.72 | 121 | 52.87 | 106 | 34-170 | 14 | 0-20 | |
| Toluene | ND | 50.00 | 65.50 | 131 | 58.89 | 118 | 80-120 | 11 | 0-20 | 3 |
| 1,2,3-Trichlorobenzene | ND | 50.00 | 57.76 | 116 | 54.81 | 110 | 41-161 | 5 | 0-20 | |
| 1,2,4-Trichlorobenzene | ND | 50.00 | 58.98 | 118 | 55.36 | 111 | 41-161 | 6 | 0-20 | |
| 1,1,1-Trichloroethane | ND | 50.00 | 67.27 | 135 | 59.13 | 118 | 75-129 | 13 | 0-20 | 3 |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 50.00 | 75.94 | 152 | 66.46 | 133 | 54-156 | 13 | 0-20 | |
| 1,1,2-Trichloroethane | ND | 50.00 | 60.53 | 121 | 58.06 | 116 | 51-147 | 4 | 0-20 | |
| Trichloroethene | ND | 50.00 | 66.25 | 133 | 59.17 | 118 | 80-120 | 11 | 0-20 | 3 |
| Trichlorofluoromethane | ND | 50.00 | 60.12 | 120 | 52.62 | 105 | 61-145 | 13 | 0-20 | |
| 1,2,3-Trichloropropane | ND | 50.00 | 56.93 | 114 | 56.85 | 114 | 51-147 | 0 | 0-20 | |
| 1,2,4-Trimethylbenzene | ND | 50.00 | 60.42 | 121 | 55.19 | 110 | 56-152 | 9 | 0-20 | |
| 1,3,5-Trimethylbenzene | ND | 50.00 | 66.37 | 133 | 60.95 | 122 | 56-158 | 9 | 0-20 | |
| Vinyl Acetate | ND | 50.00 | 71.87 | 144 | 69.70 | 139 | 35-167 | 3 | 0-20 | |
| Vinyl Chloride | ND | 50.00 | 58.36 | 117 | 51.44 | 103 | 67-133 | 13 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 117.1 | 117 | 106.5 | 106 | 51-153 | 9 | 0-20 | |
| o-Xylene | ND | 50.00 | 58.46 | 117 | 53.47 | 107 | 51-153 | 9 | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 57.98 | 116 | 55.94 | 112 | 64-130 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-06-0547-10 | Sample | Solid | GC/MS W | 06/07/15 | 06/08/15 01:40 | 150607S007 | | | | |
| 15-06-0547-10 | Matrix Spike | Solid | GC/MS W | 06/07/15 | 06/08/15 02:09 | 150607S007 | | | | |
| 15-06-0547-10 | Matrix Spike Duplicate | Solid | GC/MS W | 06/07/15 | 06/08/15 02:38 | 150607S007 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 42.85 | 86 | 45.55 | 91 | 61-127 | 6 | 0-20 | |
| Carbon Tetrachloride | ND | 50.00 | 44.95 | 90 | 47.45 | 95 | 51-135 | 5 | 0-29 | |
| Chlorobenzene | ND | 50.00 | 39.37 | 79 | 41.75 | 84 | 57-123 | 6 | 0-20 | |
| 1,2-Dibromoethane | ND | 50.00 | 42.83 | 86 | 43.60 | 87 | 64-124 | 2 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 36.73 | 73 | 37.92 | 76 | 35-131 | 3 | 0-25 | |
| 1,2-Dichloroethane | ND | 50.00 | 48.74 | 97 | 49.50 | 99 | 80-120 | 2 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 48.65 | 97 | 50.57 | 101 | 47-143 | 4 | 0-25 | |
| Ethylbenzene | ND | 50.00 | 41.88 | 84 | 44.53 | 89 | 57-129 | 6 | 0-22 | |
| Toluene | ND | 50.00 | 42.10 | 84 | 44.90 | 90 | 63-123 | 6 | 0-20 | |
| Trichloroethene | ND | 50.00 | 47.97 | 96 | 50.35 | 101 | 44-158 | 5 | 0-20 | |
| Vinyl Chloride | ND | 50.00 | 48.13 | 96 | 53.17 | 106 | 49-139 | 10 | 0-47 | |
| p/m-Xylene | ND | 100.0 | 84.06 | 84 | 88.21 | 88 | 70-130 | 5 | 0-30 | |
| o-Xylene | ND | 50.00 | 42.15 | 84 | 43.91 | 88 | 70-130 | 4 | 0-30 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 42.96 | 86 | 43.24 | 86 | 57-123 | 1 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|--------------|-------------|------------|----------------|----------------|-----------------------|
| 15-06-0498-3 | Sample | Solid | ICP/MS 04 | 06/08/15 00:00 | 06/09/15 13:52 | 150608S01 |
| 15-06-0498-3 | PDS | Solid | ICP/MS 04 | 06/08/15 00:00 | 06/09/15 13:45 | 150608S01 |
| Parameter | Sample Conc. | Spike Added | PDS Conc. | PDS %Rec. | %Rec. CL | Qualifiers |
| Antimony | ND | 25.00 | 26.69 | 107 | 75-125 | |
| Arsenic | 7.643 | 25.00 | 33.82 | 105 | 75-125 | |
| Barium | 199.5 | 25.00 | 227.8 | 4X | 75-125 | Q |
| Beryllium | ND | 25.00 | 27.15 | 109 | 75-125 | |
| Cadmium | 3.397 | 25.00 | 29.90 | 106 | 75-125 | |
| Chromium | 26.68 | 25.00 | 53.76 | 108 | 75-125 | |
| Copper | 29.47 | 25.00 | 55.21 | 103 | 75-125 | |
| Lead | 8.765 | 25.00 | 34.90 | 105 | 75-125 | |
| Nickel | 48.32 | 25.00 | 73.02 | 99 | 75-125 | |
| Selenium | ND | 25.00 | 27.82 | 111 | 75-125 | |
| Silver | ND | 12.50 | 12.69 | 102 | 75-125 | |
| Thallium | ND | 25.00 | 26.06 | 104 | 75-125 | |
| Zinc | 92.30 | 25.00 | 123.1 | 123 | 75-125 | |
| Aluminum | 11780 | 25.00 | 11860 | 4X | 75-125 | Q |
| Iron | 25280 | 25.00 | 25040 | 4X | 75-125 | Q |
| Manganese | 332.0 | 25.00 | 353.2 | 4X | 75-125 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: ASTM D-2216 (M)

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|--------|------------|----------------|----------------|------------------------|
| 15-06-0567-1 | Sample | Solid | N/A | 06/08/15 00:00 | 06/08/15 20:00 | F0608MOID6 |
| 15-06-0567-1 | Sample Duplicate | Solid | N/A | 06/08/15 00:00 | 06/08/15 20:00 | F0608MOID6 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|------------------|---------------------|------------------|------------|---------------|-------------------|
| Moisture | 71.30 | 72.00 | 1 | 0-10 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: SM 2320B M

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|--------|------------|---------------|----------------|------------------------|
| 15-06-0770-3 | Sample | Solid | PH1/BUR03 | N/A | 06/15/15 16:35 | F0615ALKD2 |
| 15-06-0770-3 | Sample Duplicate | Solid | PH1/BUR03 | N/A | 06/15/15 16:35 | F0615ALKD2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|---|---------------------|------------------|------------|---------------|-------------------|
| Alkalinity, Total (as CaCO ₃) | 200.0 | 195.0 | 3 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: SM 2320B M

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|--------|------------|----------------|----------------|------------------------|
| 15-06-0770-3 | Sample | Solid | PH1/BUR03 | 06/15/15 00:00 | 06/15/15 16:35 | F0615HCOD2 |
| 15-06-0770-3 | Sample Duplicate | Solid | PH1/BUR03 | 06/15/15 00:00 | 06/15/15 16:35 | F0615HCOD2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Bicarbonate (as CaCO ₃) | 200.0 | 195.0 | 3 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: SM 2320B M

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|--------|------------|----------------|----------------|------------------------|
| 15-06-0770-3 | Sample | Solid | PH1/BUR03 | 06/15/15 00:00 | 06/15/15 16:35 | F0615CO3D2 |
| 15-06-0770-3 | Sample Duplicate | Solid | PH1/BUR03 | 06/15/15 00:00 | 06/15/15 16:35 | F0615CO3D2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: SM 2540 C (M)

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|---------------------|------------------|-----------------------|-----------------------|------------------------|
| HCS330 | Sample | Solid | SC 2 | 06/12/15 00:00 | 06/12/15 19:00 | F0612TDSD4 |
| HCS330 | Sample Duplicate | Solid | SC 2 | 06/12/15 00:00 | 06/12/15 19:00 | F0612TDSD4 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Solids, Total Dissolved | | 17870 | 17900 | 0 | 0-10 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 099-12-922-584 | LCS | Solid | IC 10 | 06/10/15 | 06/11/15 00:37 | 150610L01P |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Fluoride | | 25.00 | 23.96 | 96 | 90-110 | |
| Chloride | | 500.0 | 537.1 | 107 | 90-110 | |
| Bromide | | 50.00 | 51.66 | 103 | 90-110 | |
| Nitrate (as N) | | 50.00 | 53.74 | 107 | 90-110 | |
| Sulfate | | 500.0 | 541.4 | 108 | 90-110 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: EPA 9060A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-06-013-1274 | LCS | Solid | TOC 5 | 06/17/15 | 06/17/15 15:52 | F0617TOCL1 | | | |
| 099-06-013-1274 | LCSD | Solid | TOC 5 | 06/17/15 | 06/17/15 15:52 | F0617TOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 6000 | 6524 | 109 | 6516 | 109 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: N/A
Method: SM 4500 P B/E (M)

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|--------|------------|---------------|----------------|-----------------------|
| 099-05-001-5419 | LCS | Solid | UV 7 | 06/10/15 | 06/11/15 16:26 | F0611TPL2 |
| 099-05-001-5419 | LCSD | Solid | UV 7 | 06/10/15 | 06/11/15 16:26 | F0611TPL2 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| Phosphorus, Total | 2.000 | 1.940 | 97 | 1.945 | 97 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|--------------------|------------------------|------------------|-----------------|-------------------|
| 097-01-002-21188 | LCS | Solid | ICP 7300 | 06/10/15 | 06/11/15 16:39 | 150610L04 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Calcium | | 25.00 | 26.08 | 104 | 80-120 | |
| Magnesium | | 25.00 | 25.28 | 101 | 80-120 | |
| Potassium | | 250.0 | 256.9 | 103 | 80-120 | |
| Sodium | | 250.0 | 250.5 | 100 | 80-120 | |
| Strontium | | 25.00 | 24.61 | 98 | 80-120 | |
| Silicon | | 25.00 | 23.68 | 95 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3050B
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-621-953 | LCS | Solid | ICP/MS 04 | 06/08/15 | 06/09/15 13:33 | 150608L01 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 25.00 | 25.40 | 102 | 80-120 | 73-127 | |
| Arsenic | 25.00 | 25.07 | 100 | 80-120 | 73-127 | |
| Barium | 25.00 | 25.73 | 103 | 80-120 | 73-127 | |
| Beryllium | 25.00 | 28.52 | 114 | 80-120 | 73-127 | |
| Cadmium | 25.00 | 26.46 | 106 | 80-120 | 73-127 | |
| Chromium | 25.00 | 27.65 | 111 | 80-120 | 73-127 | |
| Copper | 25.00 | 27.47 | 110 | 80-120 | 73-127 | |
| Lead | 25.00 | 25.94 | 104 | 80-120 | 73-127 | |
| Nickel | 25.00 | 26.52 | 106 | 80-120 | 73-127 | |
| Selenium | 25.00 | 26.62 | 106 | 80-120 | 73-127 | |
| Silver | 12.50 | 12.57 | 101 | 80-120 | 73-127 | |
| Thallium | 25.00 | 25.32 | 101 | 80-120 | 73-127 | |
| Zinc | 25.00 | 27.45 | 110 | 80-120 | 73-127 | |
| Aluminum | 25.00 | 25.57 | 102 | 80-120 | 73-127 | |
| Iron | 25.00 | 30.84 | 123 | 80-120 | 73-127 | ME |
| Manganese | 25.00 | 26.87 | 107 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 1

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 7471A Total
Method: EPA 7471A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------|-------------------|-----------------|-----------------------|------------------|
| 099-16-272-1352 | LCS | Solid | Mercury 05 | 06/16/15 | 06/16/15 14:30 | 150616L01 |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.8350 | 0.9767 | 117 | 85-121 | |


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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-12-537-2140 | LCS | Solid | GC 44 | 06/13/15 | 06/16/15 20:03 | 150613L05 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Aldrin | 25.00 | 22.10 | 88 | 50-135 | 36-149 | |
| Alpha-BHC | 25.00 | 21.00 | 84 | 50-135 | 36-149 | |
| Beta-BHC | 25.00 | 22.14 | 89 | 50-135 | 36-149 | |
| 4,4'-DDD | 25.00 | 25.08 | 100 | 50-135 | 36-149 | |
| 4,4'-DDE | 25.00 | 25.18 | 101 | 50-135 | 36-149 | |
| 4,4'-DDT | 25.00 | 24.78 | 99 | 50-135 | 36-149 | |
| Delta-BHC | 25.00 | 22.71 | 91 | 50-135 | 36-149 | |
| Dieldrin | 25.00 | 24.27 | 97 | 50-135 | 36-149 | |
| Endosulfan I | 25.00 | 22.44 | 90 | 50-135 | 36-149 | |
| Endosulfan II | 25.00 | 23.55 | 94 | 50-135 | 36-149 | |
| Endosulfan Sulfate | 25.00 | 24.34 | 97 | 50-135 | 36-149 | |
| Endrin | 25.00 | 27.85 | 111 | 50-135 | 36-149 | |
| Endrin Aldehyde | 25.00 | 21.15 | 85 | 50-135 | 36-149 | |
| Gamma-BHC | 25.00 | 22.42 | 90 | 50-135 | 36-149 | |
| Heptachlor | 25.00 | 24.40 | 98 | 50-135 | 36-149 | |
| Heptachlor Epoxide | 25.00 | 23.13 | 93 | 50-135 | 36-149 | |
| Methoxychlor | 25.00 | 27.88 | 112 | 50-135 | 36-149 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 099-12-535-3269 | LCS | Solid | GC 31 | 06/13/15 | 06/16/15 11:02 | 150613L06 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Aroclor-1016 | | 100.0 | 106.4 | 106 | 50-135 | |
| Aroclor-1260 | | 100.0 | 93.42 | 93 | 50-135 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-973-187 | LCS | Solid | GC 26 | 06/08/15 | 06/09/15 12:25 | 150608L12 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Azinphos Methyl | 4.000 | 4.834 | 121 | 30-130 | 13-147 | |
| Bolstar | 4.000 | 4.652 | 116 | 30-130 | 13-147 | |
| Chlorpyrifos | 4.000 | 4.669 | 117 | 30-130 | 13-147 | |
| Coumaphos | 4.000 | 5.010 | 125 | 30-130 | 13-147 | |
| Diazinon | 4.000 | 4.173 | 104 | 30-130 | 13-147 | |
| Disulfoton | 4.000 | 5.113 | 128 | 30-130 | 13-147 | |
| Ethoprop | 4.000 | 4.134 | 103 | 30-130 | 13-147 | |
| Fensulfothion | 4.000 | 4.946 | 124 | 30-130 | 13-147 | |
| Fenthion | 4.000 | 4.893 | 122 | 30-130 | 13-147 | |
| Merphos | 4.000 | 4.870 | 122 | 30-130 | 13-147 | |
| Methyl Parathion | 4.000 | 4.779 | 119 | 30-130 | 13-147 | |
| Phorate | 4.000 | 4.829 | 121 | 30-130 | 13-147 | |
| Ronnel | 4.000 | 5.118 | 128 | 30-130 | 13-147 | |
| Stirophos | 4.000 | 4.586 | 115 | 30-130 | 13-147 | |
| Tokuthion | 4.000 | 5.025 | 126 | 30-130 | 13-147 | |
| Trichloronate | 4.000 | 5.092 | 127 | 30-130 | 13-147 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 095-01-033-1294 | LCS | Solid | GC 40 | 06/10/15 | 06/11/15 18:57 | 150610L09 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| 2,4-D | | 400.0 | 284.1 | 71 | 30-130 | |
| 2,4,5-T | | 40.00 | 28.70 | 72 | 30-130 | |
| 2,4-DB | | 400.0 | 291.7 | 73 | 30-130 | |


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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 3545
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-12-549-3310 | LCS | Solid | GC/MS SS | 06/13/15 | 06/15/15 21:35 | 150613L07 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Acenaphthene | 10.00 | 8.591 | 86 | 51-123 | 39-135 | |
| Acenaphthylene | 10.00 | 8.478 | 85 | 52-120 | 41-131 | |
| Butyl Benzyl Phthalate | 10.00 | 8.423 | 84 | 43-139 | 27-155 | |
| 4-Chloro-3-Methylphenol | 10.00 | 7.741 | 77 | 55-121 | 44-132 | |
| 2-Chlorophenol | 10.00 | 7.668 | 77 | 58-124 | 47-135 | |
| 1,4-Dichlorobenzene | 10.00 | 6.881 | 69 | 42-132 | 27-147 | |
| Dimethyl Phthalate | 10.00 | 8.440 | 84 | 51-123 | 39-135 | |
| 2,4-Dinitrotoluene | 10.00 | 9.918 | 99 | 51-129 | 38-142 | |
| Fluorene | 10.00 | 9.052 | 91 | 54-126 | 42-138 | |
| N-Nitroso-di-n-propylamine | 10.00 | 6.254 | 63 | 40-136 | 24-152 | |
| Naphthalene | 10.00 | 7.284 | 73 | 32-146 | 13-165 | |
| 4-Nitrophenol | 10.00 | 7.807 | 78 | 24-126 | 7-143 | |
| Pentachlorophenol | 10.00 | 8.981 | 90 | 23-131 | 5-149 | |
| Phenol | 10.00 | 7.050 | 71 | 40-130 | 25-145 | |
| Pyrene | 10.00 | 7.960 | 80 | 47-143 | 31-159 | |
| 1,2,4-Trichlorobenzene | 10.00 | 7.701 | 77 | 45-129 | 31-143 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-17397 | LCS | Aqueous | GC/MS V V | 06/13/15 | 06/13/15 09:32 | 150613L025 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 50.56 | 101 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 49.57 | 99 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 52.02 | 104 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 56.81 | 114 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 52.02 | 104 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 54.92 | 110 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 56.26 | 113 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 53.93 | 108 | 80-123 | 73-130 | |
| Toluene | 50.00 | 54.97 | 110 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 54.93 | 110 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 49.17 | 98 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 97.56 | 98 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 48.99 | 98 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 51.57 | 103 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/05/15
Work Order: 15-06-0463
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-12-796-9798 | LCS | Solid | GC/MS W | 06/07/15 | 06/07/15 23:44 | 150607L024 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 51.97 | 104 | 78-120 | 71-127 | |
| Carbon Tetrachloride | 50.00 | 54.37 | 109 | 49-139 | 34-154 | |
| Chlorobenzene | 50.00 | 49.55 | 99 | 79-120 | 72-127 | |
| 1,2-Dibromoethane | 50.00 | 52.72 | 105 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 47.85 | 96 | 75-120 | 68-128 | |
| 1,2-Dichloroethane | 50.00 | 61.22 | 122 | 80-120 | 73-127 | ME |
| 1,1-Dichloroethene | 50.00 | 54.63 | 109 | 74-122 | 66-130 | |
| Ethylbenzene | 50.00 | 52.06 | 104 | 76-120 | 69-127 | |
| Toluene | 50.00 | 51.93 | 104 | 77-120 | 70-127 | |
| Trichloroethene | 50.00 | 55.04 | 110 | 80-120 | 73-127 | |
| Vinyl Chloride | 50.00 | 55.08 | 110 | 68-122 | 59-131 | |
| p/m-Xylene | 100.0 | 104.9 | 105 | 75-125 | 67-133 | |
| o-Xylene | 50.00 | 51.87 | 104 | 75-125 | 67-133 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 51.75 | 103 | 77-120 | 70-127 | |

Total number of LCS compounds: 14

Total number of ME compounds: 1

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits

Glossary of Terms and Qualifiers

Work Order: 15-06-0463

Page 1 of 1

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|---|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| CI | See case narrative. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |
| | Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time. |
| | A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations. |

9463

ORIGIN ID:SATA (210) 877-2847
SWCA INC

6200 UTSA BLVD STE 102

SAN ANTONIO, TX 782491618
UNITED STATES US

SHIP DATE: 04JUN15
ACTWGT: 33.4 LB
CAD: /OFFC1601
DIMS: 20x15x12 IN

BILL SENDER

Part # 156297-435 RIT2 12/14
REF: JUNE 11/15 8:00 PM ClearCase

TO

CALSCIENCE
7440 LINCOLN WAY

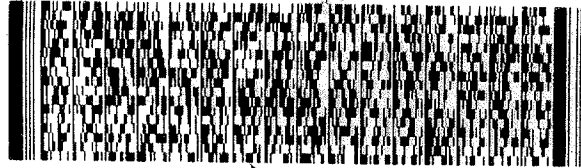
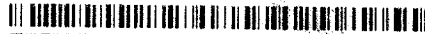
GARDEN GROVE CA 92841

(714) 895-5494

INU:
PO:

REF:

DEPT:



FedEx
Express

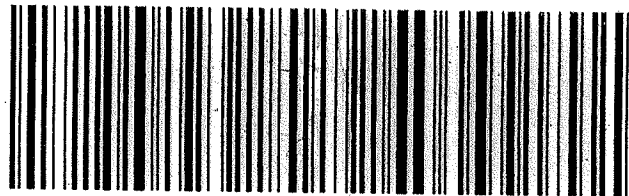


TRK# 8062 7398 5439
0215

FRI - 05 JUN AA
STANDARD OVERNIGHT

A7 APVA

DSR
92841
CA-US SNA



SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 1CLIENT: SWCADATE: 06 / 05 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 3-1 °C (w/ CF): 2.8 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ FilterChecked by: IS

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Present but Not Intact☐ Not Present☐ N/AChecked by: ISSample(s) ☐ Present and Intact☐ Present but Not Intact☒ Not Present☐ N/AChecked by: qbs

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: 150508B)Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na2} ☐ 100PJ ☐ 100PJ_{na2} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☐ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s☐ 500PB ☐ 1AGB ☐ 1AGB_{na2} ☐ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☒ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☒ 202CGJAir: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: qbss = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOHReviewed by: qbs

**WORK ORDER NUMBER: 15-06-0567***The difference is service*

AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For**Client:** SWCA Environmental Consultants**Client Project Name:** EAA 27122**Attention:** Philip Pearce
6200 UTSA Blvd.
Suite 102
San Antonio, TX 78249-1618

A handwritten signature in black ink, appearing to read "H. Burley", with the letters "FOR" written in a smaller font below it.

Approved for release on 06/25/2015 by:
Don Burley
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

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Client Project Name: EAA 27122
Work Order Number: 15-06-0567

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Work Order Narrative

Work Order: 15-06-0567

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Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 06/06/15. They were assigned to Work Order 15-06-0567.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

Solid Samples 15-06-0567-7, -8 (HSM370 and FDHSM370) for TOC analysis contained a combination of soil, sand, gravel, and water.



Calscience

Sample Summary

| | |
|--|------------------------------------|
| Client: SWCA Environmental Consultants | Work Order: 15-06-0567 |
| 6200 UTSA Blvd., Suite 102 | Project Name: EAA 27122 |
| San Antonio, TX 78249-1618 | PO Number: |
| | Date/Time Received: 06/06/15 09:30 |
| | Number of Containers: 26 |
| Attn: Philip Pearce | |

| Sample Identification | Lab Number | Collection Date and Time | Number of Containers | Matrix |
|-----------------------|---------------|--------------------------|----------------------|---------|
| HSM310 | 15-06-0567-1 | 06/05/15 10:03 | 2 | Solid |
| HSM320 | 15-06-0567-2 | 06/05/15 10:48 | 2 | Solid |
| HSM330 | 15-06-0567-3 | 06/05/15 11:03 | 2 | Solid |
| HSM340 | 15-06-0567-4 | 06/05/15 11:21 | 2 | Solid |
| HSM350 | 15-06-0567-5 | 06/05/15 12:24 | 2 | Solid |
| HSM360 | 15-06-0567-6 | 06/05/15 13:13 | 2 | Solid |
| HSM370 | 15-06-0567-7 | 06/05/15 13:51 | 2 | Solid |
| FDHSM370 | 15-06-0567-8 | 06/05/15 13:51 | 2 | Solid |
| TB08 | 15-06-0567-9 | 06/05/15 00:00 | 1 | Aqueous |
| EB01 | 15-06-0567-10 | 06/05/15 15:41 | 9 | Aqueous |

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Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB01 | 15-06-0567-10-I | 06/05/15 15:41 | Aqueous | IC 7 | N/A | 06/06/15 13:27 | 150606L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | ND | 0.10 | 0.025 | 1.00 | |
| Chloride | ND | 1.0 | 0.12 | 1.00 | |
| Bromide | ND | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | ND | 0.10 | 0.025 | 1.00 | |
| Sulfate | ND | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|--------------|-----------------|-----|---------|------|-----|----------------|-----------|
| Method Blank | 099-12-906-5827 | N/A | Aqueous | IC 7 | N/A | 06/06/15 10:26 | 150606L01 |
|--------------|-----------------|-----|---------|------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | ND | 0.10 | 0.025 | 1.00 | |
| Chloride | ND | 1.0 | 0.12 | 1.00 | |
| Bromide | ND | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | ND | 0.10 | 0.025 | 1.00 | |
| Sulfate | ND | 1.0 | 0.19 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 300.0
Units: mg/kg

Project: EAA 27122

Page 1 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM310 | 15-06-0567-1-A | 06/05/15 10:03 | Solid | IC 10 | 06/09/15 | 06/09/15 12:15 | 150609L01P |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|-----|------|------|------------|
| Fluoride | 5.5 | 3.5 | 1.1 | 1.00 | |
| Chloride | 55 | 35 | 5.1 | 1.00 | |
| Bromide | ND | 3.5 | 1.2 | 1.00 | |
| Nitrate (as N) | 1.2 | 3.5 | 0.85 | 1.00 | J |
| Sulfate | 420 | 35 | 7.4 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|
| HSM320 | 15-06-0567-2-A | 06/05/15 10:48 | Solid | IC 10 | 06/09/15 | 06/09/15 12:31 | 150609L01P |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|-----|-----|------|------------|
| Fluoride | 6.2 | 5.1 | 1.6 | 1.00 | |
| Chloride | 84 | 51 | 7.4 | 1.00 | |
| Bromide | ND | 5.1 | 1.8 | 1.00 | |
| Nitrate (as N) | 2.0 | 5.1 | 1.2 | 1.00 | J |
| Sulfate | 220 | 51 | 11 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|
| HSM330 | 15-06-0567-3-A | 06/05/15 11:03 | Solid | IC 10 | 06/09/15 | 06/09/15 12:48 | 150609L01P |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|-----|------|------|------------|
| Fluoride | 0.92 | 1.3 | 0.39 | 1.00 | J |
| Chloride | 7.7 | 13 | 1.8 | 1.00 | J |
| Bromide | ND | 1.3 | 0.45 | 1.00 | |
| Nitrate (as N) | 0.73 | 1.3 | 0.31 | 1.00 | J |
| Sulfate | 11 | 13 | 2.7 | 1.00 | J |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 300.0
Units: mg/kg

Project: EAA 27122

Page 2 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM340 | 15-06-0567-4-A | 06/05/15 11:21 | Solid | IC 10 | 06/09/15 | 06/09/15 13:04 | 150609L01P |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|-----|------|------|------------|
| Fluoride | 2.0 | 1.2 | 0.39 | 1.00 | |
| Chloride | 7.9 | 12 | 1.8 | 1.00 | J |
| Bromide | ND | 1.2 | 0.44 | 1.00 | |
| Nitrate (as N) | ND | 1.2 | 0.30 | 1.00 | |
| Sulfate | 26 | 12 | 2.6 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|
| HSM350 | 15-06-0567-5-A | 06/05/15 12:24 | Solid | IC 10 | 06/09/15 | 06/09/15 13:21 | 150609L01P |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|-----|------|------|------------|
| Fluoride | 2.8 | 2.2 | 0.68 | 1.00 | |
| Chloride | 27 | 22 | 3.2 | 1.00 | |
| Bromide | ND | 2.2 | 0.78 | 1.00 | |
| Nitrate (as N) | ND | 2.2 | 0.53 | 1.00 | |
| Sulfate | 130 | 22 | 4.6 | 1.00 | |

| | | | | | | | |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|
| HSM360 | 15-06-0567-6-A | 06/05/15 13:13 | Solid | IC 10 | 06/09/15 | 06/09/15 13:37 | 150609L01P |
|--------|----------------|----------------|-------|-------|----------|----------------|------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|-----|------|------|------------|
| Fluoride | 1.3 | 1.3 | 0.42 | 1.00 | J |
| Chloride | 8.9 | 13 | 2.0 | 1.00 | J |
| Bromide | ND | 1.3 | 0.48 | 1.00 | |
| Nitrate (as N) | 0.56 | 1.3 | 0.33 | 1.00 | J |
| Sulfate | 13 | 13 | 2.8 | 1.00 | J |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 300.0
Units: mg/kg

Project: EAA 27122

Page 3 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM370 | 15-06-0567-7-A | 06/05/15 13:51 | Solid | IC 10 | 06/09/15 | 06/09/15 13:53 | 150609L01P |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|-----|------|------|------------|
| Fluoride | 1.7 | 1.7 | 0.53 | 1.00 | |
| Chloride | 17 | 17 | 2.5 | 1.00 | J |
| Bromide | ND | 1.7 | 0.61 | 1.00 | |
| Nitrate (as N) | 0.69 | 1.7 | 0.42 | 1.00 | J |
| Sulfate | 30 | 17 | 3.6 | 1.00 | |

| | | | | | | | |
|----------|----------------|----------------|-------|-------|----------|----------------|------------|
| FDHSM370 | 15-06-0567-8-A | 06/05/15 13:51 | Solid | IC 10 | 06/09/15 | 06/09/15 14:10 | 150609L01P |
|----------|----------------|----------------|-------|-------|----------|----------------|------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|-----|------|------|------------|
| Fluoride | 2.0 | 1.8 | 0.57 | 1.00 | |
| Chloride | 16 | 18 | 2.7 | 1.00 | J |
| Bromide | ND | 1.8 | 0.66 | 1.00 | |
| Nitrate (as N) | 0.61 | 1.8 | 0.45 | 1.00 | J |
| Sulfate | 38 | 18 | 3.9 | 1.00 | |

| | | | | | | | |
|--------------|----------------|-----|-------|-------|----------|----------------|------------|
| Method Blank | 099-12-922-582 | N/A | Solid | IC 10 | 06/09/15 | 06/09/15 11:39 | 150609L01P |
|--------------|----------------|-----|-------|-------|----------|----------------|------------|

Comment(s):
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|-----|------|------|------------|
| Fluoride | ND | 1.0 | 0.31 | 1.00 | |
| Chloride | ND | 10 | 1.5 | 1.00 | |
| Bromide | ND | 1.0 | 0.36 | 1.00 | |
| Nitrate (as N) | ND | 1.0 | 0.24 | 1.00 | |
| Sulfate | ND | 10 | 2.1 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6010B
Units: mg/kg

Project: EAA 27122

Page 1 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM310 | 15-06-0567-1-BB | 06/05/15 10:03 | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:13 | 150611L09A |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|-------|-------|------------|
| Calcium | 150000 | 17.3 | 1.32 | 0.995 | |
| Magnesium | 3850 | 17.3 | 0.587 | 0.995 | |
| Potassium | 3010 | 86.7 | 6.08 | 0.995 | |
| Sodium | 334 | 86.7 | 6.29 | 0.995 | |
| Strontium | 295 | 5.20 | 0.477 | 0.995 | |
| Silicon | 3190 | 17.3 | 4.57 | 0.995 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|----------|----------|----------------|------------|
| HSM320 | 15-06-0567-2-BB | 06/05/15 10:48 | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:20 | 150611L09A |
|--------|-----------------|----------------|-------|----------|----------|----------------|------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|-------|------|------------|
| Calcium | 200000 | 26.9 | 2.04 | 1.05 | |
| Magnesium | 6770 | 26.9 | 0.909 | 1.05 | |
| Potassium | 4540 | 134 | 9.42 | 1.05 | |
| Sodium | 670 | 134 | 9.75 | 1.05 | |
| Strontium | 342 | 8.06 | 0.739 | 1.05 | |
| Silicon | 5380 | 26.9 | 7.07 | 1.05 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|----------|----------|----------------|------------|
| HSM330 | 15-06-0567-3-BB | 06/05/15 11:03 | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:22 | 150611L09A |
|--------|-----------------|----------------|-------|----------|----------|----------------|------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|------|------|------------|
| Calcium | 274000 | 65.2 | 4.96 | 10.4 | |
| Magnesium | 3500 | 65.2 | 2.21 | 10.4 | |
| Potassium | 741 | 326 | 22.9 | 10.4 | |
| Sodium | 422 | 326 | 23.7 | 10.4 | |
| Strontium | 149 | 19.6 | 1.79 | 10.4 | |
| Silicon | 965 | 65.2 | 17.2 | 10.4 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6010B
Units: mg/kg

Project: EAA 27122

Page 2 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM340 | 15-06-0567-4-BB | 06/05/15 11:21 | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:25 | 150611L09A |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|------|------|------------|
| Calcium | 224000 | 64.9 | 4.94 | 10.4 | |
| Magnesium | 3450 | 64.9 | 2.20 | 10.4 | |
| Potassium | 1120 | 325 | 22.8 | 10.4 | |
| Sodium | 344 | 325 | 23.6 | 10.4 | |
| Strontium | 151 | 19.5 | 1.79 | 10.4 | |
| Silicon | 1510 | 64.9 | 17.1 | 10.4 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|----------|----------|----------------|------------|
| HSM350 | 15-06-0567-5-BB | 06/05/15 12:24 | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:27 | 150611L09A |
|--------|-----------------|----------------|-------|----------|----------|----------------|------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|------|------|------------|
| Calcium | 158000 | 111 | 8.46 | 10.2 | |
| Magnesium | 3470 | 111 | 3.76 | 10.2 | |
| Potassium | 1500 | 556 | 39.0 | 10.2 | |
| Sodium | 245 | 556 | 40.4 | 10.2 | J |
| Strontium | 211 | 33.3 | 3.06 | 10.2 | |
| Silicon | 2500 | 111 | 29.3 | 10.2 | |

| | | | | | | | |
|--------|-----------------|----------------|-------|----------|----------|----------------|------------|
| HSM360 | 15-06-0567-6-BB | 06/05/15 13:13 | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:30 | 150611L09A |
|--------|-----------------|----------------|-------|----------|----------|----------------|------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|------|------|------------|
| Calcium | 214000 | 70.3 | 5.35 | 10.5 | |
| Magnesium | 2880 | 70.3 | 2.38 | 10.5 | |
| Potassium | 1620 | 351 | 24.7 | 10.5 | |
| Sodium | 183 | 351 | 25.5 | 10.5 | J |
| Strontium | 198 | 21.1 | 1.93 | 10.5 | |
| Silicon | 1500 | 70.3 | 18.5 | 10.5 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6010B
Units: mg/kg

Project: EAA 27122

Page 3 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM370 | 15-06-0567-7-BB | 06/05/15 13:51 | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:52 | 150611L09A |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|------|------|------------|
| Calcium | 339000 | 86.8 | 6.60 | 10.2 | |
| Magnesium | 4920 | 86.8 | 2.94 | 10.2 | |
| Potassium | 1860 | 434 | 30.5 | 10.2 | |
| Sodium | ND | 434 | 31.5 | 10.2 | |
| Strontium | 293 | 26.0 | 2.39 | 10.2 | |
| Silicon | 1860 | 86.8 | 22.9 | 10.2 | |

| | | | | | | | |
|----------|-----------------|----------------|-------|----------|----------|----------------|------------|
| FDHSM370 | 15-06-0567-8-BB | 06/05/15 13:51 | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:55 | 150611L09A |
|----------|-----------------|----------------|-------|----------|----------|----------------|------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|------|------|------------|
| Calcium | 312000 | 89.8 | 6.84 | 9.76 | |
| Magnesium | 5780 | 89.8 | 3.04 | 9.76 | |
| Potassium | 2230 | 449 | 31.5 | 9.76 | |
| Sodium | ND | 449 | 32.6 | 9.76 | |
| Strontium | 286 | 27.0 | 2.47 | 9.76 | |
| Silicon | 2050 | 89.8 | 23.7 | 9.76 | |

| | | | | | | | |
|--------------|------------------|-----|-------|----------|----------|----------------|------------|
| Method Blank | 097-01-002-21301 | N/A | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:08 | 150611L09A |
|--------------|------------------|-----|-------|----------|----------|----------------|------------|

Comment(s):
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|-------|-------|------------|
| Calcium | ND | 4.98 | 0.379 | 0.995 | |
| Magnesium | ND | 4.98 | 0.168 | 0.995 | |
| Potassium | ND | 24.9 | 1.75 | 0.995 | |
| Sodium | ND | 24.9 | 1.81 | 0.995 | |
| Strontium | ND | 1.49 | 0.137 | 0.995 | |
| Silicon | ND | 4.98 | 1.31 | 0.995 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB01 | 15-06-0567-10-F | 06/05/15 15:41 | Aqueous | ICP 7300 | 06/15/15 | 06/23/15 15:03 | 150615LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 0.0548 | 0.100 | 0.0118 | 1.00 | J |
| Magnesium | 0.0451 | 0.100 | 0.00336 | 1.00 | B,J |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | 0.150 | 0.500 | 0.103 | 1.00 | J |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|-----------------|-----|---------|----------|----------|----------------|------------|
| Method Blank | 099-15-683-1299 | N/A | Aqueous | ICP 7300 | 06/15/15 | 06/16/15 11:27 | 150615LA4F |
|--------------|-----------------|-----|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | ND | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 0.0372 | 0.100 | 0.00336 | 1.00 | J |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

Page 1 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM310 | 15-06-0567-1-BB | 06/05/15 10:03 | Solid | ICP/MS 03 | 06/12/15 | 06/12/15 18:19 | 150612L01 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|-------|------|------------|
| Antimony | ND | 6.97 | 0.454 | 1.00 | |
| Arsenic | 7.22 | 3.48 | 1.36 | 1.00 | |
| Barium | 70.1 | 3.48 | 0.390 | 1.00 | |
| Beryllium | 0.823 | 3.48 | 0.716 | 1.00 | J |
| Cadmium | 0.640 | 3.48 | 0.418 | 1.00 | J |
| Chromium | 23.4 | 6.97 | 0.407 | 1.00 | |
| Copper | 13.5 | 3.48 | 0.405 | 1.00 | |
| Lead | 21.6 | 3.48 | 0.478 | 1.00 | |
| Nickel | 15.8 | 3.48 | 0.446 | 1.00 | |
| Selenium | 2.15 | 3.48 | 0.920 | 1.00 | J |
| Silver | 0.506 | 3.48 | 0.495 | 1.00 | J |
| Thallium | ND | 3.48 | 0.366 | 1.00 | |
| Zinc | 95.1 | 17.4 | 2.76 | 1.00 | |
| Aluminum | 9790 | 87.1 | 11.3 | 1.00 | |
| Iron | 16300 | 87.1 | 14.3 | 1.00 | |
| Manganese | 401 | 8.71 | 0.455 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

Page 2 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM320 | 15-06-0567-2-BB | 06/05/15 10:48 | Solid | ICP/MS 03 | 06/12/15 | 06/12/15 18:22 | 150612L01 |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|-------|------|------------|
| Antimony | ND | 10.2 | 0.665 | 1.00 | |
| Arsenic | 7.16 | 5.10 | 1.99 | 1.00 | |
| Barium | 87.3 | 5.10 | 0.571 | 1.00 | |
| Beryllium | 1.16 | 5.10 | 1.05 | 1.00 | J |
| Cadmium | 1.17 | 5.10 | 0.612 | 1.00 | J |
| Chromium | 26.6 | 10.2 | 0.596 | 1.00 | |
| Copper | 58.3 | 5.10 | 0.594 | 1.00 | |
| Lead | 65.4 | 5.10 | 0.699 | 1.00 | |
| Nickel | 20.5 | 5.10 | 0.653 | 1.00 | |
| Selenium | 4.20 | 5.10 | 1.35 | 1.00 | J |
| Silver | 0.822 | 5.10 | 0.725 | 1.00 | J |
| Thallium | ND | 5.10 | 0.535 | 1.00 | |
| Zinc | 253 | 25.5 | 4.05 | 1.00 | |
| Aluminum | 12400 | 128 | 16.6 | 1.00 | |
| Iron | 24500 | 128 | 20.9 | 1.00 | |
| Manganese | 390 | 12.8 | 0.666 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

Page 3 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM330 | 15-06-0567-3-BB | 06/05/15 11:03 | Solid | ICP/MS 03 | 06/12/15 | 06/12/15 18:26 | 150612L01 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|-------|------|------------|
| Antimony | ND | 2.52 | 0.164 | 1.00 | |
| Arsenic | 5.99 | 1.26 | 0.490 | 1.00 | |
| Barium | 38.4 | 1.26 | 0.141 | 1.00 | |
| Beryllium | ND | 1.26 | 0.258 | 1.00 | |
| Cadmium | 0.292 | 1.26 | 0.151 | 1.00 | J |
| Chromium | 8.38 | 2.52 | 0.147 | 1.00 | |
| Copper | 3.84 | 1.26 | 0.146 | 1.00 | |
| Lead | 16.0 | 1.26 | 0.172 | 1.00 | |
| Nickel | 10.8 | 1.26 | 0.161 | 1.00 | |
| Selenium | ND | 1.26 | 0.332 | 1.00 | |
| Silver | ND | 1.26 | 0.179 | 1.00 | |
| Thallium | ND | 1.26 | 0.132 | 1.00 | |
| Zinc | 60.7 | 6.29 | 0.998 | 1.00 | |
| Aluminum | 1200 | 31.4 | 4.08 | 1.00 | |
| Iron | 9350 | 31.4 | 5.15 | 1.00 | |
| Manganese | 334 | 3.14 | 0.164 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM340 | 15-06-0567-4-BB | 06/05/15 11:21 | Solid | ICP/MS 03 | 06/12/15 | 06/12/15 18:29 | 150612L01 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|-------|------|------------|
| Antimony | 0.225 | 2.49 | 0.163 | 1.00 | J |
| Arsenic | 2.98 | 1.25 | 0.485 | 1.00 | |
| Barium | 31.0 | 1.25 | 0.139 | 1.00 | |
| Beryllium | ND | 1.25 | 0.256 | 1.00 | |
| Cadmium | 0.441 | 1.25 | 0.150 | 1.00 | J |
| Chromium | 17.0 | 2.49 | 0.146 | 1.00 | |
| Copper | 13.6 | 1.25 | 0.145 | 1.00 | |
| Lead | 63.5 | 1.25 | 0.171 | 1.00 | |
| Nickel | 9.60 | 1.25 | 0.160 | 1.00 | |
| Selenium | ND | 1.25 | 0.329 | 1.00 | |
| Silver | 0.221 | 1.25 | 0.177 | 1.00 | J |
| Thallium | ND | 1.25 | 0.131 | 1.00 | |
| Zinc | 61.2 | 6.23 | 0.989 | 1.00 | |
| Aluminum | 2840 | 31.2 | 4.05 | 1.00 | |
| Iron | 7640 | 31.2 | 5.10 | 1.00 | |
| Manganese | 152 | 3.12 | 0.163 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM350 | 15-06-0567-5-BB | 06/05/15 12:24 | Solid | ICP/MS 03 | 06/12/15 | 06/12/15 18:33 | 150612L01 |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|-------|------|------------|
| Antimony | ND | 4.36 | 0.284 | 1.00 | |
| Arsenic | 5.69 | 2.18 | 0.848 | 1.00 | |
| Barium | 45.4 | 2.18 | 0.244 | 1.00 | |
| Beryllium | 0.469 | 2.18 | 0.447 | 1.00 | J |
| Cadmium | 0.596 | 2.18 | 0.261 | 1.00 | J |
| Chromium | 15.3 | 4.36 | 0.254 | 1.00 | |
| Copper | 17.8 | 2.18 | 0.254 | 1.00 | |
| Lead | 35.5 | 2.18 | 0.299 | 1.00 | |
| Nickel | 14.5 | 2.18 | 0.279 | 1.00 | |
| Selenium | 1.18 | 2.18 | 0.575 | 1.00 | J |
| Silver | 0.384 | 2.18 | 0.310 | 1.00 | J |
| Thallium | ND | 2.18 | 0.229 | 1.00 | |
| Zinc | 84.6 | 10.9 | 1.73 | 1.00 | |
| Aluminum | 4890 | 54.5 | 7.07 | 1.00 | |
| Iron | 12100 | 54.5 | 8.91 | 1.00 | |
| Manganese | 343 | 5.45 | 0.284 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM360 | 15-06-0567-6-BB | 06/05/15 13:13 | Solid | ICP/MS 03 | 06/12/15 | 06/12/15 18:36 | 150612L01 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|-------|------|------------|
| Antimony | ND | 2.68 | 0.175 | 1.00 | |
| Arsenic | 4.11 | 1.34 | 0.523 | 1.00 | |
| Barium | 34.7 | 1.34 | 0.150 | 1.00 | |
| Beryllium | 0.343 | 1.34 | 0.276 | 1.00 | J |
| Cadmium | 0.427 | 1.34 | 0.161 | 1.00 | J |
| Chromium | 10.0 | 2.68 | 0.157 | 1.00 | |
| Copper | 5.74 | 1.34 | 0.156 | 1.00 | |
| Lead | 54.4 | 1.34 | 0.184 | 1.00 | |
| Nickel | 9.41 | 1.34 | 0.172 | 1.00 | |
| Selenium | 0.386 | 1.34 | 0.354 | 1.00 | J |
| Silver | 0.259 | 1.34 | 0.191 | 1.00 | J |
| Thallium | ND | 1.34 | 0.141 | 1.00 | |
| Zinc | 33.0 | 6.71 | 1.06 | 1.00 | |
| Aluminum | 4930 | 33.6 | 4.36 | 1.00 | |
| Iron | 9170 | 33.6 | 5.49 | 1.00 | |
| Manganese | 240 | 3.36 | 0.175 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM370 | 15-06-0567-7-BB | 06/05/15 13:51 | Solid | ICP/MS 03 | 06/12/15 | 06/12/15 18:40 | 150612L01 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|-------|------|------------|
| Antimony | ND | 3.42 | 0.223 | 1.00 | |
| Arsenic | 11.0 | 1.71 | 0.665 | 1.00 | |
| Barium | 42.7 | 1.71 | 0.191 | 1.00 | |
| Beryllium | 0.436 | 1.71 | 0.351 | 1.00 | J |
| Cadmium | 0.612 | 1.71 | 0.205 | 1.00 | J |
| Chromium | 11.6 | 3.42 | 0.200 | 1.00 | |
| Copper | 8.30 | 1.71 | 0.199 | 1.00 | |
| Lead | 17.9 | 1.71 | 0.234 | 1.00 | |
| Nickel | 18.0 | 1.71 | 0.219 | 1.00 | |
| Selenium | ND | 1.71 | 0.451 | 1.00 | |
| Silver | 0.282 | 1.71 | 0.243 | 1.00 | J |
| Thallium | ND | 1.71 | 0.179 | 1.00 | |
| Zinc | 42.6 | 8.55 | 1.36 | 1.00 | |
| Aluminum | 5760 | 42.7 | 5.55 | 1.00 | |
| Iron | 15300 | 42.7 | 6.99 | 1.00 | |
| Manganese | 439 | 4.27 | 0.223 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| FDHSM370 | 15-06-0567-8-BB | 06/05/15 13:51 | Solid | ICP/MS 03 | 06/12/15 | 06/15/15 12:49 | 150612L01 |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|-------|------|------------|
| Antimony | ND | 3.68 | 0.240 | 1.00 | |
| Arsenic | 8.30 | 1.84 | 0.717 | 1.00 | |
| Barium | 55.7 | 1.84 | 0.206 | 1.00 | |
| Beryllium | 0.540 | 1.84 | 0.378 | 1.00 | J |
| Cadmium | 0.560 | 1.84 | 0.221 | 1.00 | J |
| Chromium | 14.0 | 3.68 | 0.215 | 1.00 | |
| Copper | 8.41 | 1.84 | 0.214 | 1.00 | |
| Lead | 19.6 | 1.84 | 0.252 | 1.00 | |
| Nickel | 14.9 | 1.84 | 0.236 | 1.00 | |
| Selenium | 0.670 | 1.84 | 0.486 | 1.00 | J |
| Silver | ND | 1.84 | 0.262 | 1.00 | |
| Thallium | ND | 1.84 | 0.193 | 1.00 | |
| Zinc | 43.5 | 9.21 | 1.46 | 1.00 | |
| Aluminum | 8180 | 46.0 | 5.98 | 1.00 | |
| Iron | 13800 | 46.0 | 7.54 | 1.00 | |
| Manganese | 471 | 4.60 | 0.240 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6020
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-621-955 | N/A | Solid | ICP/MS 03 | 06/12/15 | 06/12/15 17:25 | 150612L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|------|-------|------|------------|
| Antimony | ND | 2.00 | 0.130 | 1.00 | |
| Arsenic | ND | 1.00 | 0.389 | 1.00 | |
| Barium | ND | 1.00 | 0.112 | 1.00 | |
| Beryllium | ND | 1.00 | 0.205 | 1.00 | |
| Cadmium | ND | 1.00 | 0.120 | 1.00 | |
| Chromium | ND | 2.00 | 0.117 | 1.00 | |
| Copper | ND | 1.00 | 0.116 | 1.00 | |
| Lead | ND | 1.00 | 0.137 | 1.00 | |
| Nickel | ND | 1.00 | 0.128 | 1.00 | |
| Selenium | ND | 1.00 | 0.264 | 1.00 | |
| Silver | ND | 1.00 | 0.142 | 1.00 | |
| Thallium | ND | 1.00 | 0.105 | 1.00 | |
| Zinc | ND | 5.00 | 0.793 | 1.00 | |
| Aluminum | ND | 25.0 | 3.25 | 1.00 | |
| Iron | ND | 25.0 | 4.09 | 1.00 | |
| Manganese | ND | 2.50 | 0.131 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB01 | 15-06-0567-10-F | 06/05/15 15:41 | Aqueous | ICP/MS 04 | 06/08/15 | 06/09/15 18:50 | 150608LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00104 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.000299 | 0.00100 | 0.000132 | 1.00 | J |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.000853 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.000158 | 0.00100 | 0.000139 | 1.00 | J |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-842 | N/A | Aqueous | ICP/MS 04 | 06/08/15 | 06/09/15 16:59 | 150608LA2F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB01 | 15-06-0567-10-F | 06/05/15 15:41 | Aqueous | Mercury 04 | 06/08/15 | 06/08/15 19:07 | 150608L01F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|-----|---------|------------|----------|-------------------|------------|
| Method Blank | 099-15-763-567 | N/A | Aqueous | Mercury 04 | 06/08/15 | 06/08/15 13:54 | 150608L01F |
|--------------|----------------|-----|---------|------------|----------|-------------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 7471A Total
Method: EPA 7471A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|
| HSM310 | 15-06-0567-1-BB | 06/05/15 10:03 | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:27 | 150617L03 |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | 0.0574 | 0.277 | 0.0195 | 1.00 | J |

| | | | | | | | |
|---------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|
| HSM320 | 15-06-0567-2-BB | 06/05/15 10:48 | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:30 | 150617L03 |
|---------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | 0.0735 | 0.418 | 0.0295 | 1.00 | J |

| | | | | | | | |
|---------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|
| HSM330 | 15-06-0567-3-BB | 06/05/15 11:03 | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:32 | 150617L03 |
|---------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | 0.0996 | 0.103 | 0.00726 | 1.00 | J |

| | | | | | | | |
|---------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|
| HSM340 | 15-06-0567-4-BB | 06/05/15 11:21 | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:38 | 150617L03 |
|---------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | 0.0215 | 0.0990 | 0.00697 | 1.00 | J |

| | | | | | | | |
|---------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|
| HSM350 | 15-06-0567-5-BB | 06/05/15 12:24 | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:41 | 150617L03 |
|---------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | 0.0558 | 0.179 | 0.0126 | 1.00 | J |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 7471A Total
Method: EPA 7471A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|
| HSM360 | 15-06-0567-6-BB | 06/05/15 13:13 | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:43 | 150617L03 |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | 0.0173 | 0.116 | 0.00815 | 1.00 | J |

| | | | | | | | |
|---------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|
| HSM370 | 15-06-0567-7-BB | 06/05/15 13:51 | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:45 | 150617L03 |
|---------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | 0.0166 | 0.145 | 0.0102 | 1.00 | J |

| | | | | | | | |
|-----------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|
| FDHSM370 | 15-06-0567-8-BB | 06/05/15 13:51 | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:47 | 150617L03 |
|-----------------|------------------------|-----------------------|--------------|-------------------|-----------------|-----------------------|------------------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | 0.0268 | 0.151 | 0.0106 | 1.00 | J |

| | | | | | | | |
|---------------------|------------------------|------------|--------------|-------------------|-----------------|-----------------------|------------------|
| Method Blank | 099-16-272-1355 | N/A | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:03 | 150617L03 |
|---------------------|------------------------|------------|--------------|-------------------|-----------------|-----------------------|------------------|

Comment(s):
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.0833 | 0.00587 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM310 | 15-06-0567-1-B | 06/05/15 10:03 | Solid | GC 41 | 06/10/15 | 06/17/15 14:02 | 150610L17 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-----|-----|------|------------|
| Aldrin | ND | 17 | 7.6 | 1.00 | |
| Alpha-BHC | ND | 34 | 13 | 1.00 | |
| Beta-BHC | ND | 17 | 8.6 | 1.00 | |
| Chlordane | ND | 170 | 90 | 1.00 | |
| 4,4'-DDD | ND | 17 | 8.2 | 1.00 | |
| 4,4'-DDE | ND | 17 | 7.7 | 1.00 | |
| 4,4'-DDT | ND | 17 | 7.5 | 1.00 | |
| Delta-BHC | ND | 34 | 15 | 1.00 | |
| Dieldrin | ND | 17 | 7.6 | 1.00 | |
| Endosulfan I | ND | 17 | 6.8 | 1.00 | |
| Endosulfan II | ND | 17 | 8.1 | 1.00 | |
| Endosulfan Sulfate | ND | 17 | 9.0 | 1.00 | |
| Endrin | ND | 17 | 8.1 | 1.00 | |
| Endrin Aldehyde | ND | 17 | 10 | 1.00 | |
| Endrin Ketone | ND | 17 | 8.7 | 1.00 | |
| Gamma-BHC | ND | 17 | 7.7 | 1.00 | |
| Heptachlor | ND | 17 | 7.5 | 1.00 | |
| Heptachlor Epoxide | ND | 34 | 13 | 1.00 | |
| Methoxychlor | ND | 17 | 9.4 | 1.00 | |
| Toxaphene | ND | 340 | 150 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 149 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 97 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM320 | 15-06-0567-2-B | 06/05/15 10:48 | Solid | GC 41 | 06/10/15 | 06/17/15 14:17 | 150610L17 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-----|-----|------|------------|
| Aldrin | ND | 25 | 11 | 1.00 | |
| Alpha-BHC | ND | 51 | 19 | 1.00 | |
| Beta-BHC | ND | 25 | 13 | 1.00 | |
| Chlordane | ND | 250 | 130 | 1.00 | |
| 4,4'-DDD | ND | 25 | 12 | 1.00 | |
| 4,4'-DDE | 17 | 25 | 11 | 1.00 | J |
| 4,4'-DDT | ND | 25 | 11 | 1.00 | |
| Delta-BHC | ND | 51 | 22 | 1.00 | |
| Dieldrin | ND | 25 | 11 | 1.00 | |
| Endosulfan I | ND | 25 | 10 | 1.00 | |
| Endosulfan II | ND | 25 | 12 | 1.00 | |
| Endosulfan Sulfate | ND | 25 | 13 | 1.00 | |
| Endrin | ND | 25 | 12 | 1.00 | |
| Endrin Aldehyde | ND | 25 | 15 | 1.00 | |
| Endrin Ketone | ND | 25 | 13 | 1.00 | |
| Gamma-BHC | ND | 25 | 11 | 1.00 | |
| Heptachlor | ND | 25 | 11 | 1.00 | |
| Heptachlor Epoxide | ND | 51 | 19 | 1.00 | |
| Methoxychlor | ND | 25 | 14 | 1.00 | |
| Toxaphene | ND | 510 | 230 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 106 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 96 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM330 | 15-06-0567-3-B | 06/05/15 11:03 | Solid | GC 41 | 06/10/15 | 06/17/15 14:33 | 150610L17 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-----|-----|------|------------|
| Aldrin | ND | 6.3 | 2.7 | 1.00 | |
| Alpha-BHC | ND | 13 | 4.6 | 1.00 | |
| Beta-BHC | ND | 6.3 | 3.1 | 1.00 | |
| Chlordane | ND | 63 | 33 | 1.00 | |
| 4,4'-DDD | ND | 6.3 | 3.0 | 1.00 | |
| 4,4'-DDE | ND | 6.3 | 2.8 | 1.00 | |
| 4,4'-DDT | ND | 6.3 | 2.7 | 1.00 | |
| Delta-BHC | ND | 13 | 5.5 | 1.00 | |
| Dieldrin | ND | 6.3 | 2.7 | 1.00 | |
| Endosulfan I | ND | 6.3 | 2.5 | 1.00 | |
| Endosulfan II | ND | 6.3 | 2.9 | 1.00 | |
| Endosulfan Sulfate | ND | 6.3 | 3.3 | 1.00 | |
| Endrin | ND | 6.3 | 2.9 | 1.00 | |
| Endrin Aldehyde | ND | 6.3 | 3.8 | 1.00 | |
| Endrin Ketone | ND | 6.3 | 3.1 | 1.00 | |
| Gamma-BHC | ND | 6.3 | 2.8 | 1.00 | |
| Heptachlor | ND | 6.3 | 2.7 | 1.00 | |
| Heptachlor Epoxide | ND | 13 | 4.6 | 1.00 | |
| Methoxychlor | ND | 6.3 | 3.4 | 1.00 | |
| Toxaphene | ND | 130 | 56 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 101 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 76 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM340 | 15-06-0567-4-B | 06/05/15 11:21 | Solid | GC 41 | 06/10/15 | 06/17/15 14:48 | 150610L17 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-----|-----|------|------------|
| Aldrin | ND | 6.2 | 2.7 | 1.00 | |
| Alpha-BHC | ND | 12 | 4.6 | 1.00 | |
| Beta-BHC | ND | 6.2 | 3.1 | 1.00 | |
| Chlordane | ND | 62 | 32 | 1.00 | |
| 4,4'-DDD | ND | 6.2 | 2.9 | 1.00 | |
| 4,4'-DDE | 9.2 | 6.2 | 2.7 | 1.00 | |
| 4,4'-DDT | ND | 6.2 | 2.7 | 1.00 | |
| Delta-BHC | ND | 12 | 5.4 | 1.00 | |
| Dieldrin | ND | 6.2 | 2.7 | 1.00 | |
| Endosulfan I | ND | 6.2 | 2.4 | 1.00 | |
| Endosulfan II | ND | 6.2 | 2.9 | 1.00 | |
| Endosulfan Sulfate | ND | 6.2 | 3.2 | 1.00 | |
| Endrin | ND | 6.2 | 2.9 | 1.00 | |
| Endrin Aldehyde | ND | 6.2 | 3.7 | 1.00 | |
| Endrin Ketone | ND | 6.2 | 3.1 | 1.00 | |
| Gamma-BHC | ND | 6.2 | 2.8 | 1.00 | |
| Heptachlor | ND | 6.2 | 2.7 | 1.00 | |
| Heptachlor Epoxide | ND | 12 | 4.6 | 1.00 | |
| Methoxychlor | ND | 6.2 | 3.4 | 1.00 | |
| Toxaphene | ND | 120 | 55 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 130 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 89 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM350 | 15-06-0567-5-B | 06/05/15 12:24 | Solid | GC 41 | 06/10/15 | 06/17/15 15:03 | 150610L17 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-----|-----|------|------------|
| Aldrin | ND | 11 | 4.7 | 1.00 | |
| Alpha-BHC | ND | 22 | 8.0 | 1.00 | |
| Beta-BHC | ND | 11 | 5.4 | 1.00 | |
| Chlordane | ND | 110 | 56 | 1.00 | |
| 4,4'-DDD | ND | 11 | 5.1 | 1.00 | |
| 4,4'-DDE | ND | 11 | 4.8 | 1.00 | |
| 4,4'-DDT | ND | 11 | 4.7 | 1.00 | |
| Delta-BHC | ND | 22 | 9.5 | 1.00 | |
| Dieldrin | ND | 11 | 4.7 | 1.00 | |
| Endosulfan I | ND | 11 | 4.3 | 1.00 | |
| Endosulfan II | ND | 11 | 5.1 | 1.00 | |
| Endosulfan Sulfate | ND | 11 | 5.6 | 1.00 | |
| Endrin | ND | 11 | 5.1 | 1.00 | |
| Endrin Aldehyde | ND | 11 | 6.5 | 1.00 | |
| Endrin Ketone | ND | 11 | 5.4 | 1.00 | |
| Gamma-BHC | ND | 11 | 4.8 | 1.00 | |
| Heptachlor | ND | 11 | 4.7 | 1.00 | |
| Heptachlor Epoxide | ND | 22 | 8.0 | 1.00 | |
| Methoxychlor | ND | 11 | 5.9 | 1.00 | |
| Toxaphene | ND | 220 | 97 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 197 | 24-168 | 2,7 |
| 2,4,5,6-Tetrachloro-m-Xylene | 97 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM360 | 15-06-0567-6-B | 06/05/15 13:13 | Solid | GC 41 | 06/10/15 | 06/17/15 15:18 | 150610L17 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-----|-----|------|------------|
| Aldrin | ND | 6.7 | 2.9 | 1.00 | |
| Alpha-BHC | ND | 13 | 5.0 | 1.00 | |
| Beta-BHC | ND | 6.7 | 3.3 | 1.00 | |
| Chlordane | ND | 67 | 35 | 1.00 | |
| 4,4'-DDD | ND | 6.7 | 3.2 | 1.00 | |
| 4,4'-DDE | ND | 6.7 | 3.0 | 1.00 | |
| 4,4'-DDT | ND | 6.7 | 2.9 | 1.00 | |
| Delta-BHC | ND | 13 | 5.9 | 1.00 | |
| Dieldrin | ND | 6.7 | 2.9 | 1.00 | |
| Endosulfan I | ND | 6.7 | 2.7 | 1.00 | |
| Endosulfan II | ND | 6.7 | 3.2 | 1.00 | |
| Endosulfan Sulfate | ND | 6.7 | 3.5 | 1.00 | |
| Endrin | ND | 6.7 | 3.2 | 1.00 | |
| Endrin Aldehyde | ND | 6.7 | 4.1 | 1.00 | |
| Endrin Ketone | ND | 6.7 | 3.4 | 1.00 | |
| Gamma-BHC | ND | 6.7 | 3.0 | 1.00 | |
| Heptachlor | ND | 6.7 | 2.9 | 1.00 | |
| Heptachlor Epoxide | ND | 13 | 5.0 | 1.00 | |
| Methoxychlor | ND | 6.7 | 3.6 | 1.00 | |
| Toxaphene | ND | 130 | 60 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 95 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 94 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM370 | 15-06-0567-7-B | 06/05/15 13:51 | Solid | GC 41 | 06/10/15 | 06/17/15 15:34 | 150610L17 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-----|-----|------|------------|
| Aldrin | ND | 8.5 | 3.7 | 1.00 | |
| Alpha-BHC | ND | 17 | 6.3 | 1.00 | |
| Beta-BHC | ND | 8.5 | 4.2 | 1.00 | |
| Chlordane | ND | 85 | 44 | 1.00 | |
| 4,4'-DDD | ND | 8.5 | 4.0 | 1.00 | |
| 4,4'-DDE | ND | 8.5 | 3.8 | 1.00 | |
| 4,4'-DDT | ND | 8.5 | 3.7 | 1.00 | |
| Delta-BHC | ND | 17 | 7.5 | 1.00 | |
| Dieldrin | ND | 8.5 | 3.7 | 1.00 | |
| Endosulfan I | ND | 8.5 | 3.4 | 1.00 | |
| Endosulfan II | ND | 8.5 | 4.0 | 1.00 | |
| Endosulfan Sulfate | ND | 8.5 | 4.4 | 1.00 | |
| Endrin | ND | 8.5 | 4.0 | 1.00 | |
| Endrin Aldehyde | ND | 8.5 | 5.1 | 1.00 | |
| Endrin Ketone | ND | 8.5 | 4.3 | 1.00 | |
| Gamma-BHC | ND | 8.5 | 3.8 | 1.00 | |
| Heptachlor | ND | 8.5 | 3.7 | 1.00 | |
| Heptachlor Epoxide | ND | 17 | 6.3 | 1.00 | |
| Methoxychlor | ND | 8.5 | 4.6 | 1.00 | |
| Toxaphene | ND | 170 | 76 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 94 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 102 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

Page 8 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| FDHSM370 | 15-06-0567-8-B | 06/05/15 13:51 | Solid | GC 41 | 06/10/15 | 06/17/15 15:49 | 150610L17 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-----|-----|------|------------|
| Aldrin | ND | 9.2 | 4.0 | 1.00 | |
| Alpha-BHC | ND | 18 | 6.8 | 1.00 | |
| Beta-BHC | ND | 9.2 | 4.6 | 1.00 | |
| Chlordane | ND | 92 | 48 | 1.00 | |
| 4,4'-DDD | ND | 9.2 | 4.4 | 1.00 | |
| 4,4'-DDE | ND | 9.2 | 4.1 | 1.00 | |
| 4,4'-DDT | ND | 9.2 | 4.0 | 1.00 | |
| Delta-BHC | ND | 18 | 8.1 | 1.00 | |
| Dieldrin | ND | 9.2 | 4.0 | 1.00 | |
| Endosulfan I | ND | 9.2 | 3.7 | 1.00 | |
| Endosulfan II | ND | 9.2 | 4.3 | 1.00 | |
| Endosulfan Sulfate | ND | 9.2 | 4.8 | 1.00 | |
| Endrin | ND | 9.2 | 4.3 | 1.00 | |
| Endrin Aldehyde | ND | 9.2 | 5.6 | 1.00 | |
| Endrin Ketone | ND | 9.2 | 4.6 | 1.00 | |
| Gamma-BHC | ND | 9.2 | 4.1 | 1.00 | |
| Heptachlor | ND | 9.2 | 4.0 | 1.00 | |
| Heptachlor Epoxide | ND | 18 | 6.8 | 1.00 | |
| Methoxychlor | ND | 9.2 | 5.0 | 1.00 | |
| Toxaphene | ND | 180 | 83 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 96 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 94 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8081A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-537-2135 | N/A | Solid | GC 41 | 06/10/15 | 06/11/15 18:47 | 150610L17 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-----|-----|------|------------|
| Aldrin | ND | 5.0 | 2.2 | 1.00 | |
| Alpha-BHC | ND | 10 | 3.7 | 1.00 | |
| Beta-BHC | ND | 5.0 | 2.5 | 1.00 | |
| Chlordane | ND | 50 | 26 | 1.00 | |
| 4,4'-DDD | ND | 5.0 | 2.4 | 1.00 | |
| 4,4'-DDE | ND | 5.0 | 2.2 | 1.00 | |
| 4,4'-DDT | ND | 5.0 | 2.2 | 1.00 | |
| Delta-BHC | ND | 10 | 4.4 | 1.00 | |
| Dieldrin | ND | 5.0 | 2.2 | 1.00 | |
| Endosulfan I | ND | 5.0 | 2.0 | 1.00 | |
| Endosulfan II | ND | 5.0 | 2.4 | 1.00 | |
| Endosulfan Sulfate | ND | 5.0 | 2.6 | 1.00 | |
| Endrin | ND | 5.0 | 2.4 | 1.00 | |
| Endrin Aldehyde | ND | 5.0 | 3.0 | 1.00 | |
| Endrin Ketone | ND | 5.0 | 2.5 | 1.00 | |
| Gamma-BHC | ND | 5.0 | 2.2 | 1.00 | |
| Heptachlor | ND | 5.0 | 2.2 | 1.00 | |
| Heptachlor Epoxide | ND | 10 | 3.7 | 1.00 | |
| Methoxychlor | ND | 5.0 | 2.7 | 1.00 | |
| Toxaphene | ND | 100 | 45 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 107 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 79 | 25-145 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

Page 1 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB01 | 15-06-0567-10-I | 06/05/15 15:41 | Aqueous | GC 44 | 06/08/15 | 06/10/15 15:08 | 150608L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 102 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 111 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

Page 2 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-813 | N/A | Aqueous | GC 44 | 06/08/15 | 06/10/15 14:26 | 150608L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 102 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 111 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8082
Units: ug/kg

Project: EAA 27122

Page 1 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM310 | 15-06-0567-1-B | 06/05/15 10:03 | Solid | GC 58 | 06/10/15 | 06/13/15 11:49 | 150610L18 |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|-----|------|------------|
| Aroclor-1016 | ND | 170 | 72 | 1.00 | |
| Aroclor-1221 | ND | 170 | 150 | 1.00 | |
| Aroclor-1232 | ND | 170 | 87 | 1.00 | |
| Aroclor-1242 | ND | 170 | 130 | 1.00 | |
| Aroclor-1248 | ND | 170 | 110 | 1.00 | |
| Aroclor-1254 | ND | 170 | 110 | 1.00 | |
| Aroclor-1260 | ND | 170 | 100 | 1.00 | |
| Aroclor-1262 | ND | 170 | 120 | 1.00 | |
| Aroclor-1268 | ND | 170 | 120 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 141 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 112 | 25-145 | |

| | | | | | | | |
|--------|----------------|-------------------|-------|-------|----------|-------------------|-----------|
| HSM320 | 15-06-0567-2-B | 06/05/15 10:48 | Solid | GC 58 | 06/10/15 | 06/13/15 14:36 | 150610L18 |
|--------|----------------|-------------------|-------|-------|----------|-------------------|-----------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|-----|------|------------|
| Aroclor-1016 | ND | 250 | 110 | 1.00 | |
| Aroclor-1221 | ND | 250 | 210 | 1.00 | |
| Aroclor-1232 | ND | 250 | 130 | 1.00 | |
| Aroclor-1242 | ND | 250 | 190 | 1.00 | |
| Aroclor-1248 | ND | 250 | 160 | 1.00 | |
| Aroclor-1254 | ND | 250 | 160 | 1.00 | |
| Aroclor-1260 | ND | 250 | 150 | 1.00 | |
| Aroclor-1262 | ND | 250 | 180 | 1.00 | |
| Aroclor-1268 | ND | 250 | 170 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 122 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 105 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8082
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM330 | 15-06-0567-3-B | 06/05/15 11:03 | Solid | GC 58 | 06/10/15 | 06/13/15 12:25 | 150610L18 |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|----|-----|------|------------|
| Aroclor-1016 | ND | 63 | 26 | 1.00 | |
| Aroclor-1221 | ND | 63 | 53 | 1.00 | |
| Aroclor-1232 | ND | 63 | 31 | 1.00 | |
| Aroclor-1242 | ND | 63 | 46 | 1.00 | |
| Aroclor-1248 | ND | 63 | 40 | 1.00 | |
| Aroclor-1254 | ND | 63 | 40 | 1.00 | |
| Aroclor-1260 | ND | 63 | 38 | 1.00 | |
| Aroclor-1262 | ND | 63 | 43 | 1.00 | |
| Aroclor-1268 | ND | 63 | 42 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 98 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 85 | 25-145 | |

| | | | | | | | |
|--------|----------------|-------------------|-------|-------|----------|-------------------|-----------|
| HSM340 | 15-06-0567-4-B | 06/05/15 11:21 | Solid | GC 58 | 06/10/15 | 06/13/15 12:43 | 150610L18 |
|--------|----------------|-------------------|-------|-------|----------|-------------------|-----------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|----|-----|------|------------|
| Aroclor-1016 | ND | 62 | 26 | 1.00 | |
| Aroclor-1221 | ND | 62 | 52 | 1.00 | |
| Aroclor-1232 | ND | 62 | 31 | 1.00 | |
| Aroclor-1242 | ND | 62 | 46 | 1.00 | |
| Aroclor-1248 | ND | 62 | 39 | 1.00 | |
| Aroclor-1254 | 41 | 62 | 39 | 1.00 | J |
| Aroclor-1260 | 52 | 62 | 38 | 1.00 | J |
| Aroclor-1262 | ND | 62 | 43 | 1.00 | |
| Aroclor-1268 | ND | 62 | 41 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 136 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 100 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8082
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM350 | 15-06-0567-5-B | 06/05/15 12:24 | Solid | GC 58 | 06/10/15 | 06/13/15 13:01 | 150610L18 |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|-----|------|------------|
| Aroclor-1016 | ND | 110 | 45 | 1.00 | |
| Aroclor-1221 | ND | 110 | 91 | 1.00 | |
| Aroclor-1232 | ND | 110 | 54 | 1.00 | |
| Aroclor-1242 | ND | 110 | 80 | 1.00 | |
| Aroclor-1248 | ND | 110 | 69 | 1.00 | |
| Aroclor-1254 | ND | 110 | 68 | 1.00 | |
| Aroclor-1260 | ND | 110 | 66 | 1.00 | |
| Aroclor-1262 | ND | 110 | 75 | 1.00 | |
| Aroclor-1268 | ND | 110 | 72 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 136 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 121 | 25-145 | |

| | | | | | | | |
|--------|----------------|-------------------|-------|-------|----------|-------------------|-----------|
| HSM360 | 15-06-0567-6-B | 06/05/15 13:13 | Solid | GC 58 | 06/10/15 | 06/13/15 13:23 | 150610L18 |
|--------|----------------|-------------------|-------|-------|----------|-------------------|-----------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|----|-----|------|------------|
| Aroclor-1016 | ND | 67 | 28 | 1.00 | |
| Aroclor-1221 | ND | 67 | 57 | 1.00 | |
| Aroclor-1232 | ND | 67 | 34 | 1.00 | |
| Aroclor-1242 | ND | 67 | 50 | 1.00 | |
| Aroclor-1248 | ND | 67 | 43 | 1.00 | |
| Aroclor-1254 | ND | 67 | 43 | 1.00 | |
| Aroclor-1260 | ND | 67 | 41 | 1.00 | |
| Aroclor-1262 | ND | 67 | 47 | 1.00 | |
| Aroclor-1268 | ND | 67 | 45 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 134 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 105 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8082
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM370 | 15-06-0567-7-B | 06/05/15 13:51 | Solid | GC 58 | 06/10/15 | 06/13/15 13:41 | 150610L18 |

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|----|-----|------|------------|
| Aroclor-1016 | ND | 85 | 35 | 1.00 | |
| Aroclor-1221 | ND | 85 | 72 | 1.00 | |
| Aroclor-1232 | ND | 85 | 43 | 1.00 | |
| Aroclor-1242 | ND | 85 | 63 | 1.00 | |
| Aroclor-1248 | ND | 85 | 54 | 1.00 | |
| Aroclor-1254 | ND | 85 | 54 | 1.00 | |
| Aroclor-1260 | ND | 85 | 52 | 1.00 | |
| Aroclor-1262 | ND | 85 | 59 | 1.00 | |
| Aroclor-1268 | ND | 85 | 57 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 136 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 116 | 25-145 | |

| FDHSM370 | 15-06-0567-8-B | 06/05/15 13:51 | Solid | GC 58 | 06/10/15 | 06/13/15 13:59 | 150610L18 |
|----------|----------------|-------------------|-------|-------|----------|-------------------|-----------|
|----------|----------------|-------------------|-------|-------|----------|-------------------|-----------|

Comment(s):
- Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|----|-----|------|------------|
| Aroclor-1016 | ND | 92 | 38 | 1.00 | |
| Aroclor-1221 | ND | 92 | 78 | 1.00 | |
| Aroclor-1232 | ND | 92 | 46 | 1.00 | |
| Aroclor-1242 | ND | 92 | 68 | 1.00 | |
| Aroclor-1248 | ND | 92 | 59 | 1.00 | |
| Aroclor-1254 | ND | 92 | 58 | 1.00 | |
| Aroclor-1260 | ND | 92 | 56 | 1.00 | |
| Aroclor-1262 | ND | 92 | 64 | 1.00 | |
| Aroclor-1268 | ND | 92 | 62 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 130 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 103 | 25-145 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8082
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-535-3266 | N/A | Solid | GC 58 | 06/10/15 | 06/11/15 22:22 | 150610L18 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|----|-----|------|------------|
| Aroclor-1016 | ND | 50 | 21 | 1.00 | |
| Aroclor-1221 | ND | 50 | 42 | 1.00 | |
| Aroclor-1232 | ND | 50 | 25 | 1.00 | |
| Aroclor-1242 | ND | 50 | 37 | 1.00 | |
| Aroclor-1248 | ND | 50 | 32 | 1.00 | |
| Aroclor-1254 | ND | 50 | 32 | 1.00 | |
| Aroclor-1260 | ND | 50 | 30 | 1.00 | |
| Aroclor-1262 | ND | 50 | 35 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 150 | 24-168 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 117 | 25-145 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB01 | 15-06-0567-10-I | 06/05/15 15:41 | Aqueous | GC 31 | 06/08/15 | 06/09/15 19:39 | 150608L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 107 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 111 | 50-135 | |

| Method Blank | 099-12-533-1050 | N/A | Aqueous | GC 31 | 06/08/15 | 06/09/15 19:20 | 150608L03 |
|--------------|-----------------|-----|---------|-------|----------|-------------------|-----------|
|--------------|-----------------|-----|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 107 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 108 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 1 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB01 | 15-06-0567-10-I | 06/05/15 15:41 | Aqueous | GC 26 | 06/08/15 | 06/10/15 15:05 | 150608L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 103 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-97 | N/A | Aqueous | GC 26 | 06/08/15 | 06/10/15 12:08 | 150608L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 97 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

Page 1 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM310 | 15-06-0567-1-B | 06/05/15 10:03 | Solid | GC 26 | 06/09/15 | 06/15/15 10:41 | 150609L20 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|-----|------|------|------------|
| Demeton-o/s | ND | 1.7 | 0.98 | 1.00 | |
| Azinphos Methyl | ND | 1.7 | 0.78 | 1.00 | |
| Bolstar | ND | 1.7 | 0.61 | 1.00 | |
| Chlorpyrifos | ND | 1.7 | 0.57 | 1.00 | |
| Coumaphos | ND | 1.7 | 0.52 | 1.00 | |
| Diazinon | ND | 1.7 | 0.55 | 1.00 | |
| Dichlorvos | ND | 1.7 | 0.95 | 1.00 | |
| Disulfoton | ND | 1.7 | 0.53 | 1.00 | |
| Ethoprop | ND | 1.7 | 0.57 | 1.00 | |
| Fensulfothion | ND | 1.7 | 0.62 | 1.00 | |
| Fenthion | ND | 1.7 | 0.64 | 1.00 | |
| Merphos | ND | 1.7 | 0.66 | 1.00 | |
| Methyl Parathion | ND | 1.7 | 0.73 | 1.00 | |
| Mevinphos | ND | 1.7 | 0.83 | 1.00 | |
| Naled | ND | 14 | 6.0 | 1.00 | |
| Phorate | ND | 1.7 | 0.78 | 1.00 | |
| Ronnel | ND | 1.7 | 0.54 | 1.00 | |
| Stirophos | ND | 6.9 | 2.1 | 1.00 | |
| Tokuthion | ND | 1.7 | 0.58 | 1.00 | |
| Trichloronate | ND | 1.7 | 0.52 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 91 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM320 | 15-06-0567-2-B | 06/05/15 10:48 | Solid | GC 26 | 06/09/15 | 06/15/15 11:25 | 150609L20 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|-----|------|------|------------|
| Demeton-o/s | ND | 2.5 | 1.4 | 1.00 | |
| Azinphos Methyl | ND | 2.5 | 1.1 | 1.00 | |
| Bolstar | ND | 2.5 | 0.89 | 1.00 | |
| Chlorpyrifos | ND | 2.5 | 0.83 | 1.00 | |
| Coumaphos | ND | 2.5 | 0.75 | 1.00 | |
| Diazinon | ND | 2.5 | 0.80 | 1.00 | |
| Dichlorvos | ND | 2.5 | 1.4 | 1.00 | |
| Disulfoton | ND | 2.5 | 0.76 | 1.00 | |
| Ethoprop | ND | 2.5 | 0.83 | 1.00 | |
| Fensulfothion | ND | 2.5 | 0.91 | 1.00 | |
| Fenthion | ND | 2.5 | 0.93 | 1.00 | |
| Merphos | ND | 2.5 | 0.96 | 1.00 | |
| Methyl Parathion | ND | 2.5 | 1.1 | 1.00 | |
| Mevinphos | ND | 2.5 | 1.2 | 1.00 | |
| Naled | ND | 20 | 8.6 | 1.00 | |
| Phorate | ND | 2.5 | 1.1 | 1.00 | |
| Ronnel | ND | 2.5 | 0.78 | 1.00 | |
| Stirophos | ND | 10 | 3.0 | 1.00 | |
| Tokuthion | ND | 2.5 | 0.84 | 1.00 | |
| Trichloronate | ND | 2.5 | 0.76 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 92 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM330 | 15-06-0567-3-B | 06/05/15 11:03 | Solid | GC 26 | 06/09/15 | 06/15/15 12:09 | 150609L20 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|------|------|------|------------|
| Demeton-o/s | ND | 0.61 | 0.35 | 1.00 | |
| Azinphos Methyl | ND | 0.61 | 0.28 | 1.00 | |
| Bolstar | ND | 0.61 | 0.22 | 1.00 | |
| Chlorpyrifos | ND | 0.61 | 0.20 | 1.00 | |
| Coumaphos | ND | 0.61 | 0.18 | 1.00 | |
| Diazinon | ND | 0.61 | 0.20 | 1.00 | |
| Dichlorvos | ND | 0.61 | 0.34 | 1.00 | |
| Disulfoton | ND | 0.61 | 0.19 | 1.00 | |
| Ethoprop | ND | 0.61 | 0.20 | 1.00 | |
| Fensulfothion | ND | 0.61 | 0.22 | 1.00 | |
| Fenthion | ND | 0.61 | 0.23 | 1.00 | |
| Merphos | ND | 0.61 | 0.24 | 1.00 | |
| Methyl Parathion | ND | 0.61 | 0.26 | 1.00 | |
| Mevinphos | ND | 0.61 | 0.30 | 1.00 | |
| Naled | ND | 4.9 | 2.1 | 1.00 | |
| Phorate | ND | 0.61 | 0.28 | 1.00 | |
| Ronnel | ND | 0.61 | 0.19 | 1.00 | |
| Stirophos | ND | 2.5 | 0.73 | 1.00 | |
| Tokuthion | ND | 0.61 | 0.21 | 1.00 | |
| Trichloronate | ND | 0.61 | 0.19 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 97 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM340 | 15-06-0567-4-B | 06/05/15 11:21 | Solid | GC 26 | 06/09/15 | 06/15/15 12:53 | 150609L20 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|------|------|------|------------|
| Demeton-o/s | ND | 0.62 | 0.35 | 1.00 | |
| Azinphos Methyl | ND | 0.62 | 0.28 | 1.00 | |
| Bolstar | ND | 0.62 | 0.22 | 1.00 | |
| Chlorpyrifos | ND | 0.62 | 0.21 | 1.00 | |
| Coumaphos | ND | 0.62 | 0.18 | 1.00 | |
| Diazinon | ND | 0.62 | 0.20 | 1.00 | |
| Dichlorvos | ND | 0.62 | 0.34 | 1.00 | |
| Disulfoton | ND | 0.62 | 0.19 | 1.00 | |
| Ethoprop | ND | 0.62 | 0.21 | 1.00 | |
| Fensulfothion | ND | 0.62 | 0.22 | 1.00 | |
| Fenthion | ND | 0.62 | 0.23 | 1.00 | |
| Merphos | ND | 0.62 | 0.24 | 1.00 | |
| Methyl Parathion | ND | 0.62 | 0.26 | 1.00 | |
| Mevinphos | ND | 0.62 | 0.30 | 1.00 | |
| Naled | ND | 4.9 | 2.1 | 1.00 | |
| Phorate | ND | 0.62 | 0.28 | 1.00 | |
| Ronnel | ND | 0.62 | 0.19 | 1.00 | |
| Stirophos | ND | 2.5 | 0.74 | 1.00 | |
| Tokuthion | ND | 0.62 | 0.21 | 1.00 | |
| Trichloronate | ND | 0.62 | 0.19 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 98 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM350 | 15-06-0567-5-B | 06/05/15 12:24 | Solid | GC 26 | 06/09/15 | 06/15/15 13:38 | 150609L20 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|-----|------|------|------------|
| Demeton-o/s | ND | 1.1 | 0.62 | 1.00 | |
| Azinphos Methyl | ND | 1.1 | 0.49 | 1.00 | |
| Bolstar | ND | 1.1 | 0.39 | 1.00 | |
| Chlorpyrifos | ND | 1.1 | 0.36 | 1.00 | |
| Coumaphos | ND | 1.1 | 0.33 | 1.00 | |
| Diazinon | ND | 1.1 | 0.35 | 1.00 | |
| Dichlorvos | ND | 1.1 | 0.60 | 1.00 | |
| Disulfoton | ND | 1.1 | 0.33 | 1.00 | |
| Ethoprop | ND | 1.1 | 0.36 | 1.00 | |
| Fensulfothion | ND | 1.1 | 0.39 | 1.00 | |
| Fenthion | ND | 1.1 | 0.41 | 1.00 | |
| Merphos | ND | 1.1 | 0.42 | 1.00 | |
| Methyl Parathion | ND | 1.1 | 0.46 | 1.00 | |
| Mevinphos | ND | 1.1 | 0.53 | 1.00 | |
| Naled | ND | 8.7 | 3.8 | 1.00 | |
| Phorate | ND | 1.1 | 0.49 | 1.00 | |
| Ronnel | ND | 1.1 | 0.34 | 1.00 | |
| Stirophos | ND | 4.4 | 1.3 | 1.00 | |
| Tokuthion | ND | 1.1 | 0.37 | 1.00 | |
| Trichloronate | ND | 1.1 | 0.33 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 110 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM360 | 15-06-0567-6-B | 06/05/15 13:13 | Solid | GC 26 | 06/09/15 | 06/15/15 14:22 | 150609L20 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|------|------|------|------------|
| Demeton-o/s | ND | 0.66 | 0.38 | 1.00 | |
| Azinphos Methyl | ND | 0.66 | 0.30 | 1.00 | |
| Bolstar | ND | 0.66 | 0.23 | 1.00 | |
| Chlorpyrifos | ND | 0.66 | 0.22 | 1.00 | |
| Coumaphos | ND | 0.66 | 0.20 | 1.00 | |
| Diazinon | ND | 0.66 | 0.21 | 1.00 | |
| Dichlorvos | ND | 0.66 | 0.36 | 1.00 | |
| Disulfoton | ND | 0.66 | 0.20 | 1.00 | |
| Ethoprop | ND | 0.66 | 0.22 | 1.00 | |
| Fensulfothion | ND | 0.66 | 0.24 | 1.00 | |
| Fenthion | ND | 0.66 | 0.24 | 1.00 | |
| Merphos | ND | 0.66 | 0.25 | 1.00 | |
| Methyl Parathion | ND | 0.66 | 0.28 | 1.00 | |
| Mevinphos | ND | 0.66 | 0.32 | 1.00 | |
| Naled | ND | 5.3 | 2.3 | 1.00 | |
| Phorate | ND | 0.66 | 0.30 | 1.00 | |
| Ronnel | ND | 0.66 | 0.20 | 1.00 | |
| Stirophos | ND | 2.6 | 0.79 | 1.00 | |
| Tokuthion | ND | 0.66 | 0.22 | 1.00 | |
| Trichloronate | ND | 0.66 | 0.20 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 110 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM370 | 15-06-0567-7-B | 06/05/15 13:51 | Solid | GC 26 | 06/09/15 | 06/15/15 15:06 | 150609L20 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|------|------|------|------------|
| Demeton-o/s | ND | 0.86 | 0.49 | 1.00 | |
| Azinphos Methyl | ND | 0.86 | 0.39 | 1.00 | |
| Bolstar | ND | 0.86 | 0.30 | 1.00 | |
| Chlorpyrifos | ND | 0.86 | 0.29 | 1.00 | |
| Coumaphos | ND | 0.86 | 0.26 | 1.00 | |
| Diazinon | ND | 0.86 | 0.27 | 1.00 | |
| Dichlorvos | ND | 0.86 | 0.47 | 1.00 | |
| Disulfoton | ND | 0.86 | 0.26 | 1.00 | |
| Ethoprop | ND | 0.86 | 0.29 | 1.00 | |
| Fensulfothion | ND | 0.86 | 0.31 | 1.00 | |
| Fenthion | ND | 0.86 | 0.32 | 1.00 | |
| Merphos | ND | 0.86 | 0.33 | 1.00 | |
| Methyl Parathion | ND | 0.86 | 0.37 | 1.00 | |
| Mevinphos | ND | 0.86 | 0.41 | 1.00 | |
| Naled | ND | 6.9 | 3.0 | 1.00 | |
| Phorate | ND | 0.86 | 0.39 | 1.00 | |
| Ronnel | ND | 0.86 | 0.27 | 1.00 | |
| Stirophos | ND | 3.4 | 1.0 | 1.00 | |
| Tokuthion | ND | 0.86 | 0.29 | 1.00 | |
| Trichloronate | ND | 0.86 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 101 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| FDHSM370 | 15-06-0567-8-B | 06/05/15 13:51 | Solid | GC 26 | 06/09/15 | 06/15/15 15:50 | 150609L20 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|------|------|------|------------|
| Demeton-o/s | ND | 0.90 | 0.52 | 1.00 | |
| Azinphos Methyl | ND | 0.90 | 0.41 | 1.00 | |
| Bolstar | ND | 0.90 | 0.32 | 1.00 | |
| Chlorpyrifos | ND | 0.90 | 0.30 | 1.00 | |
| Coumaphos | ND | 0.90 | 0.27 | 1.00 | |
| Diazinon | ND | 0.90 | 0.29 | 1.00 | |
| Dichlorvos | ND | 0.90 | 0.50 | 1.00 | |
| Disulfoton | ND | 0.90 | 0.28 | 1.00 | |
| Ethoprop | ND | 0.90 | 0.30 | 1.00 | |
| Fensulfothion | ND | 0.90 | 0.33 | 1.00 | |
| Fenthion | ND | 0.90 | 0.34 | 1.00 | |
| Merphos | ND | 0.90 | 0.35 | 1.00 | |
| Methyl Parathion | ND | 0.90 | 0.38 | 1.00 | |
| Mevinphos | ND | 0.90 | 0.44 | 1.00 | |
| Naled | ND | 7.2 | 3.1 | 1.00 | |
| Phorate | ND | 0.90 | 0.41 | 1.00 | |
| Ronnel | ND | 0.90 | 0.28 | 1.00 | |
| Stirophos | ND | 3.6 | 1.1 | 1.00 | |
| Tokuthion | ND | 0.90 | 0.30 | 1.00 | |
| Trichloronate | ND | 0.90 | 0.27 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 106 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8141A
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-973-188 | N/A | Solid | GC 26 | 06/09/15 | 06/12/15 11:24 | 150609L20 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|------|------|------|------------|
| Demeton-o/s | ND | 0.50 | 0.29 | 1.00 | |
| Azinphos Methyl | ND | 0.50 | 0.23 | 1.00 | |
| Bolstar | ND | 0.50 | 0.18 | 1.00 | |
| Chlorpyrifos | ND | 0.50 | 0.17 | 1.00 | |
| Coumaphos | ND | 0.50 | 0.15 | 1.00 | |
| Diazinon | ND | 0.50 | 0.16 | 1.00 | |
| Dichlorvos | ND | 0.50 | 0.28 | 1.00 | |
| Disulfoton | ND | 0.50 | 0.15 | 1.00 | |
| Ethoprop | ND | 0.50 | 0.17 | 1.00 | |
| Fensulfothion | ND | 0.50 | 0.18 | 1.00 | |
| Fenthion | ND | 0.50 | 0.19 | 1.00 | |
| Merphos | ND | 0.50 | 0.19 | 1.00 | |
| Methyl Parathion | ND | 0.50 | 0.21 | 1.00 | |
| Mevinphos | ND | 0.50 | 0.24 | 1.00 | |
| Naled | ND | 4.0 | 1.7 | 1.00 | |
| Phorate | ND | 0.50 | 0.22 | 1.00 | |
| Ronnel | ND | 0.50 | 0.16 | 1.00 | |
| Stirophos | ND | 2.0 | 0.60 | 1.00 | |
| Tokuthion | ND | 0.50 | 0.17 | 1.00 | |
| Trichloronate | ND | 0.50 | 0.15 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 97 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/kg

Project: EAA 27122

Page 1 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM310 | 15-06-0567-1-B | 06/05/15 10:03 | Solid | GC 40 | 06/10/15 | 06/17/15 13:35 | 150610L09 |

Comment(s): - Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|-------|-------|------|------------|
| Dalapon | ND | 870 | 250 | 1.00 | |
| Dicamba | ND | 35 | 4.3 | 1.00 | |
| MCP | ND | 35000 | 11000 | 1.00 | |
| MCPA | ND | 35000 | 7100 | 1.00 | |
| Dichlorprop | ND | 350 | 58 | 1.00 | |
| 2,4-D | ND | 350 | 49 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 35 | 8.1 | 1.00 | |
| 2,4,5-T | ND | 35 | 4.0 | 1.00 | |
| 2,4-DB | ND | 350 | 55 | 1.00 | |
| Dinoseb | ND | 170 | 18 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 69 | 30-130 | |

| HSM320 | 15-06-0567-2-B | 06/05/15 10:48 | Solid | GC 40 | 06/10/15 | 06/17/15 13:59 | 150610L09 |
|--------|----------------|-------------------|-------|-------|----------|-------------------|-----------|
|--------|----------------|-------------------|-------|-------|----------|-------------------|-----------|

Comment(s): - Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|-------|-------|------|------------|
| Dalapon | ND | 1300 | 370 | 1.00 | |
| Dicamba | ND | 51 | 6.3 | 1.00 | |
| MCP | ND | 51000 | 16000 | 1.00 | |
| MCPA | ND | 51000 | 10000 | 1.00 | |
| Dichlorprop | ND | 510 | 85 | 1.00 | |
| 2,4-D | ND | 510 | 73 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 51 | 12 | 1.00 | |
| 2,4,5-T | ND | 51 | 5.8 | 1.00 | |
| 2,4-DB | ND | 510 | 81 | 1.00 | |
| Dinoseb | ND | 250 | 26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 66 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM330 | 15-06-0567-3-B | 06/05/15 11:03 | Solid | GC 40 | 06/10/15 | 06/17/15 14:22 | 150610L09 |

Comment(s): - Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|-------|------|------|------------|
| Dalapon | ND | 310 | 92 | 1.00 | |
| Dicamba | ND | 13 | 1.6 | 1.00 | |
| MCP | ND | 13000 | 4000 | 1.00 | |
| MCPA | ND | 13000 | 2600 | 1.00 | |
| Dichlorprop | ND | 130 | 21 | 1.00 | |
| 2,4-D | ND | 130 | 18 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 13 | 3.0 | 1.00 | |
| 2,4,5-T | ND | 13 | 1.4 | 1.00 | |
| 2,4-DB | ND | 130 | 20 | 1.00 | |
| Dinoseb | ND | 63 | 6.4 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 65 | 30-130 | |

| | | | | | | | |
|--------|----------------|-------------------|-------|-------|----------|-------------------|-----------|
| HSM340 | 15-06-0567-4-B | 06/05/15 11:21 | Solid | GC 40 | 06/10/15 | 06/17/15 14:45 | 150610L09 |
|--------|----------------|-------------------|-------|-------|----------|-------------------|-----------|

Comment(s): - Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|-------|------|------|------------|
| Dalapon | ND | 310 | 90 | 1.00 | |
| Dicamba | ND | 12 | 1.5 | 1.00 | |
| MCP | ND | 12000 | 3900 | 1.00 | |
| MCPA | ND | 12000 | 2500 | 1.00 | |
| Dichlorprop | ND | 120 | 21 | 1.00 | |
| 2,4-D | ND | 120 | 18 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 12 | 2.9 | 1.00 | |
| 2,4,5-T | ND | 12 | 1.4 | 1.00 | |
| 2,4-DB | ND | 120 | 20 | 1.00 | |
| Dinoseb | ND | 62 | 6.3 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 70 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM350 | 15-06-0567-5-B | 06/05/15 12:24 | Solid | GC 40 | 06/10/15 | 06/17/15 15:08 | 150610L09 |

Comment(s): - Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|-------|------|------|------------|
| Dalapon | ND | 550 | 160 | 1.00 | |
| Dicamba | ND | 22 | 2.7 | 1.00 | |
| MCP | ND | 22000 | 6900 | 1.00 | |
| MCPA | ND | 22000 | 4500 | 1.00 | |
| Dichlorprop | ND | 220 | 36 | 1.00 | |
| 2,4-D | ND | 220 | 31 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 22 | 5.1 | 1.00 | |
| 2,4,5-T | ND | 22 | 2.5 | 1.00 | |
| 2,4-DB | ND | 220 | 35 | 1.00 | |
| Dinoseb | ND | 110 | 11 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 72 | 30-130 | |

| | | | | | | | |
|--------|----------------|-------------------|-------|-------|----------|-------------------|-----------|
| HSM360 | 15-06-0567-6-B | 06/05/15 13:13 | Solid | GC 40 | 06/10/15 | 06/17/15 15:31 | 150610L09 |
|--------|----------------|-------------------|-------|-------|----------|-------------------|-----------|

Comment(s): - Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|-------|------|------|------------|
| Dalapon | ND | 330 | 97 | 1.00 | |
| Dicamba | ND | 13 | 1.7 | 1.00 | |
| MCP | ND | 13000 | 4200 | 1.00 | |
| MCPA | ND | 13000 | 2700 | 1.00 | |
| Dichlorprop | ND | 130 | 22 | 1.00 | |
| 2,4-D | ND | 130 | 19 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 13 | 3.1 | 1.00 | |
| 2,4,5-T | ND | 13 | 1.5 | 1.00 | |
| 2,4-DB | ND | 130 | 21 | 1.00 | |
| Dinoseb | ND | 67 | 6.8 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 60 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM370 | 15-06-0567-7-B | 06/05/15 13:51 | Solid | GC 40 | 06/10/15 | 06/17/15 15:55 | 150610L09 |

Comment(s): - Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|-------|------|------|------------|
| Dalapon | ND | 430 | 120 | 1.00 | |
| Dicamba | ND | 17 | 2.1 | 1.00 | |
| MCP | ND | 17000 | 5400 | 1.00 | |
| MCPA | ND | 17000 | 3500 | 1.00 | |
| Dichlorprop | ND | 170 | 28 | 1.00 | |
| 2,4-D | ND | 170 | 24 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 17 | 4.0 | 1.00 | |
| 2,4,5-T | ND | 17 | 1.9 | 1.00 | |
| 2,4-DB | ND | 170 | 27 | 1.00 | |
| Dinoseb | ND | 85 | 8.6 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 84 | 30-130 | |

| FDHSM370 | 15-06-0567-8-B | 06/05/15 13:51 | Solid | GC 40 | 06/10/15 | 06/17/15 16:18 | 150610L09 |
|----------|----------------|----------------|-------|-------|----------|----------------|-----------|
|----------|----------------|----------------|-------|-------|----------|----------------|-----------|

Comment(s): - Results are reported on a dry weight basis.
- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|-------|------|------|------------|
| Dalapon | ND | 460 | 130 | 1.00 | |
| Dicamba | ND | 18 | 2.3 | 1.00 | |
| MCP | ND | 18000 | 5800 | 1.00 | |
| MCPA | ND | 18000 | 3700 | 1.00 | |
| Dichlorprop | ND | 180 | 30 | 1.00 | |
| 2,4-D | ND | 180 | 26 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 18 | 4.3 | 1.00 | |
| 2,4,5-T | ND | 18 | 2.1 | 1.00 | |
| 2,4-DB | ND | 180 | 29 | 1.00 | |
| Dinoseb | ND | 91 | 9.3 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 85 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-033-1294 | N/A | Solid | GC 40 | 06/10/15 | 06/11/15 19:20 | 150610L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|-------|------|------|------------|
| Dalapon | ND | 250 | 73 | 1.00 | |
| Dicamba | ND | 10 | 1.2 | 1.00 | |
| MCPD | ND | 10000 | 3200 | 1.00 | |
| MCPA | ND | 10000 | 2100 | 1.00 | |
| Dichlorprop | ND | 100 | 17 | 1.00 | |
| 2,4-D | ND | 100 | 14 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 10 | 2.3 | 1.00 | |
| 2,4,5-T | ND | 10 | 1.1 | 1.00 | |
| 2,4-DB | ND | 100 | 16 | 1.00 | |
| Dinoseb | ND | 50 | 5.1 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 67 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB01 | 15-06-0567-10-I | 06/05/15 15:41 | Aqueous | GC 40 | 06/08/15 | 06/16/15 20:19 | 150608L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 73 | 0-123 | |

| Method Blank | 095-01-034-647 | N/A | Aqueous | GC 40 | 06/08/15 | 06/12/15 04:34 | 150608L06 |
|--------------|----------------|-----|---------|-------|----------|-------------------|-----------|
|--------------|----------------|-----|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 118 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB01 | 15-06-0567-10-H | 06/05/15 15:41 | Aqueous | GC/MS TT | 06/08/15 | 06/08/15 22:48 | 150608L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 3 of 6

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 73 | 33-120 | |
| 2-Fluorophenol | 45 | 24-120 | |
| Nitrobenzene-d5 | 79 | 38-120 | |
| p-Terphenyl-d14 | 76 | 41-137 | |
| Phenol-d6 | 27 | 16-120 | |
| 2,4,6-Tribromophenol | 91 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 4 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-4051 | N/A | Aqueous | GC/MS TT | 06/08/15 | 06/08/15 16:52 | 150608L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 6 of 6

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 77 | 33-120 | |
| 2-Fluorophenol | 52 | 24-120 | |
| Nitrobenzene-d5 | 81 | 38-120 | |
| p-Terphenyl-d14 | 81 | 41-137 | |
| Phenol-d6 | 31 | 16-120 | |
| 2,4,6-Tribromophenol | 91 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

Page 1 of 27

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM310 | 15-06-0567-1-B | 06/05/15 10:03 | Solid | GC/MS CCC | 06/11/15 | 06/12/15 12:55 | 150611L02 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|------|------|------------|
| Acenaphthene | ND | 1.7 | 0.22 | 1.00 | |
| Acenaphthylene | ND | 1.7 | 0.21 | 1.00 | |
| Aniline | ND | 1.7 | 0.21 | 1.00 | |
| Anthracene | ND | 1.7 | 0.22 | 1.00 | |
| Azobenzene | ND | 1.7 | 0.19 | 1.00 | |
| Benzidine | ND | 34 | 2.8 | 1.00 | |
| Benzo (a) Anthracene | ND | 1.7 | 0.20 | 1.00 | |
| Benzo (a) Pyrene | ND | 1.7 | 0.19 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 1.7 | 0.22 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 1.7 | 0.19 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 1.7 | 0.22 | 1.00 | |
| Benzoic Acid | ND | 8.6 | 1.7 | 1.00 | |
| Benzyl Alcohol | ND | 1.7 | 0.23 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 1.7 | 0.19 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 8.6 | 1.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 1.7 | 0.20 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 1.7 | 0.18 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 1.7 | 0.21 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 1.7 | 0.19 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 1.7 | 0.23 | 1.00 | |
| 4-Chloroaniline | ND | 1.7 | 0.22 | 1.00 | |
| 2-Chloronaphthalene | ND | 1.7 | 0.20 | 1.00 | |
| 2-Chlorophenol | ND | 1.7 | 0.23 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 1.7 | 0.22 | 1.00 | |
| Chrysene | ND | 1.7 | 0.22 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 1.7 | 0.21 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 1.7 | 0.35 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 1.7 | 0.16 | 1.00 | |
| Dibenzofuran | ND | 1.7 | 0.21 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.7 | 0.22 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.7 | 0.25 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.7 | 0.26 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 34 | 1.3 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|------|------|------------|
| 2,4-Dichlorophenol | ND | 1.7 | 0.20 | 1.00 | |
| Diethyl Phthalate | ND | 1.7 | 0.20 | 1.00 | |
| Dimethyl Phthalate | ND | 1.7 | 0.39 | 1.00 | |
| 2,4-Dimethylphenol | ND | 1.7 | 0.91 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 8.6 | 2.2 | 1.00 | |
| 2,4-Dinitrophenol | ND | 8.6 | 1.2 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 1.7 | 0.23 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 1.7 | 0.25 | 1.00 | |
| Fluoranthene | ND | 1.7 | 0.21 | 1.00 | |
| Fluorene | ND | 1.7 | 0.22 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 1.7 | 0.22 | 1.00 | |
| Hexachlorobenzene | ND | 1.7 | 0.23 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 8.6 | 1.7 | 1.00 | |
| Hexachloroethane | ND | 1.7 | 0.27 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 1.7 | 0.18 | 1.00 | |
| Isophorone | ND | 1.7 | 0.20 | 1.00 | |
| 2-Methylnaphthalene | ND | 1.7 | 0.21 | 1.00 | |
| 1-Methylnaphthalene | ND | 1.7 | 0.19 | 1.00 | |
| 2-Methylphenol | ND | 1.7 | 0.30 | 1.00 | |
| 3/4-Methylphenol | ND | 1.7 | 0.56 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 1.7 | 0.29 | 1.00 | |
| N-Nitrosodimethylamine | ND | 1.7 | 0.16 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 1.7 | 0.47 | 1.00 | |
| Naphthalene | ND | 1.7 | 0.20 | 1.00 | |
| 4-Nitroaniline | ND | 1.7 | 0.22 | 1.00 | |
| 3-Nitroaniline | ND | 1.7 | 0.24 | 1.00 | |
| 2-Nitroaniline | ND | 1.7 | 0.18 | 1.00 | |
| Nitrobenzene | ND | 8.6 | 1.1 | 1.00 | |
| 4-Nitrophenol | ND | 1.7 | 0.19 | 1.00 | |
| 2-Nitrophenol | ND | 1.7 | 0.23 | 1.00 | |
| Pentachlorophenol | ND | 8.6 | 1.3 | 1.00 | |
| Phenanthrene | ND | 1.7 | 0.24 | 1.00 | |
| Phenol | ND | 1.7 | 0.17 | 1.00 | |
| Pyrene | ND | 1.7 | 0.26 | 1.00 | |
| Pyridine | ND | 1.7 | 0.19 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.7 | 0.20 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 1.7 | 0.26 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 1.7 | 0.22 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 56 | 27-120 | |
| 2-Fluorophenol | 62 | 25-120 | |
| Nitrobenzene-d5 | 55 | 33-123 | |
| p-Terphenyl-d14 | 75 | 27-159 | |
| Phenol-d6 | 63 | 26-122 | |
| 2,4,6-Tribromophenol | 82 | 18-138 | |



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

Page 4 of 27

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM320 | 15-06-0567-2-B | 06/05/15 10:48 | Solid | GC/MS CCC | 06/11/15 | 06/12/15 13:13 | 150611L02 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|------|------|------------|
| Acenaphthene | ND | 2.6 | 0.32 | 1.00 | |
| Acenaphthylene | ND | 2.6 | 0.30 | 1.00 | |
| Aniline | ND | 2.6 | 0.30 | 1.00 | |
| Anthracene | ND | 2.6 | 0.32 | 1.00 | |
| Azobenzene | ND | 2.6 | 0.28 | 1.00 | |
| Benzidine | ND | 51 | 4.2 | 1.00 | |
| Benzo (a) Anthracene | 0.51 | 2.6 | 0.29 | 1.00 | J |
| Benzo (a) Pyrene | 0.40 | 2.6 | 0.28 | 1.00 | J |
| Benzo (b) Fluoranthene | ND | 2.6 | 0.33 | 1.00 | |
| Benzo (g,h,i) Perylene | 0.33 | 2.6 | 0.28 | 1.00 | J |
| Benzo (k) Fluoranthene | ND | 2.6 | 0.33 | 1.00 | |
| Benzoic Acid | ND | 13 | 2.5 | 1.00 | |
| Benzyl Alcohol | ND | 2.6 | 0.34 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 2.6 | 0.29 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 13 | 2.1 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 2.6 | 0.29 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 2.6 | 0.27 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 2.6 | 0.31 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 2.6 | 0.28 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 2.6 | 0.34 | 1.00 | |
| 4-Chloroaniline | ND | 2.6 | 0.33 | 1.00 | |
| 2-Chloronaphthalene | ND | 2.6 | 0.30 | 1.00 | |
| 2-Chlorophenol | ND | 2.6 | 0.34 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 2.6 | 0.33 | 1.00 | |
| Chrysene | 0.59 | 2.6 | 0.33 | 1.00 | J |
| Di-n-Butyl Phthalate | ND | 2.6 | 0.31 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 2.6 | 0.51 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 2.6 | 0.24 | 1.00 | |
| Dibenzofuran | ND | 2.6 | 0.31 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 2.6 | 0.33 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 2.6 | 0.37 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 2.6 | 0.38 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 51 | 1.9 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|------|------|------------|
| 2,4-Dichlorophenol | ND | 2.6 | 0.30 | 1.00 | |
| Diethyl Phthalate | ND | 2.6 | 0.30 | 1.00 | |
| Dimethyl Phthalate | ND | 2.6 | 0.58 | 1.00 | |
| 2,4-Dimethylphenol | ND | 2.6 | 1.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 13 | 3.2 | 1.00 | |
| 2,4-Dinitrophenol | ND | 13 | 1.7 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 2.6 | 0.34 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 2.6 | 0.37 | 1.00 | |
| Fluoranthene | 0.84 | 2.6 | 0.32 | 1.00 | J |
| Fluorene | ND | 2.6 | 0.32 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 2.6 | 0.32 | 1.00 | |
| Hexachlorobenzene | ND | 2.6 | 0.34 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 13 | 2.6 | 1.00 | |
| Hexachloroethane | ND | 2.6 | 0.40 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | 0.30 | 2.6 | 0.27 | 1.00 | J |
| Isophorone | ND | 2.6 | 0.29 | 1.00 | |
| 2-Methylnaphthalene | ND | 2.6 | 0.31 | 1.00 | |
| 1-Methylnaphthalene | ND | 2.6 | 0.29 | 1.00 | |
| 2-Methylphenol | ND | 2.6 | 0.44 | 1.00 | |
| 3/4-Methylphenol | ND | 2.6 | 0.83 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 2.6 | 0.43 | 1.00 | |
| N-Nitrosodimethylamine | ND | 2.6 | 0.24 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 2.6 | 0.70 | 1.00 | |
| Naphthalene | ND | 2.6 | 0.30 | 1.00 | |
| 4-Nitroaniline | ND | 2.6 | 0.33 | 1.00 | |
| 3-Nitroaniline | ND | 2.6 | 0.36 | 1.00 | |
| 2-Nitroaniline | ND | 2.6 | 0.26 | 1.00 | |
| Nitrobenzene | ND | 13 | 1.7 | 1.00 | |
| 4-Nitrophenol | ND | 2.6 | 0.28 | 1.00 | |
| 2-Nitrophenol | ND | 2.6 | 0.34 | 1.00 | |
| Pentachlorophenol | ND | 13 | 2.0 | 1.00 | |
| Phenanthrene | ND | 2.6 | 0.35 | 1.00 | |
| Phenol | ND | 2.6 | 0.24 | 1.00 | |
| Pyrene | 0.79 | 2.6 | 0.39 | 1.00 | J |
| Pyridine | ND | 2.6 | 0.28 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 2.6 | 0.30 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 2.6 | 0.39 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 2.6 | 0.32 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 55 | 27-120 | |
| 2-Fluorophenol | 67 | 25-120 | |
| Nitrobenzene-d5 | 56 | 33-123 | |
| p-Terphenyl-d14 | 79 | 27-159 | |
| Phenol-d6 | 66 | 26-122 | |
| 2,4,6-Tribromophenol | 84 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM330 | 15-06-0567-3-B | 06/05/15 11:03 | Solid | GC/MS CCC | 06/11/15 | 06/12/15 13:31 | 150611L02 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|------|-------|------|------------|
| Acenaphthene | ND | 0.62 | 0.079 | 1.00 | |
| Acenaphthylene | ND | 0.62 | 0.074 | 1.00 | |
| Aniline | ND | 0.62 | 0.074 | 1.00 | |
| Anthracene | ND | 0.62 | 0.079 | 1.00 | |
| Azobenzene | ND | 0.62 | 0.067 | 1.00 | |
| Benzidine | ND | 12 | 1.0 | 1.00 | |
| Benzo (a) Anthracene | 0.65 | 0.62 | 0.071 | 1.00 | |
| Benzo (a) Pyrene | 0.57 | 0.62 | 0.068 | 1.00 | J |
| Benzo (b) Fluoranthene | 0.72 | 0.62 | 0.079 | 1.00 | |
| Benzo (g,h,i) Perylene | 0.49 | 0.62 | 0.069 | 1.00 | J |
| Benzo (k) Fluoranthene | 0.61 | 0.62 | 0.081 | 1.00 | J |
| Benzoic Acid | ND | 3.1 | 0.62 | 1.00 | |
| Benzyl Alcohol | ND | 0.62 | 0.082 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 0.62 | 0.070 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 3.1 | 0.51 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 0.62 | 0.072 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | 0.12 | 0.62 | 0.066 | 1.00 | J |
| 4-Bromophenyl-Phenyl Ether | ND | 0.62 | 0.076 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 0.62 | 0.067 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 0.62 | 0.083 | 1.00 | |
| 4-Chloroaniline | ND | 0.62 | 0.081 | 1.00 | |
| 2-Chloronaphthalene | ND | 0.62 | 0.074 | 1.00 | |
| 2-Chlorophenol | ND | 0.62 | 0.083 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 0.62 | 0.080 | 1.00 | |
| Chrysene | 0.88 | 0.62 | 0.080 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 0.62 | 0.075 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 0.62 | 0.13 | 1.00 | |
| Dibenz (a,h) Anthracene | 0.15 | 0.62 | 0.058 | 1.00 | J |
| Dibenzofuran | ND | 0.62 | 0.075 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 0.62 | 0.080 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 0.62 | 0.091 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 0.62 | 0.094 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 12 | 0.45 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|------|-------|------|------------|
| 2,4-Dichlorophenol | ND | 0.62 | 0.072 | 1.00 | |
| Diethyl Phthalate | ND | 0.62 | 0.072 | 1.00 | |
| Dimethyl Phthalate | ND | 0.62 | 0.14 | 1.00 | |
| 2,4-Dimethylphenol | ND | 0.62 | 0.33 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 3.1 | 0.79 | 1.00 | |
| 2,4-Dinitrophenol | ND | 3.1 | 0.42 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 0.62 | 0.083 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 0.62 | 0.091 | 1.00 | |
| Fluoranthene | 1.5 | 0.62 | 0.078 | 1.00 | |
| Fluorene | ND | 0.62 | 0.079 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 0.62 | 0.079 | 1.00 | |
| Hexachlorobenzene | ND | 0.62 | 0.083 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 3.1 | 0.63 | 1.00 | |
| Hexachloroethane | ND | 0.62 | 0.098 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | 0.43 | 0.62 | 0.067 | 1.00 | J |
| Isophorone | ND | 0.62 | 0.071 | 1.00 | |
| 2-Methylnaphthalene | ND | 0.62 | 0.075 | 1.00 | |
| 1-Methylnaphthalene | ND | 0.62 | 0.070 | 1.00 | |
| 2-Methylphenol | ND | 0.62 | 0.11 | 1.00 | |
| 3/4-Methylphenol | ND | 0.62 | 0.20 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 0.62 | 0.10 | 1.00 | |
| N-Nitrosodimethylamine | ND | 0.62 | 0.059 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 0.62 | 0.17 | 1.00 | |
| Naphthalene | ND | 0.62 | 0.073 | 1.00 | |
| 4-Nitroaniline | ND | 0.62 | 0.081 | 1.00 | |
| 3-Nitroaniline | ND | 0.62 | 0.087 | 1.00 | |
| 2-Nitroaniline | ND | 0.62 | 0.065 | 1.00 | |
| Nitrobenzene | ND | 3.1 | 0.40 | 1.00 | |
| 4-Nitrophenol | ND | 0.62 | 0.068 | 1.00 | |
| 2-Nitrophenol | ND | 0.62 | 0.084 | 1.00 | |
| Pentachlorophenol | ND | 3.1 | 0.48 | 1.00 | |
| Phenanthrene | 0.59 | 0.62 | 0.085 | 1.00 | J |
| Phenol | ND | 0.62 | 0.060 | 1.00 | |
| Pyrene | 1.4 | 0.62 | 0.095 | 1.00 | |
| Pyridine | ND | 0.62 | 0.068 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 0.62 | 0.074 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 0.62 | 0.096 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 0.62 | 0.079 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 73 | 27-120 | |
| 2-Fluorophenol | 74 | 25-120 | |
| Nitrobenzene-d5 | 64 | 33-123 | |
| p-Terphenyl-d14 | 95 | 27-159 | |
| Phenol-d6 | 76 | 26-122 | |
| 2,4,6-Tribromophenol | 98 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM340 | 15-06-0567-4-B | 06/05/15 11:21 | Solid | GC/MS CCC | 06/11/15 | 06/12/15 13:49 | 150611L02 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|------|------|------------|
| Acenaphthene | 0.27 | 1.2 | 0.16 | 2.00 | J |
| Acenaphthylene | ND | 1.2 | 0.15 | 2.00 | |
| Aniline | ND | 1.2 | 0.15 | 2.00 | |
| Anthracene | 0.68 | 1.2 | 0.16 | 2.00 | J |
| Azobenzene | ND | 1.2 | 0.13 | 2.00 | |
| Benzidine | ND | 25 | 2.0 | 2.00 | |
| Benzo (a) Anthracene | 4.7 | 1.2 | 0.14 | 2.00 | |
| Benzo (a) Pyrene | 3.8 | 1.2 | 0.14 | 2.00 | |
| Benzo (b) Fluoranthene | 4.8 | 1.2 | 0.16 | 2.00 | |
| Benzo (g,h,i) Perylene | 2.9 | 1.2 | 0.14 | 2.00 | |
| Benzo (k) Fluoranthene | 3.8 | 1.2 | 0.16 | 2.00 | |
| Benzoic Acid | ND | 6.2 | 1.2 | 2.00 | |
| Benzyl Alcohol | ND | 1.2 | 0.16 | 2.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 1.2 | 0.14 | 2.00 | |
| Bis(2-Chloroethyl) Ether | ND | 6.2 | 1.0 | 2.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 1.2 | 0.14 | 2.00 | |
| Bis(2-Ethylhexyl) Phthalate | 0.74 | 1.2 | 0.13 | 2.00 | J |
| 4-Bromophenyl-Phenyl Ether | ND | 1.2 | 0.15 | 2.00 | |
| Butyl Benzyl Phthalate | ND | 1.2 | 0.13 | 2.00 | |
| 4-Chloro-3-Methylphenol | ND | 1.2 | 0.16 | 2.00 | |
| 4-Chloroaniline | ND | 1.2 | 0.16 | 2.00 | |
| 2-Chloronaphthalene | ND | 1.2 | 0.15 | 2.00 | |
| 2-Chlorophenol | ND | 1.2 | 0.17 | 2.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 1.2 | 0.16 | 2.00 | |
| Chrysene | 6.2 | 1.2 | 0.16 | 2.00 | |
| Di-n-Butyl Phthalate | ND | 1.2 | 0.15 | 2.00 | |
| Di-n-Octyl Phthalate | ND | 1.2 | 0.25 | 2.00 | |
| Dibenz (a,h) Anthracene | 0.88 | 1.2 | 0.12 | 2.00 | J |
| Dibenzofuran | ND | 1.2 | 0.15 | 2.00 | |
| 1,2-Dichlorobenzene | ND | 1.2 | 0.16 | 2.00 | |
| 1,3-Dichlorobenzene | ND | 1.2 | 0.18 | 2.00 | |
| 1,4-Dichlorobenzene | ND | 1.2 | 0.19 | 2.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 0.90 | 2.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|------|------|------------|
| 2,4-Dichlorophenol | ND | 1.2 | 0.14 | 2.00 | |
| Diethyl Phthalate | ND | 1.2 | 0.14 | 2.00 | |
| Dimethyl Phthalate | ND | 1.2 | 0.28 | 2.00 | |
| 2,4-Dimethylphenol | ND | 1.2 | 0.66 | 2.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 6.2 | 1.6 | 2.00 | |
| 2,4-Dinitrophenol | ND | 6.2 | 0.84 | 2.00 | |
| 2,4-Dinitrotoluene | ND | 1.2 | 0.17 | 2.00 | |
| 2,6-Dinitrotoluene | ND | 1.2 | 0.18 | 2.00 | |
| Fluoranthene | 13 | 1.2 | 0.15 | 2.00 | |
| Fluorene | 0.31 | 1.2 | 0.16 | 2.00 | J |
| Hexachloro-1,3-Butadiene | ND | 1.2 | 0.16 | 2.00 | |
| Hexachlorobenzene | ND | 1.2 | 0.17 | 2.00 | |
| Hexachlorocyclopentadiene | ND | 6.2 | 1.3 | 2.00 | |
| Hexachloroethane | ND | 1.2 | 0.19 | 2.00 | |
| Indeno (1,2,3-c,d) Pyrene | 2.7 | 1.2 | 0.13 | 2.00 | |
| Isophorone | ND | 1.2 | 0.14 | 2.00 | |
| 2-Methylnaphthalene | ND | 1.2 | 0.15 | 2.00 | |
| 1-Methylnaphthalene | ND | 1.2 | 0.14 | 2.00 | |
| 2-Methylphenol | ND | 1.2 | 0.21 | 2.00 | |
| 3/4-Methylphenol | ND | 1.2 | 0.40 | 2.00 | |
| N-Nitroso-di-n-propylamine | ND | 1.2 | 0.21 | 2.00 | |
| N-Nitrosodimethylamine | ND | 1.2 | 0.12 | 2.00 | |
| N-Nitrosodiphenylamine | ND | 1.2 | 0.34 | 2.00 | |
| Naphthalene | ND | 1.2 | 0.15 | 2.00 | |
| 4-Nitroaniline | ND | 1.2 | 0.16 | 2.00 | |
| 3-Nitroaniline | ND | 1.2 | 0.17 | 2.00 | |
| 2-Nitroaniline | ND | 1.2 | 0.13 | 2.00 | |
| Nitrobenzene | ND | 6.2 | 0.80 | 2.00 | |
| 4-Nitrophenol | ND | 1.2 | 0.13 | 2.00 | |
| 2-Nitrophenol | ND | 1.2 | 0.17 | 2.00 | |
| Pentachlorophenol | ND | 6.2 | 0.96 | 2.00 | |
| Phenanthrene | 6.6 | 1.2 | 0.17 | 2.00 | |
| Phenol | ND | 1.2 | 0.12 | 2.00 | |
| Pyrene | 12 | 1.2 | 0.19 | 2.00 | |
| Pyridine | ND | 1.2 | 0.14 | 2.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.2 | 0.15 | 2.00 | |
| 2,4,6-Trichlorophenol | ND | 1.2 | 0.19 | 2.00 | |
| 2,4,5-Trichlorophenol | ND | 1.2 | 0.16 | 2.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 75 | 27-120 | |
| 2-Fluorophenol | 62 | 25-120 | |
| Nitrobenzene-d5 | 64 | 33-123 | |
| p-Terphenyl-d14 | 92 | 27-159 | |
| Phenol-d6 | 66 | 26-122 | |
| 2,4,6-Tribromophenol | 82 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM350 | 15-06-0567-5-B | 06/05/15 12:24 | Solid | GC/MS CCC | 06/11/15 | 06/12/15 14:07 | 150611L02 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|------|------|------------|
| Acenaphthene | ND | 1.1 | 0.14 | 1.00 | |
| Acenaphthylene | ND | 1.1 | 0.13 | 1.00 | |
| Aniline | ND | 1.1 | 0.13 | 1.00 | |
| Anthracene | ND | 1.1 | 0.14 | 1.00 | |
| Azobenzene | ND | 1.1 | 0.12 | 1.00 | |
| Benzidine | ND | 22 | 1.8 | 1.00 | |
| Benzo (a) Anthracene | 0.15 | 1.1 | 0.12 | 1.00 | J |
| Benzo (a) Pyrene | 0.15 | 1.1 | 0.12 | 1.00 | J |
| Benzo (b) Fluoranthene | 0.21 | 1.1 | 0.14 | 1.00 | J |
| Benzo (g,h,i) Perylene | 0.13 | 1.1 | 0.12 | 1.00 | J |
| Benzo (k) Fluoranthene | 0.15 | 1.1 | 0.14 | 1.00 | J |
| Benzoic Acid | ND | 5.4 | 1.1 | 1.00 | |
| Benzyl Alcohol | ND | 1.1 | 0.14 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 1.1 | 0.12 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 5.4 | 0.88 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 1.1 | 0.12 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | 0.24 | 1.1 | 0.12 | 1.00 | J |
| 4-Bromophenyl-Phenyl Ether | ND | 1.1 | 0.13 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 1.1 | 0.12 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 1.1 | 0.14 | 1.00 | |
| 4-Chloroaniline | ND | 1.1 | 0.14 | 1.00 | |
| 2-Chloronaphthalene | ND | 1.1 | 0.13 | 1.00 | |
| 2-Chlorophenol | ND | 1.1 | 0.14 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 1.1 | 0.14 | 1.00 | |
| Chrysene | 0.21 | 1.1 | 0.14 | 1.00 | J |
| Di-n-Butyl Phthalate | ND | 1.1 | 0.13 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 1.1 | 0.22 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 1.1 | 0.10 | 1.00 | |
| Dibenzofuran | ND | 1.1 | 0.13 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.1 | 0.14 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.1 | 0.16 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.1 | 0.16 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 22 | 0.79 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|------|------|------------|
| 2,4-Dichlorophenol | ND | 1.1 | 0.13 | 1.00 | |
| Diethyl Phthalate | ND | 1.1 | 0.13 | 1.00 | |
| Dimethyl Phthalate | ND | 1.1 | 0.25 | 1.00 | |
| 2,4-Dimethylphenol | ND | 1.1 | 0.57 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 5.4 | 1.4 | 1.00 | |
| 2,4-Dinitrophenol | ND | 5.4 | 0.73 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 1.1 | 0.14 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 1.1 | 0.16 | 1.00 | |
| Fluoranthene | 0.30 | 1.1 | 0.13 | 1.00 | J |
| Fluorene | ND | 1.1 | 0.14 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 1.1 | 0.14 | 1.00 | |
| Hexachlorobenzene | ND | 1.1 | 0.14 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 5.4 | 1.1 | 1.00 | |
| Hexachloroethane | ND | 1.1 | 0.17 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | 0.12 | 1.1 | 0.12 | 1.00 | J |
| Isophorone | ND | 1.1 | 0.12 | 1.00 | |
| 2-Methylnaphthalene | ND | 1.1 | 0.13 | 1.00 | |
| 1-Methylnaphthalene | ND | 1.1 | 0.12 | 1.00 | |
| 2-Methylphenol | ND | 1.1 | 0.19 | 1.00 | |
| 3/4-Methylphenol | ND | 1.1 | 0.35 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 1.1 | 0.18 | 1.00 | |
| N-Nitrosodimethylamine | ND | 1.1 | 0.10 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 1.1 | 0.30 | 1.00 | |
| Naphthalene | ND | 1.1 | 0.13 | 1.00 | |
| 4-Nitroaniline | ND | 1.1 | 0.14 | 1.00 | |
| 3-Nitroaniline | ND | 1.1 | 0.15 | 1.00 | |
| 2-Nitroaniline | ND | 1.1 | 0.11 | 1.00 | |
| Nitrobenzene | ND | 5.4 | 0.70 | 1.00 | |
| 4-Nitrophenol | ND | 1.1 | 0.12 | 1.00 | |
| 2-Nitrophenol | ND | 1.1 | 0.15 | 1.00 | |
| Pentachlorophenol | ND | 5.4 | 0.84 | 1.00 | |
| Phenanthrene | ND | 1.1 | 0.15 | 1.00 | |
| Phenol | ND | 1.1 | 0.10 | 1.00 | |
| Pyrene | 0.66 | 1.1 | 0.17 | 1.00 | J |
| Pyridine | ND | 1.1 | 0.12 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.1 | 0.13 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 1.1 | 0.17 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 1.1 | 0.14 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 54 | 27-120 | |
| 2-Fluorophenol | 50 | 25-120 | |
| Nitrobenzene-d5 | 42 | 33-123 | |
| p-Terphenyl-d14 | 67 | 27-159 | |
| Phenol-d6 | 51 | 26-122 | |
| 2,4,6-Tribromophenol | 71 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM360 | 15-06-0567-6-B | 06/05/15 13:13 | Solid | GC/MS CCC | 06/11/15 | 06/12/15 14:25 | 150611L02 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|------|-------|------|------------|
| Acenaphthene | ND | 0.66 | 0.084 | 1.00 | |
| Acenaphthylene | ND | 0.66 | 0.079 | 1.00 | |
| Aniline | ND | 0.66 | 0.079 | 1.00 | |
| Anthracene | ND | 0.66 | 0.084 | 1.00 | |
| Azobenzene | ND | 0.66 | 0.072 | 1.00 | |
| Benzidine | ND | 13 | 1.1 | 1.00 | |
| Benzo (a) Anthracene | ND | 0.66 | 0.076 | 1.00 | |
| Benzo (a) Pyrene | ND | 0.66 | 0.072 | 1.00 | |
| Benzo (b) Fluoranthene | 0.090 | 0.66 | 0.085 | 1.00 | J |
| Benzo (g,h,i) Perylene | ND | 0.66 | 0.073 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 0.66 | 0.086 | 1.00 | |
| Benzoic Acid | ND | 3.3 | 0.66 | 1.00 | |
| Benzyl Alcohol | ND | 0.66 | 0.088 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 0.66 | 0.075 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 3.3 | 0.54 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 0.66 | 0.076 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 0.66 | 0.071 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 0.66 | 0.081 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 0.66 | 0.072 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 0.66 | 0.088 | 1.00 | |
| 4-Chloroaniline | ND | 0.66 | 0.087 | 1.00 | |
| 2-Chloronaphthalene | ND | 0.66 | 0.079 | 1.00 | |
| 2-Chlorophenol | ND | 0.66 | 0.089 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 0.66 | 0.085 | 1.00 | |
| Chrysene | 0.087 | 0.66 | 0.085 | 1.00 | J |
| Di-n-Butyl Phthalate | ND | 0.66 | 0.080 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 0.66 | 0.13 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 0.66 | 0.062 | 1.00 | |
| Dibenzofuran | ND | 0.66 | 0.080 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 0.66 | 0.085 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 0.66 | 0.097 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 0.66 | 0.10 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 13 | 0.48 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|------|-------|------|------------|
| 2,4-Dichlorophenol | ND | 0.66 | 0.077 | 1.00 | |
| Diethyl Phthalate | ND | 0.66 | 0.077 | 1.00 | |
| Dimethyl Phthalate | ND | 0.66 | 0.15 | 1.00 | |
| 2,4-Dimethylphenol | ND | 0.66 | 0.35 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 3.3 | 0.84 | 1.00 | |
| 2,4-Dinitrophenol | ND | 3.3 | 0.45 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 0.66 | 0.089 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 0.66 | 0.097 | 1.00 | |
| Fluoranthene | 0.11 | 0.66 | 0.083 | 1.00 | J |
| Fluorene | ND | 0.66 | 0.084 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 0.66 | 0.084 | 1.00 | |
| Hexachlorobenzene | ND | 0.66 | 0.089 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 3.3 | 0.67 | 1.00 | |
| Hexachloroethane | ND | 0.66 | 0.10 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 0.66 | 0.071 | 1.00 | |
| Isophorone | ND | 0.66 | 0.076 | 1.00 | |
| 2-Methylnaphthalene | ND | 0.66 | 0.080 | 1.00 | |
| 1-Methylnaphthalene | ND | 0.66 | 0.075 | 1.00 | |
| 2-Methylphenol | ND | 0.66 | 0.12 | 1.00 | |
| 3/4-Methylphenol | ND | 0.66 | 0.22 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 0.66 | 0.11 | 1.00 | |
| N-Nitrosodimethylamine | ND | 0.66 | 0.063 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 0.66 | 0.18 | 1.00 | |
| Naphthalene | ND | 0.66 | 0.078 | 1.00 | |
| 4-Nitroaniline | ND | 0.66 | 0.087 | 1.00 | |
| 3-Nitroaniline | ND | 0.66 | 0.093 | 1.00 | |
| 2-Nitroaniline | ND | 0.66 | 0.069 | 1.00 | |
| Nitrobenzene | ND | 3.3 | 0.43 | 1.00 | |
| 4-Nitrophenol | ND | 0.66 | 0.072 | 1.00 | |
| 2-Nitrophenol | ND | 0.66 | 0.090 | 1.00 | |
| Pentachlorophenol | ND | 3.3 | 0.51 | 1.00 | |
| Phenanthrene | ND | 0.66 | 0.091 | 1.00 | |
| Phenol | ND | 0.66 | 0.064 | 1.00 | |
| Pyrene | 0.11 | 0.66 | 0.10 | 1.00 | J |
| Pyridine | ND | 0.66 | 0.073 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 0.66 | 0.079 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 0.66 | 0.10 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 0.66 | 0.085 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 82 | 27-120 | |
| 2-Fluorophenol | 77 | 25-120 | |
| Nitrobenzene-d5 | 78 | 33-123 | |
| p-Terphenyl-d14 | 104 | 27-159 | |
| Phenol-d6 | 77 | 26-122 | |
| 2,4,6-Tribromophenol | 98 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM370 | 15-06-0567-7-B | 06/05/15 13:51 | Solid | GC/MS CCC | 06/11/15 | 06/12/15 15:01 | 150611L02 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|------|-------|------|------------|
| Acenaphthene | ND | 0.85 | 0.11 | 1.00 | |
| Acenaphthylene | ND | 0.85 | 0.10 | 1.00 | |
| Aniline | ND | 0.85 | 0.10 | 1.00 | |
| Anthracene | ND | 0.85 | 0.11 | 1.00 | |
| Azobenzene | ND | 0.85 | 0.092 | 1.00 | |
| Benzidine | ND | 17 | 1.4 | 1.00 | |
| Benzo (a) Anthracene | ND | 0.85 | 0.097 | 1.00 | |
| Benzo (a) Pyrene | ND | 0.85 | 0.093 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 0.85 | 0.11 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 0.85 | 0.094 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 0.85 | 0.11 | 1.00 | |
| Benzoic Acid | ND | 4.3 | 0.84 | 1.00 | |
| Benzyl Alcohol | ND | 0.85 | 0.11 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 0.85 | 0.096 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 4.3 | 0.69 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 0.85 | 0.098 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 0.85 | 0.091 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 0.85 | 0.10 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 0.85 | 0.092 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 0.85 | 0.11 | 1.00 | |
| 4-Chloroaniline | ND | 0.85 | 0.11 | 1.00 | |
| 2-Chloronaphthalene | ND | 0.85 | 0.10 | 1.00 | |
| 2-Chlorophenol | ND | 0.85 | 0.11 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 0.85 | 0.11 | 1.00 | |
| Chrysene | ND | 0.85 | 0.11 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 0.85 | 0.10 | 1.00 | |
| Di-n-Octyl Phthalate | 0.48 | 0.85 | 0.17 | 1.00 | J |
| Dibenz (a,h) Anthracene | ND | 0.85 | 0.079 | 1.00 | |
| Dibenzofuran | ND | 0.85 | 0.10 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 0.85 | 0.11 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 0.85 | 0.12 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 0.85 | 0.13 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 17 | 0.62 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|------|-------|------|------------|
| 2,4-Dichlorophenol | ND | 0.85 | 0.099 | 1.00 | |
| Diethyl Phthalate | ND | 0.85 | 0.099 | 1.00 | |
| Dimethyl Phthalate | ND | 0.85 | 0.19 | 1.00 | |
| 2,4-Dimethylphenol | ND | 0.85 | 0.45 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 4.3 | 1.1 | 1.00 | |
| 2,4-Dinitrophenol | ND | 4.3 | 0.58 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 0.85 | 0.11 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 0.85 | 0.12 | 1.00 | |
| Fluoranthene | 0.11 | 0.85 | 0.11 | 1.00 | J |
| Fluorene | ND | 0.85 | 0.11 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 0.85 | 0.11 | 1.00 | |
| Hexachlorobenzene | ND | 0.85 | 0.11 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 4.3 | 0.86 | 1.00 | |
| Hexachloroethane | ND | 0.85 | 0.13 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 0.85 | 0.091 | 1.00 | |
| Isophorone | ND | 0.85 | 0.097 | 1.00 | |
| 2-Methylnaphthalene | 0.15 | 0.85 | 0.10 | 1.00 | J |
| 1-Methylnaphthalene | ND | 0.85 | 0.096 | 1.00 | |
| 2-Methylphenol | ND | 0.85 | 0.15 | 1.00 | |
| 3/4-Methylphenol | ND | 0.85 | 0.28 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 0.85 | 0.14 | 1.00 | |
| N-Nitrosodimethylamine | ND | 0.85 | 0.080 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 0.85 | 0.23 | 1.00 | |
| Naphthalene | ND | 0.85 | 0.10 | 1.00 | |
| 4-Nitroaniline | ND | 0.85 | 0.11 | 1.00 | |
| 3-Nitroaniline | ND | 0.85 | 0.12 | 1.00 | |
| 2-Nitroaniline | ND | 0.85 | 0.088 | 1.00 | |
| Nitrobenzene | ND | 4.3 | 0.55 | 1.00 | |
| 4-Nitrophenol | ND | 0.85 | 0.092 | 1.00 | |
| 2-Nitrophenol | ND | 0.85 | 0.11 | 1.00 | |
| Pentachlorophenol | ND | 4.3 | 0.66 | 1.00 | |
| Phenanthrene | ND | 0.85 | 0.12 | 1.00 | |
| Phenol | ND | 0.85 | 0.081 | 1.00 | |
| Pyrene | ND | 0.85 | 0.13 | 1.00 | |
| Pyridine | ND | 0.85 | 0.093 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 0.85 | 0.10 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 0.85 | 0.13 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 0.85 | 0.11 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 65 | 27-120 | |
| 2-Fluorophenol | 58 | 25-120 | |
| Nitrobenzene-d5 | 54 | 33-123 | |
| p-Terphenyl-d14 | 79 | 27-159 | |
| Phenol-d6 | 54 | 26-122 | |
| 2,4,6-Tribromophenol | 96 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| FDHSM370 | 15-06-0567-8-B | 06/05/15 13:51 | Solid | GC/MS CCC | 06/11/15 | 06/12/15 15:19 | 150611L02 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|------|-------|------|------------|
| Acenaphthene | ND | 0.92 | 0.12 | 1.00 | |
| Acenaphthylene | ND | 0.92 | 0.11 | 1.00 | |
| Aniline | ND | 0.92 | 0.11 | 1.00 | |
| Anthracene | ND | 0.92 | 0.12 | 1.00 | |
| Azobenzene | ND | 0.92 | 0.10 | 1.00 | |
| Benzidine | ND | 18 | 1.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 0.92 | 0.10 | 1.00 | |
| Benzo (a) Pyrene | ND | 0.92 | 0.10 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 0.92 | 0.12 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 0.92 | 0.10 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 0.92 | 0.12 | 1.00 | |
| Benzoic Acid | ND | 4.6 | 0.91 | 1.00 | |
| Benzyl Alcohol | ND | 0.92 | 0.12 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 0.92 | 0.10 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 4.6 | 0.75 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 0.92 | 0.11 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 0.92 | 0.098 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 0.92 | 0.11 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 0.92 | 0.099 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 0.92 | 0.12 | 1.00 | |
| 4-Chloroaniline | ND | 0.92 | 0.12 | 1.00 | |
| 2-Chloronaphthalene | ND | 0.92 | 0.11 | 1.00 | |
| 2-Chlorophenol | ND | 0.92 | 0.12 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 0.92 | 0.12 | 1.00 | |
| Chrysene | ND | 0.92 | 0.12 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 0.92 | 0.11 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 0.92 | 0.19 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 0.92 | 0.085 | 1.00 | |
| Dibenzofuran | ND | 0.92 | 0.11 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 0.92 | 0.12 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 0.92 | 0.14 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 0.92 | 0.14 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 18 | 0.67 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|------|-------|------|------------|
| 2,4-Dichlorophenol | ND | 0.92 | 0.11 | 1.00 | |
| Diethyl Phthalate | ND | 0.92 | 0.11 | 1.00 | |
| Dimethyl Phthalate | ND | 0.92 | 0.21 | 1.00 | |
| 2,4-Dimethylphenol | ND | 0.92 | 0.49 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 4.6 | 1.2 | 1.00 | |
| 2,4-Dinitrophenol | ND | 4.6 | 0.62 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 0.92 | 0.12 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 0.92 | 0.13 | 1.00 | |
| Fluoranthene | ND | 0.92 | 0.11 | 1.00 | |
| Fluorene | ND | 0.92 | 0.12 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 0.92 | 0.12 | 1.00 | |
| Hexachlorobenzene | ND | 0.92 | 0.12 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 4.6 | 0.93 | 1.00 | |
| Hexachloroethane | ND | 0.92 | 0.14 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 0.92 | 0.098 | 1.00 | |
| Isophorone | ND | 0.92 | 0.10 | 1.00 | |
| 2-Methylnaphthalene | ND | 0.92 | 0.11 | 1.00 | |
| 1-Methylnaphthalene | ND | 0.92 | 0.10 | 1.00 | |
| 2-Methylphenol | ND | 0.92 | 0.16 | 1.00 | |
| 3/4-Methylphenol | ND | 0.92 | 0.30 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 0.92 | 0.15 | 1.00 | |
| N-Nitrosodimethylamine | ND | 0.92 | 0.087 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 0.92 | 0.25 | 1.00 | |
| Naphthalene | ND | 0.92 | 0.11 | 1.00 | |
| 4-Nitroaniline | ND | 0.92 | 0.12 | 1.00 | |
| 3-Nitroaniline | ND | 0.92 | 0.13 | 1.00 | |
| 2-Nitroaniline | ND | 0.92 | 0.095 | 1.00 | |
| Nitrobenzene | ND | 4.6 | 0.60 | 1.00 | |
| 4-Nitrophenol | ND | 0.92 | 0.10 | 1.00 | |
| 2-Nitrophenol | ND | 0.92 | 0.12 | 1.00 | |
| Pentachlorophenol | ND | 4.6 | 0.71 | 1.00 | |
| Phenanthrene | ND | 0.92 | 0.13 | 1.00 | |
| Phenol | ND | 0.92 | 0.088 | 1.00 | |
| Pyrene | ND | 0.92 | 0.14 | 1.00 | |
| Pyridine | ND | 0.92 | 0.10 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 0.92 | 0.11 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 0.92 | 0.14 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 0.92 | 0.12 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 75 | 27-120 | |
| 2-Fluorophenol | 67 | 25-120 | |
| Nitrobenzene-d5 | 64 | 33-123 | |
| p-Terphenyl-d14 | 93 | 27-159 | |
| Phenol-d6 | 61 | 26-122 | |
| 2,4,6-Tribromophenol | 103 | 18-138 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-549-3303 | N/A | Solid | GC/MS TT | 06/11/15 | 06/11/15 16:16 | 150611L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|------|-------|------|------------|
| Acenaphthene | ND | 0.50 | 0.063 | 1.00 | |
| Acenaphthylene | ND | 0.50 | 0.060 | 1.00 | |
| Aniline | ND | 0.50 | 0.060 | 1.00 | |
| Anthracene | ND | 0.50 | 0.063 | 1.00 | |
| Azobenzene | ND | 0.50 | 0.054 | 1.00 | |
| Benzidine | ND | 10 | 0.83 | 1.00 | |
| Benzo (a) Anthracene | ND | 0.50 | 0.057 | 1.00 | |
| Benzo (a) Pyrene | ND | 0.50 | 0.054 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 0.50 | 0.064 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 0.50 | 0.055 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 0.50 | 0.065 | 1.00 | |
| Benzoic Acid | ND | 2.5 | 0.50 | 1.00 | |
| Benzyl Alcohol | ND | 0.50 | 0.066 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 0.50 | 0.056 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 2.5 | 0.41 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 0.50 | 0.057 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 0.50 | 0.053 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 0.50 | 0.061 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 0.50 | 0.054 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 0.50 | 0.066 | 1.00 | |
| 4-Chloroaniline | ND | 0.50 | 0.065 | 1.00 | |
| 2-Chloronaphthalene | ND | 0.50 | 0.059 | 1.00 | |
| 2-Chlorophenol | ND | 0.50 | 0.067 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 0.50 | 0.064 | 1.00 | |
| Chrysene | ND | 0.50 | 0.064 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 0.50 | 0.060 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 0.50 | 0.10 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 0.50 | 0.046 | 1.00 | |
| Dibenzofuran | ND | 0.50 | 0.060 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 0.50 | 0.064 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 0.50 | 0.073 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 0.50 | 0.075 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 10 | 0.36 | 1.00 | |
| 2,4-Dichlorophenol | ND | 0.50 | 0.058 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|------|-------|------|------------|
| Diethyl Phthalate | ND | 0.50 | 0.058 | 1.00 | |
| Dimethyl Phthalate | ND | 0.50 | 0.11 | 1.00 | |
| 2,4-Dimethylphenol | ND | 0.50 | 0.27 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 2.5 | 0.63 | 1.00 | |
| 2,4-Dinitrophenol | ND | 2.5 | 0.34 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 0.50 | 0.067 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 0.50 | 0.073 | 1.00 | |
| Fluoranthene | ND | 0.50 | 0.062 | 1.00 | |
| Fluorene | ND | 0.50 | 0.063 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 0.50 | 0.063 | 1.00 | |
| Hexachlorobenzene | ND | 0.50 | 0.067 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 2.5 | 0.50 | 1.00 | |
| Hexachloroethane | ND | 0.50 | 0.078 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 0.50 | 0.053 | 1.00 | |
| Isophorone | ND | 0.50 | 0.057 | 1.00 | |
| 2-Methylnaphthalene | ND | 0.50 | 0.060 | 1.00 | |
| 1-Methylnaphthalene | ND | 0.50 | 0.056 | 1.00 | |
| 2-Methylphenol | ND | 0.50 | 0.087 | 1.00 | |
| 3/4-Methylphenol | ND | 0.50 | 0.16 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 0.50 | 0.084 | 1.00 | |
| N-Nitrosodimethylamine | ND | 0.50 | 0.047 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 0.50 | 0.14 | 1.00 | |
| Naphthalene | ND | 0.50 | 0.059 | 1.00 | |
| 4-Nitroaniline | ND | 0.50 | 0.065 | 1.00 | |
| 3-Nitroaniline | ND | 0.50 | 0.070 | 1.00 | |
| 2-Nitroaniline | ND | 0.50 | 0.052 | 1.00 | |
| Nitrobenzene | ND | 2.5 | 0.32 | 1.00 | |
| 4-Nitrophenol | ND | 0.50 | 0.054 | 1.00 | |
| 2-Nitrophenol | ND | 0.50 | 0.067 | 1.00 | |
| Pentachlorophenol | ND | 2.5 | 0.39 | 1.00 | |
| Phenanthrene | ND | 0.50 | 0.069 | 1.00 | |
| Phenol | ND | 0.50 | 0.048 | 1.00 | |
| Pyrene | ND | 0.50 | 0.077 | 1.00 | |
| Pyridine | ND | 0.50 | 0.055 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 0.50 | 0.059 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 0.50 | 0.077 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 0.50 | 0.064 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C
Units: mg/kg

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 85 | 27-120 | |
| 2-Fluorophenol | 86 | 25-120 | |
| Nitrobenzene-d5 | 85 | 33-123 | |
| p-Terphenyl-d14 | 95 | 27-159 | |
| Phenol-d6 | 80 | 26-122 | |
| 2,4,6-Tribromophenol | 113 | 18-138 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 1 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| TB08 | 15-06-0567-9-A | 06/05/15 00:00 | Aqueous | GC/MS WW | 06/13/15 | 06/13/15 16:15 | 150613L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 2 of 6

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 111 | 78-126 | |
| 1,2-Dichloroethane-d4 | 110 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 3 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB01 | 15-06-0567-10-A | 06/05/15 15:41 | Aqueous | GC/MS WW | 06/13/15 | 06/13/15 16:48 | 150613L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | 23 | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | 0.27 | 1.0 | 0.24 | 1.00 | J |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 95 | 80-120 | |
| Dibromofluoromethane | 111 | 78-126 | |
| 1,2-Dichloroethane-d4 | 110 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-17386 | N/A | Aqueous | GC/MS WW | 06/13/15 | 06/13/15 15:44 | 150613L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | 0.42 | 10 | 0.41 | 1.00 | J |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 106 | 78-126 | |
| 1,2-Dichloroethane-d4 | 107 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM310 | 15-06-0567-1-A | 06/05/15 10:03 | Solid | GC/MS BB | 06/07/15 | 06/09/15 01:29 | 150608L032 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|-----|------|------|------------|
| Acetone | ND | 450 | 22 | 1.00 | |
| Benzene | ND | 18 | 0.47 | 1.00 | |
| Bromobenzene | ND | 18 | 0.75 | 1.00 | |
| Bromochloromethane | ND | 18 | 2.5 | 1.00 | |
| Bromodichloromethane | ND | 18 | 0.83 | 1.00 | |
| Bromoform | ND | 18 | 2.8 | 1.00 | |
| Bromomethane | ND | 90 | 34 | 1.00 | |
| 2-Butanone | ND | 180 | 14 | 1.00 | |
| n-Butylbenzene | ND | 18 | 0.56 | 1.00 | |
| sec-Butylbenzene | ND | 18 | 2.1 | 1.00 | |
| tert-Butylbenzene | ND | 18 | 0.54 | 1.00 | |
| Carbon Disulfide | 2.2 | 180 | 1.1 | 1.00 | J |
| Carbon Tetrachloride | ND | 18 | 1.0 | 1.00 | |
| Chlorobenzene | ND | 18 | 0.80 | 1.00 | |
| Chloroethane | ND | 18 | 5.3 | 1.00 | |
| Chloroform | ND | 18 | 0.86 | 1.00 | |
| Chloromethane | ND | 90 | 1.1 | 1.00 | |
| 2-Chlorotoluene | ND | 18 | 0.83 | 1.00 | |
| 4-Chlorotoluene | ND | 18 | 0.76 | 1.00 | |
| Dibromochloromethane | ND | 18 | 2.0 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 36 | 6.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 18 | 0.92 | 1.00 | |
| Dibromomethane | ND | 18 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 18 | 0.82 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 18 | 0.63 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 18 | 0.80 | 1.00 | |
| Dichlorodifluoromethane | ND | 18 | 1.6 | 1.00 | |
| 1,1-Dichloroethane | ND | 18 | 0.76 | 1.00 | |
| 1,2-Dichloroethane | ND | 18 | 1.1 | 1.00 | |
| 1,1-Dichloroethene | ND | 18 | 1.2 | 1.00 | |
| c-1,2-Dichloroethene | ND | 18 | 1.0 | 1.00 | |
| t-1,2-Dichloroethene | ND | 18 | 1.8 | 1.00 | |
| 1,2-Dichloropropane | ND | 18 | 1.6 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|-----|------|------|------------|
| 1,3-Dichloropropane | ND | 18 | 0.91 | 1.00 | |
| 2,2-Dichloropropane | ND | 18 | 1.2 | 1.00 | |
| 1,1-Dichloropropene | ND | 18 | 1.2 | 1.00 | |
| c-1,3-Dichloropropene | ND | 18 | 0.91 | 1.00 | |
| t-1,3-Dichloropropene | ND | 18 | 2.2 | 1.00 | |
| Ethylbenzene | ND | 18 | 0.54 | 1.00 | |
| 2-Hexanone | ND | 180 | 6.3 | 1.00 | |
| Isopropylbenzene | ND | 18 | 2.0 | 1.00 | |
| p-Isopropyltoluene | ND | 18 | 2.3 | 1.00 | |
| Methylene Chloride | ND | 180 | 4.8 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 180 | 15 | 1.00 | |
| Naphthalene | ND | 180 | 2.9 | 1.00 | |
| n-Propylbenzene | ND | 18 | 1.8 | 1.00 | |
| Styrene | ND | 18 | 2.2 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 18 | 0.86 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 18 | 1.2 | 1.00 | |
| Tetrachloroethene | ND | 18 | 0.75 | 1.00 | |
| Toluene | ND | 18 | 1.8 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 36 | 3.3 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 18 | 1.1 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 18 | 0.81 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 18 | 1.3 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 180 | 1.3 | 1.00 | |
| Trichloroethene | ND | 18 | 1.1 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 18 | 3.0 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 18 | 2.1 | 1.00 | |
| Trichlorofluoromethane | ND | 180 | 1.3 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 18 | 2.0 | 1.00 | |
| Vinyl Acetate | ND | 180 | 17 | 1.00 | |
| Vinyl Chloride | ND | 18 | 1.8 | 1.00 | |
| p/m-Xylene | ND | 18 | 0.96 | 1.00 | |
| o-Xylene | ND | 18 | 2.0 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 18 | 1.1 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 95 | 60-132 | |
| Dibromofluoromethane | 105 | 63-141 | |
| 1,2-Dichloroethane-d4 | 110 | 62-146 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM320 | 15-06-0567-2-A | 06/05/15 10:48 | Solid | GC/MS Q | 06/07/15 | 06/09/15 05:12 | 150608L054 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|-----|------|------|------------|
| Acetone | ND | 650 | 32 | 1.00 | |
| Benzene | ND | 26 | 0.67 | 1.00 | |
| Bromobenzene | ND | 26 | 1.1 | 1.00 | |
| Bromochloromethane | ND | 26 | 3.6 | 1.00 | |
| Bromodichloromethane | ND | 26 | 1.2 | 1.00 | |
| Bromoform | ND | 26 | 4.1 | 1.00 | |
| Bromomethane | ND | 130 | 49 | 1.00 | |
| 2-Butanone | ND | 260 | 19 | 1.00 | |
| n-Butylbenzene | ND | 26 | 0.81 | 1.00 | |
| sec-Butylbenzene | ND | 26 | 3.0 | 1.00 | |
| tert-Butylbenzene | ND | 26 | 0.78 | 1.00 | |
| Carbon Disulfide | ND | 260 | 1.6 | 1.00 | |
| Carbon Tetrachloride | ND | 26 | 1.5 | 1.00 | |
| Chlorobenzene | ND | 26 | 1.2 | 1.00 | |
| Chloroethane | ND | 26 | 7.7 | 1.00 | |
| Chloroform | ND | 26 | 1.2 | 1.00 | |
| Chloromethane | ND | 130 | 1.6 | 1.00 | |
| 2-Chlorotoluene | ND | 26 | 1.2 | 1.00 | |
| 4-Chlorotoluene | ND | 26 | 1.1 | 1.00 | |
| Dibromochloromethane | ND | 26 | 2.9 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 52 | 9.0 | 1.00 | |
| 1,2-Dibromoethane | ND | 26 | 1.3 | 1.00 | |
| Dibromomethane | ND | 26 | 4.0 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 26 | 1.2 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 26 | 0.91 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 26 | 1.1 | 1.00 | |
| Dichlorodifluoromethane | ND | 26 | 2.3 | 1.00 | |
| 1,1-Dichloroethane | ND | 26 | 1.1 | 1.00 | |
| 1,2-Dichloroethane | ND | 26 | 1.6 | 1.00 | |
| 1,1-Dichloroethene | ND | 26 | 1.8 | 1.00 | |
| c-1,2-Dichloroethene | ND | 26 | 1.4 | 1.00 | |
| t-1,2-Dichloroethene | ND | 26 | 2.6 | 1.00 | |
| 1,2-Dichloropropane | ND | 26 | 2.3 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 1,3-Dichloropropane | ND | 26 | 1.3 | 1.00 | |
| 2,2-Dichloropropane | ND | 26 | 1.7 | 1.00 | |
| 1,1-Dichloropropene | ND | 26 | 1.7 | 1.00 | |
| c-1,3-Dichloropropene | ND | 26 | 1.3 | 1.00 | |
| t-1,3-Dichloropropene | ND | 26 | 3.1 | 1.00 | |
| Ethylbenzene | ND | 26 | 0.78 | 1.00 | |
| 2-Hexanone | ND | 260 | 9.1 | 1.00 | |
| Isopropylbenzene | ND | 26 | 2.8 | 1.00 | |
| p-Isopropyltoluene | ND | 26 | 3.3 | 1.00 | |
| Methylene Chloride | ND | 260 | 6.9 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 260 | 22 | 1.00 | |
| Naphthalene | ND | 260 | 4.2 | 1.00 | |
| n-Propylbenzene | ND | 26 | 2.6 | 1.00 | |
| Styrene | ND | 26 | 3.1 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 26 | 1.2 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 26 | 1.8 | 1.00 | |
| Tetrachloroethene | ND | 26 | 1.1 | 1.00 | |
| Toluene | ND | 26 | 2.7 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 52 | 4.7 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 26 | 1.6 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 26 | 1.2 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 26 | 1.8 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 260 | 1.8 | 1.00 | |
| Trichloroethene | ND | 26 | 1.6 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 26 | 4.3 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 26 | 3.0 | 1.00 | |
| Trichlorofluoromethane | ND | 260 | 1.9 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 26 | 2.8 | 1.00 | |
| Vinyl Acetate | ND | 260 | 25 | 1.00 | |
| Vinyl Chloride | ND | 26 | 2.6 | 1.00 | |
| p/m-Xylene | ND | 26 | 1.4 | 1.00 | |
| o-Xylene | ND | 26 | 2.9 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 26 | 1.5 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 60-132 | |
| Dibromofluoromethane | 97 | 63-141 | |
| 1,2-Dichloroethane-d4 | 93 | 62-146 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM330 | 15-06-0567-3-A | 06/05/15 11:03 | Solid | GC/MS Q | 06/07/15 | 06/09/15 05:39 | 150608L054 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|-----|------|------|------------|
| Acetone | ND | 160 | 7.9 | 1.00 | |
| Benzene | ND | 6.3 | 0.16 | 1.00 | |
| Bromobenzene | ND | 6.3 | 0.26 | 1.00 | |
| Bromochloromethane | ND | 6.3 | 0.87 | 1.00 | |
| Bromodichloromethane | ND | 6.3 | 0.29 | 1.00 | |
| Bromoform | ND | 6.3 | 1.0 | 1.00 | |
| Bromomethane | ND | 32 | 12 | 1.00 | |
| 2-Butanone | ND | 63 | 4.8 | 1.00 | |
| n-Butylbenzene | ND | 6.3 | 0.20 | 1.00 | |
| sec-Butylbenzene | ND | 6.3 | 0.73 | 1.00 | |
| tert-Butylbenzene | ND | 6.3 | 0.19 | 1.00 | |
| Carbon Disulfide | ND | 63 | 0.39 | 1.00 | |
| Carbon Tetrachloride | ND | 6.3 | 0.36 | 1.00 | |
| Chlorobenzene | ND | 6.3 | 0.28 | 1.00 | |
| Chloroethane | ND | 6.3 | 1.9 | 1.00 | |
| Chloroform | ND | 6.3 | 0.30 | 1.00 | |
| Chloromethane | ND | 32 | 0.38 | 1.00 | |
| 2-Chlorotoluene | ND | 6.3 | 0.29 | 1.00 | |
| 4-Chlorotoluene | ND | 6.3 | 0.27 | 1.00 | |
| Dibromochloromethane | ND | 6.3 | 0.72 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 13 | 2.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 6.3 | 0.32 | 1.00 | |
| Dibromomethane | ND | 6.3 | 0.98 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 6.3 | 0.29 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 6.3 | 0.22 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 6.3 | 0.28 | 1.00 | |
| Dichlorodifluoromethane | ND | 6.3 | 0.56 | 1.00 | |
| 1,1-Dichloroethane | ND | 6.3 | 0.27 | 1.00 | |
| 1,2-Dichloroethane | ND | 6.3 | 0.40 | 1.00 | |
| 1,1-Dichloroethene | ND | 6.3 | 0.44 | 1.00 | |
| c-1,2-Dichloroethene | ND | 6.3 | 0.35 | 1.00 | |
| t-1,2-Dichloroethene | ND | 6.3 | 0.64 | 1.00 | |
| 1,2-Dichloropropane | ND | 6.3 | 0.55 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|-----|------|------|------------|
| 1,3-Dichloropropane | ND | 6.3 | 0.32 | 1.00 | |
| 2,2-Dichloropropane | ND | 6.3 | 0.42 | 1.00 | |
| 1,1-Dichloropropene | ND | 6.3 | 0.41 | 1.00 | |
| c-1,3-Dichloropropene | ND | 6.3 | 0.32 | 1.00 | |
| t-1,3-Dichloropropene | ND | 6.3 | 0.77 | 1.00 | |
| Ethylbenzene | ND | 6.3 | 0.19 | 1.00 | |
| 2-Hexanone | ND | 63 | 2.2 | 1.00 | |
| Isopropylbenzene | ND | 6.3 | 0.69 | 1.00 | |
| p-Isopropyltoluene | ND | 6.3 | 0.80 | 1.00 | |
| Methylene Chloride | ND | 63 | 1.7 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 63 | 5.5 | 1.00 | |
| Naphthalene | ND | 63 | 1.0 | 1.00 | |
| n-Propylbenzene | ND | 6.3 | 0.63 | 1.00 | |
| Styrene | ND | 6.3 | 0.76 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 6.3 | 0.30 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 6.3 | 0.44 | 1.00 | |
| Tetrachloroethene | ND | 6.3 | 0.26 | 1.00 | |
| Toluene | ND | 6.3 | 0.65 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 13 | 1.2 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 6.3 | 0.39 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 6.3 | 0.28 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 6.3 | 0.45 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 63 | 0.44 | 1.00 | |
| Trichloroethene | ND | 6.3 | 0.38 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 6.3 | 1.0 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 6.3 | 0.74 | 1.00 | |
| Trichlorofluoromethane | ND | 63 | 0.47 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 6.3 | 0.69 | 1.00 | |
| Vinyl Acetate | ND | 63 | 6.0 | 1.00 | |
| Vinyl Chloride | ND | 6.3 | 0.64 | 1.00 | |
| p/m-Xylene | ND | 6.3 | 0.34 | 1.00 | |
| o-Xylene | ND | 6.3 | 0.70 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 6.3 | 0.37 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 95 | 60-132 | |
| Dibromofluoromethane | 96 | 63-141 | |
| 1,2-Dichloroethane-d4 | 90 | 62-146 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM340 | 15-06-0567-4-A | 06/05/15 11:21 | Solid | GC/MS Q | 06/07/15 | 06/09/15 06:05 | 150608L054 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|-----|------|------|------------|
| Acetone | 11 | 160 | 7.7 | 1.00 | J |
| Benzene | ND | 6.2 | 0.16 | 1.00 | |
| Bromobenzene | ND | 6.2 | 0.26 | 1.00 | |
| Bromochloromethane | ND | 6.2 | 0.86 | 1.00 | |
| Bromodichloromethane | ND | 6.2 | 0.29 | 1.00 | |
| Bromoform | ND | 6.2 | 0.99 | 1.00 | |
| Bromomethane | ND | 31 | 12 | 1.00 | |
| 2-Butanone | ND | 62 | 4.7 | 1.00 | |
| n-Butylbenzene | ND | 6.2 | 0.19 | 1.00 | |
| sec-Butylbenzene | ND | 6.2 | 0.72 | 1.00 | |
| tert-Butylbenzene | ND | 6.2 | 0.19 | 1.00 | |
| Carbon Disulfide | ND | 62 | 0.38 | 1.00 | |
| Carbon Tetrachloride | ND | 6.2 | 0.35 | 1.00 | |
| Chlorobenzene | ND | 6.2 | 0.28 | 1.00 | |
| Chloroethane | ND | 6.2 | 1.9 | 1.00 | |
| Chloroform | ND | 6.2 | 0.30 | 1.00 | |
| Chloromethane | ND | 31 | 0.38 | 1.00 | |
| 2-Chlorotoluene | ND | 6.2 | 0.29 | 1.00 | |
| 4-Chlorotoluene | ND | 6.2 | 0.26 | 1.00 | |
| Dibromochloromethane | ND | 6.2 | 0.71 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 12 | 2.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 6.2 | 0.32 | 1.00 | |
| Dibromomethane | ND | 6.2 | 0.96 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 6.2 | 0.28 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 6.2 | 0.22 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 6.2 | 0.28 | 1.00 | |
| Dichlorodifluoromethane | ND | 6.2 | 0.55 | 1.00 | |
| 1,1-Dichloroethane | ND | 6.2 | 0.26 | 1.00 | |
| 1,2-Dichloroethane | ND | 6.2 | 0.39 | 1.00 | |
| 1,1-Dichloroethene | ND | 6.2 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 6.2 | 0.35 | 1.00 | |
| t-1,2-Dichloroethene | ND | 6.2 | 0.63 | 1.00 | |
| 1,2-Dichloropropane | ND | 6.2 | 0.54 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|----------|----------------|------------|------|------------|
| 1,3-Dichloropropane | ND | 6.2 | 0.31 | 1.00 | |
| 2,2-Dichloropropane | ND | 6.2 | 0.41 | 1.00 | |
| 1,1-Dichloropropene | ND | 6.2 | 0.41 | 1.00 | |
| c-1,3-Dichloropropene | ND | 6.2 | 0.32 | 1.00 | |
| t-1,3-Dichloropropene | ND | 6.2 | 0.75 | 1.00 | |
| Ethylbenzene | ND | 6.2 | 0.19 | 1.00 | |
| 2-Hexanone | ND | 62 | 2.2 | 1.00 | |
| Isopropylbenzene | ND | 6.2 | 0.68 | 1.00 | |
| p-Isopropyltoluene | ND | 6.2 | 0.78 | 1.00 | |
| Methylene Chloride | ND | 62 | 1.7 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 62 | 5.4 | 1.00 | |
| Naphthalene | ND | 62 | 1.0 | 1.00 | |
| n-Propylbenzene | ND | 6.2 | 0.62 | 1.00 | |
| Styrene | ND | 6.2 | 0.75 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 6.2 | 0.30 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 6.2 | 0.43 | 1.00 | |
| Tetrachloroethene | ND | 6.2 | 0.26 | 1.00 | |
| Toluene | ND | 6.2 | 0.64 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 12 | 1.1 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 6.2 | 0.39 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 6.2 | 0.28 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 6.2 | 0.44 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 62 | 0.44 | 1.00 | |
| Trichloroethene | ND | 6.2 | 0.37 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 6.2 | 1.0 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 6.2 | 0.73 | 1.00 | |
| Trichlorofluoromethane | ND | 62 | 0.47 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 6.2 | 0.68 | 1.00 | |
| Vinyl Acetate | ND | 62 | 5.9 | 1.00 | |
| Vinyl Chloride | ND | 6.2 | 0.62 | 1.00 | |
| p/m-Xylene | ND | 6.2 | 0.33 | 1.00 | |
| o-Xylene | ND | 6.2 | 0.69 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 6.2 | 0.37 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| 1,4-Bromofluorobenzene | 94 | 60-132 | | | |
| Dibromofluoromethane | 97 | 63-141 | | | |
| 1,2-Dichloroethane-d4 | 92 | 62-146 | | | |
| Toluene-d8 | 97 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM350 | 15-06-0567-5-A | 06/05/15 12:24 | Solid | GC/MS Q | 06/07/15 | 06/09/15 06:32 | 150608L054 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|-----|------|------|------------|
| Acetone | ND | 270 | 14 | 1.00 | |
| Benzene | ND | 11 | 0.28 | 1.00 | |
| Bromobenzene | ND | 11 | 0.46 | 1.00 | |
| Bromochloromethane | ND | 11 | 1.5 | 1.00 | |
| Bromodichloromethane | ND | 11 | 0.51 | 1.00 | |
| Bromoform | ND | 11 | 1.7 | 1.00 | |
| Bromomethane | ND | 54 | 21 | 1.00 | |
| 2-Butanone | ND | 110 | 8.2 | 1.00 | |
| n-Butylbenzene | ND | 11 | 0.34 | 1.00 | |
| sec-Butylbenzene | ND | 11 | 1.3 | 1.00 | |
| tert-Butylbenzene | ND | 11 | 0.33 | 1.00 | |
| Carbon Disulfide | ND | 110 | 0.67 | 1.00 | |
| Carbon Tetrachloride | ND | 11 | 0.62 | 1.00 | |
| Chlorobenzene | ND | 11 | 0.49 | 1.00 | |
| Chloroethane | ND | 11 | 3.3 | 1.00 | |
| Chloroform | ND | 11 | 0.52 | 1.00 | |
| Chloromethane | ND | 54 | 0.66 | 1.00 | |
| 2-Chlorotoluene | ND | 11 | 0.50 | 1.00 | |
| 4-Chlorotoluene | ND | 11 | 0.46 | 1.00 | |
| Dibromochloromethane | ND | 11 | 1.2 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 22 | 3.8 | 1.00 | |
| 1,2-Dibromoethane | ND | 11 | 0.56 | 1.00 | |
| Dibromomethane | ND | 11 | 1.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 11 | 0.50 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 11 | 0.38 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 11 | 0.48 | 1.00 | |
| Dichlorodifluoromethane | ND | 11 | 0.97 | 1.00 | |
| 1,1-Dichloroethane | ND | 11 | 0.46 | 1.00 | |
| 1,2-Dichloroethane | ND | 11 | 0.68 | 1.00 | |
| 1,1-Dichloroethene | ND | 11 | 0.75 | 1.00 | |
| c-1,2-Dichloroethene | ND | 11 | 0.61 | 1.00 | |
| t-1,2-Dichloroethene | ND | 11 | 1.1 | 1.00 | |
| 1,2-Dichloropropane | ND | 11 | 0.95 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|-----|------|------|------------|
| 1,3-Dichloropropane | ND | 11 | 0.55 | 1.00 | |
| 2,2-Dichloropropane | ND | 11 | 0.72 | 1.00 | |
| 1,1-Dichloropropene | ND | 11 | 0.72 | 1.00 | |
| c-1,3-Dichloropropene | ND | 11 | 0.55 | 1.00 | |
| t-1,3-Dichloropropene | ND | 11 | 1.3 | 1.00 | |
| Ethylbenzene | ND | 11 | 0.33 | 1.00 | |
| 2-Hexanone | ND | 110 | 3.8 | 1.00 | |
| Isopropylbenzene | ND | 11 | 1.2 | 1.00 | |
| p-Isopropyltoluene | ND | 11 | 1.4 | 1.00 | |
| Methylene Chloride | ND | 110 | 2.9 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 110 | 9.4 | 1.00 | |
| Naphthalene | ND | 110 | 1.8 | 1.00 | |
| n-Propylbenzene | ND | 11 | 1.1 | 1.00 | |
| Styrene | ND | 11 | 1.3 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 11 | 0.52 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 11 | 0.75 | 1.00 | |
| Tetrachloroethene | ND | 11 | 0.46 | 1.00 | |
| Toluene | ND | 11 | 1.1 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 22 | 2.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 11 | 0.68 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 11 | 0.49 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 11 | 0.77 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 110 | 0.77 | 1.00 | |
| Trichloroethene | ND | 11 | 0.65 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 11 | 1.8 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 11 | 1.3 | 1.00 | |
| Trichlorofluoromethane | ND | 110 | 0.82 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 11 | 1.2 | 1.00 | |
| Vinyl Acetate | ND | 110 | 10 | 1.00 | |
| Vinyl Chloride | ND | 11 | 1.1 | 1.00 | |
| p/m-Xylene | ND | 11 | 0.58 | 1.00 | |
| o-Xylene | ND | 11 | 1.2 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 11 | 0.64 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 95 | 60-132 | |
| Dibromofluoromethane | 98 | 63-141 | |
| 1,2-Dichloroethane-d4 | 95 | 62-146 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM360 | 15-06-0567-6-A | 06/05/15 13:13 | Solid | GC/MS Q | 06/07/15 | 06/09/15 06:58 | 150608L054 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|-----|------|------|------------|
| Acetone | ND | 170 | 8.6 | 1.00 | |
| Benzene | ND | 6.9 | 0.18 | 1.00 | |
| Bromobenzene | ND | 6.9 | 0.29 | 1.00 | |
| Bromochloromethane | ND | 6.9 | 0.95 | 1.00 | |
| Bromodichloromethane | ND | 6.9 | 0.32 | 1.00 | |
| Bromoform | ND | 6.9 | 1.1 | 1.00 | |
| Bromomethane | ND | 34 | 13 | 1.00 | |
| 2-Butanone | ND | 69 | 5.2 | 1.00 | |
| n-Butylbenzene | ND | 6.9 | 0.22 | 1.00 | |
| sec-Butylbenzene | ND | 6.9 | 0.80 | 1.00 | |
| tert-Butylbenzene | ND | 6.9 | 0.21 | 1.00 | |
| Carbon Disulfide | ND | 69 | 0.42 | 1.00 | |
| Carbon Tetrachloride | ND | 6.9 | 0.39 | 1.00 | |
| Chlorobenzene | ND | 6.9 | 0.31 | 1.00 | |
| Chloroethane | ND | 6.9 | 2.1 | 1.00 | |
| Chloroform | ND | 6.9 | 0.33 | 1.00 | |
| Chloromethane | ND | 34 | 0.42 | 1.00 | |
| 2-Chlorotoluene | ND | 6.9 | 0.32 | 1.00 | |
| 4-Chlorotoluene | ND | 6.9 | 0.29 | 1.00 | |
| Dibromochloromethane | ND | 6.9 | 0.79 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 14 | 2.4 | 1.00 | |
| 1,2-Dibromoethane | ND | 6.9 | 0.35 | 1.00 | |
| Dibromomethane | ND | 6.9 | 1.1 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 6.9 | 0.32 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 6.9 | 0.24 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 6.9 | 0.31 | 1.00 | |
| Dichlorodifluoromethane | ND | 6.9 | 0.61 | 1.00 | |
| 1,1-Dichloroethane | ND | 6.9 | 0.29 | 1.00 | |
| 1,2-Dichloroethane | ND | 6.9 | 0.43 | 1.00 | |
| 1,1-Dichloroethene | ND | 6.9 | 0.48 | 1.00 | |
| c-1,2-Dichloroethene | ND | 6.9 | 0.39 | 1.00 | |
| t-1,2-Dichloroethene | ND | 6.9 | 0.70 | 1.00 | |
| 1,2-Dichloropropane | ND | 6.9 | 0.60 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|-----|------|------|------------|
| 1,3-Dichloropropane | ND | 6.9 | 0.35 | 1.00 | |
| 2,2-Dichloropropane | ND | 6.9 | 0.46 | 1.00 | |
| 1,1-Dichloropropene | ND | 6.9 | 0.45 | 1.00 | |
| c-1,3-Dichloropropene | ND | 6.9 | 0.35 | 1.00 | |
| t-1,3-Dichloropropene | ND | 6.9 | 0.83 | 1.00 | |
| Ethylbenzene | ND | 6.9 | 0.21 | 1.00 | |
| 2-Hexanone | ND | 69 | 2.4 | 1.00 | |
| Isopropylbenzene | ND | 6.9 | 0.75 | 1.00 | |
| p-Isopropyltoluene | ND | 6.9 | 0.87 | 1.00 | |
| Methylene Chloride | ND | 69 | 1.8 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 69 | 6.0 | 1.00 | |
| Naphthalene | ND | 69 | 1.1 | 1.00 | |
| n-Propylbenzene | ND | 6.9 | 0.69 | 1.00 | |
| Styrene | ND | 6.9 | 0.83 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 6.9 | 0.33 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 6.9 | 0.48 | 1.00 | |
| Tetrachloroethene | ND | 6.9 | 0.29 | 1.00 | |
| Toluene | ND | 6.9 | 0.71 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 14 | 1.3 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 6.9 | 0.43 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 6.9 | 0.31 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 6.9 | 0.49 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 69 | 0.48 | 1.00 | |
| Trichloroethene | ND | 6.9 | 0.41 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 6.9 | 1.1 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 6.9 | 0.81 | 1.00 | |
| Trichlorofluoromethane | ND | 69 | 0.52 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 6.9 | 0.76 | 1.00 | |
| Vinyl Acetate | ND | 69 | 6.5 | 1.00 | |
| Vinyl Chloride | ND | 6.9 | 0.69 | 1.00 | |
| p/m-Xylene | ND | 6.9 | 0.37 | 1.00 | |
| o-Xylene | ND | 6.9 | 0.77 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 6.9 | 0.41 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 95 | 60-132 | |
| Dibromofluoromethane | 96 | 63-141 | |
| 1,2-Dichloroethane-d4 | 91 | 62-146 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| HSM370 | 15-06-0567-7-A | 06/05/15 13:51 | Solid | GC/MS Q | 06/07/15 | 06/09/15 07:25 | 150608L054 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|-----|------|------|------------|
| Acetone | 14 | 220 | 11 | 1.00 | J |
| Benzene | ND | 8.7 | 0.23 | 1.00 | |
| Bromobenzene | ND | 8.7 | 0.36 | 1.00 | |
| Bromochloromethane | ND | 8.7 | 1.2 | 1.00 | |
| Bromodichloromethane | ND | 8.7 | 0.40 | 1.00 | |
| Bromoform | ND | 8.7 | 1.4 | 1.00 | |
| Bromomethane | ND | 43 | 16 | 1.00 | |
| 2-Butanone | ND | 87 | 6.5 | 1.00 | |
| n-Butylbenzene | ND | 8.7 | 0.27 | 1.00 | |
| sec-Butylbenzene | ND | 8.7 | 1.0 | 1.00 | |
| tert-Butylbenzene | ND | 8.7 | 0.26 | 1.00 | |
| Carbon Disulfide | ND | 87 | 0.53 | 1.00 | |
| Carbon Tetrachloride | ND | 8.7 | 0.49 | 1.00 | |
| Chlorobenzene | ND | 8.7 | 0.39 | 1.00 | |
| Chloroethane | ND | 8.7 | 2.6 | 1.00 | |
| Chloroform | ND | 8.7 | 0.41 | 1.00 | |
| Chloromethane | ND | 43 | 0.53 | 1.00 | |
| 2-Chlorotoluene | ND | 8.7 | 0.40 | 1.00 | |
| 4-Chlorotoluene | ND | 8.7 | 0.37 | 1.00 | |
| Dibromochloromethane | ND | 8.7 | 0.99 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 17 | 3.0 | 1.00 | |
| 1,2-Dibromoethane | ND | 8.7 | 0.44 | 1.00 | |
| Dibromomethane | ND | 8.7 | 1.3 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 8.7 | 0.40 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 8.7 | 0.31 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 8.7 | 0.39 | 1.00 | |
| Dichlorodifluoromethane | ND | 8.7 | 0.77 | 1.00 | |
| 1,1-Dichloroethane | ND | 8.7 | 0.37 | 1.00 | |
| 1,2-Dichloroethane | ND | 8.7 | 0.54 | 1.00 | |
| 1,1-Dichloroethene | ND | 8.7 | 0.60 | 1.00 | |
| c-1,2-Dichloroethene | ND | 8.7 | 0.49 | 1.00 | |
| t-1,2-Dichloroethene | ND | 8.7 | 0.88 | 1.00 | |
| 1,2-Dichloropropane | ND | 8.7 | 0.76 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

Page 14 of 20

| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|----------|----------------|------------|------|------------|
| 1,3-Dichloropropane | ND | 8.7 | 0.44 | 1.00 | |
| 2,2-Dichloropropane | ND | 8.7 | 0.57 | 1.00 | |
| 1,1-Dichloropropene | ND | 8.7 | 0.57 | 1.00 | |
| c-1,3-Dichloropropene | ND | 8.7 | 0.44 | 1.00 | |
| t-1,3-Dichloropropene | ND | 8.7 | 1.1 | 1.00 | |
| Ethylbenzene | ND | 8.7 | 0.26 | 1.00 | |
| 2-Hexanone | ND | 87 | 3.1 | 1.00 | |
| Isopropylbenzene | ND | 8.7 | 0.95 | 1.00 | |
| p-Isopropyltoluene | ND | 8.7 | 1.1 | 1.00 | |
| Methylene Chloride | ND | 87 | 2.3 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 87 | 7.5 | 1.00 | |
| Naphthalene | ND | 87 | 1.4 | 1.00 | |
| n-Propylbenzene | ND | 8.7 | 0.87 | 1.00 | |
| Styrene | ND | 8.7 | 1.0 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 8.7 | 0.42 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 8.7 | 0.60 | 1.00 | |
| Tetrachloroethene | ND | 8.7 | 0.36 | 1.00 | |
| Toluene | ND | 8.7 | 0.89 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 17 | 1.6 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 8.7 | 0.54 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 8.7 | 0.39 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 8.7 | 0.61 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 87 | 0.61 | 1.00 | |
| Trichloroethene | ND | 8.7 | 0.52 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 8.7 | 1.4 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 8.7 | 1.0 | 1.00 | |
| Trichlorofluoromethane | ND | 87 | 0.65 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 8.7 | 0.95 | 1.00 | |
| Vinyl Acetate | ND | 87 | 8.2 | 1.00 | |
| Vinyl Chloride | ND | 8.7 | 0.87 | 1.00 | |
| p/m-Xylene | ND | 8.7 | 0.46 | 1.00 | |
| o-Xylene | ND | 8.7 | 0.96 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 8.7 | 0.51 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| 1,4-Bromofluorobenzene | 95 | 60-132 | | | |
| Dibromofluoromethane | 97 | 63-141 | | | |
| 1,2-Dichloroethane-d4 | 92 | 62-146 | | | |
| Toluene-d8 | 97 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| FDHSM370 | 15-06-0567-8-A | 06/05/15 13:51 | Solid | GC/MS Q | 06/07/15 | 06/09/15 07:51 | 150608L054 |

Comment(s): - Results are reported on a dry weight basis.

- Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|-----|------|------|------------|
| Acetone | ND | 240 | 12 | 1.00 | |
| Benzene | ND | 9.4 | 0.24 | 1.00 | |
| Bromobenzene | ND | 9.4 | 0.39 | 1.00 | |
| Bromochloromethane | ND | 9.4 | 1.3 | 1.00 | |
| Bromodichloromethane | ND | 9.4 | 0.44 | 1.00 | |
| Bromoform | ND | 9.4 | 1.5 | 1.00 | |
| Bromomethane | ND | 47 | 18 | 1.00 | |
| 2-Butanone | ND | 94 | 7.1 | 1.00 | |
| n-Butylbenzene | ND | 9.4 | 0.29 | 1.00 | |
| sec-Butylbenzene | ND | 9.4 | 1.1 | 1.00 | |
| tert-Butylbenzene | ND | 9.4 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 94 | 0.58 | 1.00 | |
| Carbon Tetrachloride | ND | 9.4 | 0.53 | 1.00 | |
| Chlorobenzene | ND | 9.4 | 0.42 | 1.00 | |
| Chloroethane | ND | 9.4 | 2.8 | 1.00 | |
| Chloroform | ND | 9.4 | 0.45 | 1.00 | |
| Chloromethane | ND | 47 | 0.57 | 1.00 | |
| 2-Chlorotoluene | ND | 9.4 | 0.44 | 1.00 | |
| 4-Chlorotoluene | ND | 9.4 | 0.40 | 1.00 | |
| Dibromochloromethane | ND | 9.4 | 1.1 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 19 | 3.3 | 1.00 | |
| 1,2-Dibromoethane | ND | 9.4 | 0.48 | 1.00 | |
| Dibromomethane | ND | 9.4 | 1.5 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.4 | 0.43 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.4 | 0.33 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.4 | 0.42 | 1.00 | |
| Dichlorodifluoromethane | ND | 9.4 | 0.84 | 1.00 | |
| 1,1-Dichloroethane | ND | 9.4 | 0.40 | 1.00 | |
| 1,2-Dichloroethane | ND | 9.4 | 0.59 | 1.00 | |
| 1,1-Dichloroethene | ND | 9.4 | 0.65 | 1.00 | |
| c-1,2-Dichloroethene | ND | 9.4 | 0.53 | 1.00 | |
| t-1,2-Dichloroethene | ND | 9.4 | 0.95 | 1.00 | |
| 1,2-Dichloropropane | ND | 9.4 | 0.83 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|-----|------|------|------------|
| 1,3-Dichloropropane | ND | 9.4 | 0.48 | 1.00 | |
| 2,2-Dichloropropane | ND | 9.4 | 0.62 | 1.00 | |
| 1,1-Dichloropropene | ND | 9.4 | 0.62 | 1.00 | |
| c-1,3-Dichloropropene | ND | 9.4 | 0.48 | 1.00 | |
| t-1,3-Dichloropropene | ND | 9.4 | 1.1 | 1.00 | |
| Ethylbenzene | ND | 9.4 | 0.29 | 1.00 | |
| 2-Hexanone | ND | 94 | 3.3 | 1.00 | |
| Isopropylbenzene | ND | 9.4 | 1.0 | 1.00 | |
| p-Isopropyltoluene | 1.4 | 9.4 | 1.2 | 1.00 | J |
| Methylene Chloride | ND | 94 | 2.5 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 94 | 8.1 | 1.00 | |
| Naphthalene | ND | 94 | 1.5 | 1.00 | |
| n-Propylbenzene | ND | 9.4 | 0.94 | 1.00 | |
| Styrene | ND | 9.4 | 1.1 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 9.4 | 0.45 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 9.4 | 0.65 | 1.00 | |
| Tetrachloroethene | ND | 9.4 | 0.40 | 1.00 | |
| Toluene | 2.2 | 9.4 | 0.97 | 1.00 | J |
| 1,2,3-Trichlorobenzene | ND | 19 | 1.7 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.4 | 0.58 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 9.4 | 0.42 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 9.4 | 0.67 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 94 | 0.66 | 1.00 | |
| Trichloroethene | ND | 9.4 | 0.57 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 9.4 | 1.6 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 9.4 | 1.1 | 1.00 | |
| Trichlorofluoromethane | ND | 94 | 0.71 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 9.4 | 1.0 | 1.00 | |
| Vinyl Acetate | ND | 94 | 8.9 | 1.00 | |
| Vinyl Chloride | ND | 9.4 | 0.95 | 1.00 | |
| p/m-Xylene | ND | 9.4 | 0.50 | 1.00 | |
| o-Xylene | ND | 9.4 | 1.0 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 9.4 | 0.56 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 95 | 60-132 | |
| Dibromofluoromethane | 97 | 63-141 | |
| 1,2-Dichloroethane-d4 | 93 | 62-146 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-796-9792 | N/A | Solid | GC/MS BB | 06/08/15 | 06/09/15 00:32 | 150608L032 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|-----|------|------|------------|
| Acetone | ND | 120 | 6.2 | 1.00 | |
| Benzene | ND | 5.0 | 0.13 | 1.00 | |
| Bromobenzene | ND | 5.0 | 0.21 | 1.00 | |
| Bromochloromethane | ND | 5.0 | 0.69 | 1.00 | |
| Bromodichloromethane | ND | 5.0 | 0.23 | 1.00 | |
| Bromoform | ND | 5.0 | 0.79 | 1.00 | |
| Bromomethane | ND | 25 | 9.4 | 1.00 | |
| 2-Butanone | ND | 50 | 3.8 | 1.00 | |
| n-Butylbenzene | ND | 5.0 | 0.16 | 1.00 | |
| sec-Butylbenzene | ND | 5.0 | 0.58 | 1.00 | |
| tert-Butylbenzene | ND | 5.0 | 0.15 | 1.00 | |
| Carbon Disulfide | ND | 50 | 0.31 | 1.00 | |
| Carbon Tetrachloride | ND | 5.0 | 0.28 | 1.00 | |
| Chlorobenzene | ND | 5.0 | 0.22 | 1.00 | |
| Chloroethane | ND | 5.0 | 1.5 | 1.00 | |
| Chloroform | ND | 5.0 | 0.24 | 1.00 | |
| Chloromethane | ND | 25 | 0.30 | 1.00 | |
| 2-Chlorotoluene | ND | 5.0 | 0.23 | 1.00 | |
| 4-Chlorotoluene | ND | 5.0 | 0.21 | 1.00 | |
| Dibromochloromethane | ND | 5.0 | 0.57 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 10 | 1.7 | 1.00 | |
| 1,2-Dibromoethane | ND | 5.0 | 0.26 | 1.00 | |
| Dibromomethane | ND | 5.0 | 0.77 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 5.0 | 0.23 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 5.0 | 0.18 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 5.0 | 0.22 | 1.00 | |
| Dichlorodifluoromethane | ND | 5.0 | 0.44 | 1.00 | |
| 1,1-Dichloroethane | ND | 5.0 | 0.21 | 1.00 | |
| 1,2-Dichloroethane | ND | 5.0 | 0.31 | 1.00 | |
| 1,1-Dichloroethene | ND | 5.0 | 0.35 | 1.00 | |
| c-1,2-Dichloroethene | ND | 5.0 | 0.28 | 1.00 | |
| t-1,2-Dichloroethene | ND | 5.0 | 0.51 | 1.00 | |
| 1,2-Dichloropropane | ND | 5.0 | 0.44 | 1.00 | |
| 1,3-Dichloropropane | ND | 5.0 | 0.25 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 5.0 | 0.33 | 1.00 | |
| 1,1-Dichloropropene | ND | 5.0 | 0.33 | 1.00 | |
| c-1,3-Dichloropropene | ND | 5.0 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 5.0 | 0.61 | 1.00 | |
| Ethylbenzene | ND | 5.0 | 0.15 | 1.00 | |
| 2-Hexanone | ND | 50 | 1.8 | 1.00 | |
| Isopropylbenzene | ND | 5.0 | 0.55 | 1.00 | |
| p-Isopropyltoluene | ND | 5.0 | 0.63 | 1.00 | |
| Methylene Chloride | ND | 50 | 1.3 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 50 | 4.3 | 1.00 | |
| Naphthalene | ND | 50 | 0.81 | 1.00 | |
| n-Propylbenzene | ND | 5.0 | 0.50 | 1.00 | |
| Styrene | ND | 5.0 | 0.60 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 5.0 | 0.24 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 5.0 | 0.35 | 1.00 | |
| Tetrachloroethene | ND | 5.0 | 0.21 | 1.00 | |
| Toluene | ND | 5.0 | 0.52 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 10 | 0.91 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 5.0 | 0.31 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 5.0 | 0.23 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 5.0 | 0.35 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 50 | 0.35 | 1.00 | |
| Trichloroethene | ND | 5.0 | 0.30 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.83 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 5.0 | 0.59 | 1.00 | |
| Trichlorofluoromethane | ND | 50 | 0.38 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 5.0 | 0.55 | 1.00 | |
| Vinyl Acetate | ND | 50 | 4.7 | 1.00 | |
| Vinyl Chloride | ND | 5.0 | 0.50 | 1.00 | |
| p/m-Xylene | ND | 5.0 | 0.27 | 1.00 | |
| o-Xylene | ND | 5.0 | 0.56 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 5.0 | 0.30 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 96 | 60-132 | |
| Dibromofluoromethane | 105 | 63-141 | |
| 1,2-Dichloroethane-d4 | 110 | 62-146 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|--------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-796-9818 | N/A | Solid | GC/MS Q | 06/08/15 | 06/09/15 02:06 | 150608L054 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|-----|------|------|------------|
| Acetone | ND | 120 | 6.2 | 1.00 | |
| Benzene | ND | 5.0 | 0.13 | 1.00 | |
| Bromobenzene | ND | 5.0 | 0.21 | 1.00 | |
| Bromochloromethane | ND | 5.0 | 0.69 | 1.00 | |
| Bromodichloromethane | ND | 5.0 | 0.23 | 1.00 | |
| Bromoform | ND | 5.0 | 0.79 | 1.00 | |
| Bromomethane | ND | 25 | 9.4 | 1.00 | |
| 2-Butanone | ND | 50 | 3.8 | 1.00 | |
| n-Butylbenzene | ND | 5.0 | 0.16 | 1.00 | |
| sec-Butylbenzene | ND | 5.0 | 0.58 | 1.00 | |
| tert-Butylbenzene | ND | 5.0 | 0.15 | 1.00 | |
| Carbon Disulfide | ND | 50 | 0.31 | 1.00 | |
| Carbon Tetrachloride | ND | 5.0 | 0.28 | 1.00 | |
| Chlorobenzene | ND | 5.0 | 0.22 | 1.00 | |
| Chloroethane | ND | 5.0 | 1.5 | 1.00 | |
| Chloroform | ND | 5.0 | 0.24 | 1.00 | |
| Chloromethane | ND | 25 | 0.30 | 1.00 | |
| 2-Chlorotoluene | ND | 5.0 | 0.23 | 1.00 | |
| 4-Chlorotoluene | ND | 5.0 | 0.21 | 1.00 | |
| Dibromochloromethane | ND | 5.0 | 0.57 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 10 | 1.7 | 1.00 | |
| 1,2-Dibromoethane | ND | 5.0 | 0.26 | 1.00 | |
| Dibromomethane | ND | 5.0 | 0.77 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 5.0 | 0.23 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 5.0 | 0.18 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 5.0 | 0.22 | 1.00 | |
| Dichlorodifluoromethane | ND | 5.0 | 0.44 | 1.00 | |
| 1,1-Dichloroethane | ND | 5.0 | 0.21 | 1.00 | |
| 1,2-Dichloroethane | ND | 5.0 | 0.31 | 1.00 | |
| 1,1-Dichloroethene | ND | 5.0 | 0.35 | 1.00 | |
| c-1,2-Dichloroethene | ND | 5.0 | 0.28 | 1.00 | |
| t-1,2-Dichloroethene | ND | 5.0 | 0.51 | 1.00 | |
| 1,2-Dichloropropane | ND | 5.0 | 0.44 | 1.00 | |
| 1,3-Dichloropropane | ND | 5.0 | 0.25 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/kg

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 5.0 | 0.33 | 1.00 | |
| 1,1-Dichloropropene | ND | 5.0 | 0.33 | 1.00 | |
| c-1,3-Dichloropropene | ND | 5.0 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 5.0 | 0.61 | 1.00 | |
| Ethylbenzene | ND | 5.0 | 0.15 | 1.00 | |
| 2-Hexanone | ND | 50 | 1.8 | 1.00 | |
| Isopropylbenzene | ND | 5.0 | 0.55 | 1.00 | |
| p-Isopropyltoluene | ND | 5.0 | 0.63 | 1.00 | |
| Methylene Chloride | ND | 50 | 1.3 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 50 | 4.3 | 1.00 | |
| Naphthalene | ND | 50 | 0.81 | 1.00 | |
| n-Propylbenzene | ND | 5.0 | 0.50 | 1.00 | |
| Styrene | ND | 5.0 | 0.60 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 5.0 | 0.24 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 5.0 | 0.35 | 1.00 | |
| Tetrachloroethene | ND | 5.0 | 0.21 | 1.00 | |
| Toluene | ND | 5.0 | 0.52 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 10 | 0.91 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 5.0 | 0.31 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 5.0 | 0.23 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 5.0 | 0.35 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 50 | 0.35 | 1.00 | |
| Trichloroethene | ND | 5.0 | 0.30 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.83 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 5.0 | 0.59 | 1.00 | |
| Trichlorofluoromethane | ND | 50 | 0.38 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 5.0 | 0.55 | 1.00 | |
| Vinyl Acetate | ND | 50 | 4.7 | 1.00 | |
| Vinyl Chloride | ND | 5.0 | 0.50 | 1.00 | |
| p/m-Xylene | ND | 5.0 | 0.27 | 1.00 | |
| o-Xylene | ND | 5.0 | 0.56 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 5.0 | 0.30 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 96 | 60-132 | |
| Dibromofluoromethane | 95 | 63-141 | |
| 1,2-Dichloroethane-d4 | 94 | 62-146 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

06/06/15

Work Order:

15-06-0567

Page 1 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|--------------|
| HSM310 | 15-06-0567-1 | 06/05/15 10:03 | Solid |

Comment(s): (9) - Results are reported on a dry weight basis.

(24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 71.3 | 0.100 | | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 7.71 | 0.01 | | 1.00 | | pH units | 06/06/15 | 06/06/15 | EPA 9045D |
| Carbon, Total Organic (9, 24) | 55000 | 1700 | 420 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | EPA 9060A |
| Alkalinity, Total (as CaCO3) (24) | 430 | 5.0 | 3.2 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO3) (24) | 430 | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO3) (24) | ND | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved (24) | 1400 | 10.0 | 1.00 | 1.00 | | mg/kg | 06/16/15 | 06/16/15 | SM 2540 C (M) |
| Phosphorus, Total (9, 24) | 23 | 3.5 | 0.82 | 2.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

| | | | |
|---------------|---------------------|-----------------------|--------------|
| HSM320 | 15-06-0567-2 | 06/05/15 10:48 | Solid |
|---------------|---------------------|-----------------------|--------------|

Comment(s): (9) - Results are reported on a dry weight basis.

(24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 80.4 | 0.100 | | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 7.51 | 0.01 | | 1.00 | | pH units | 06/06/15 | 06/06/15 | EPA 9045D |
| Carbon, Total Organic (9, 24) | 110000 | 2600 | 620 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | EPA 9060A |
| Alkalinity, Total (as CaCO3) (24) | 520 | 5.0 | 3.2 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO3) (24) | 520 | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO3) (24) | ND | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved (24) | 2260 | 10.0 | 1.00 | 1.00 | | mg/kg | 06/16/15 | 06/16/15 | SM 2540 C (M) |
| Phosphorus, Total (9, 24) | 200 | 51 | 12 | 20.0 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

| | | | |
|---------------|---------------------|-----------------------|--------------|
| HSM330 | 15-06-0567-3 | 06/05/15 11:03 | Solid |
|---------------|---------------------|-----------------------|--------------|

Comment(s): (9) - Results are reported on a dry weight basis.

(24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 20.5 | 0.100 | | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 8.14 | 0.01 | | 1.00 | | pH units | 06/06/15 | 06/06/15 | EPA 9045D |
| Carbon, Total Organic (9, 24) | 10000 | 630 | 150 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | EPA 9060A |
| Alkalinity, Total (as CaCO3) (24) | 300 | 5.0 | 3.2 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO3) (24) | 300 | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO3) (24) | ND | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved (24) | 6490 | 10.0 | 1.00 | 1.00 | | mg/kg | 06/16/15 | 06/16/15 | SM 2540 C (M) |
| Phosphorus, Total (9, 24) | 3.1 | 0.63 | 0.15 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

06/06/15

Work Order:

15-06-0567

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|--------------|
| HSM340 | 15-06-0567-4 | 06/05/15 11:21 | Solid |

Comment(s): (9) - Results are reported on a dry weight basis.

(24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 19.8 | 0.100 | | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 7.53 | 0.01 | | 1.00 | | pH units | 06/06/15 | 06/06/15 | EPA 9045D |
| Carbon, Total Organic (9, 24) | 14000 | 620 | 150 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | EPA 9060A |
| Alkalinity, Total (as CaCO3) (24) | 390 | 5.0 | 3.2 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO3) (24) | 390 | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO3) (24) | ND | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved (24) | 2780 | 10.0 | 1.00 | 1.00 | | mg/kg | 06/16/15 | 06/16/15 | SM 2540 C (M) |
| Phosphorus, Total (9, 24) | 0.84 | 0.62 | 0.15 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

| | | | |
|---------------|---------------------|-----------------------|--------------|
| HSM350 | 15-06-0567-5 | 06/05/15 12:24 | Solid |
|---------------|---------------------|-----------------------|--------------|

Comment(s): (9) - Results are reported on a dry weight basis.

(24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 54.1 | 0.100 | | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 7.32 | 0.01 | | 1.00 | | pH units | 06/06/15 | 06/06/15 | EPA 9045D |
| Carbon, Total Organic (9, 24) | 59000 | 1100 | 260 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | EPA 9060A |
| Alkalinity, Total (as CaCO3) (24) | 640 | 5.0 | 3.2 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO3) (24) | 640 | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO3) (24) | ND | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved (24) | 2090 | 10.0 | 1.00 | 1.00 | | mg/kg | 06/16/15 | 06/16/15 | SM 2540 C (M) |
| Phosphorus, Total (9, 24) | 4.8 | 1.1 | 0.26 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

| | | | |
|---------------|---------------------|-----------------------|--------------|
| HSM360 | 15-06-0567-6 | 06/05/15 13:13 | Solid |
|---------------|---------------------|-----------------------|--------------|

Comment(s): (9) - Results are reported on a dry weight basis.

(24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 25.5 | 0.100 | | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 7.71 | 0.01 | | 1.00 | | pH units | 06/06/15 | 06/06/15 | EPA 9045D |
| Carbon, Total Organic (9, 24) | 8700 | 670 | 160 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | EPA 9060A |
| Alkalinity, Total (as CaCO3) (24) | 400 | 5.0 | 3.2 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO3) (24) | 400 | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO3) (24) | ND | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved (24) | 9140 | 10.0 | 1.00 | 1.00 | | mg/kg | 06/16/15 | 06/16/15 | SM 2540 C (M) |
| Phosphorus, Total (9, 24) | 2.0 | 0.67 | 0.16 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

06/06/15

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15-06-0567

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|--------------|
| HSM370 | 15-06-0567-7 | 06/05/15 13:51 | Solid |

Comment(s): (9) - Results are reported on a dry weight basis.
 (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.
 (306) - Please see Work Order Narrative, additional comments section.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 41.5 | 0.100 | | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 7.46 | 0.01 | | 1.00 | | pH units | 06/06/15 | 06/06/15 | EPA 9045D |
| Carbon, Total Organic (9, 24, 306) | 25000 | 850 | 210 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | EPA 9060A |
| Alkalinity, Total (as CaCO ₃) (24) | 440 | 5.0 | 3.2 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO ₃) (24) | 440 | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO ₃) (24) | ND | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved (24) | 5500 | 10.0 | 1.00 | 1.00 | | mg/kg | 06/16/15 | 06/16/15 | SM 2540 C (M) |
| Phosphorus, Total (9, 24) | 3.1 | 0.85 | 0.20 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

| | | | |
|-----------------|---------------------|-----------------------|--------------|
| FDHSM370 | 15-06-0567-8 | 06/05/15 13:51 | Solid |
|-----------------|---------------------|-----------------------|--------------|

Comment(s): (9) - Results are reported on a dry weight basis.
 (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.
 (306) - Please see Work Order Narrative, additional comments section.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|------|------|------------|----------|---------------|---------------|-------------------|
| Moisture | 45.7 | 0.100 | | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| pH | 7.49 | 0.01 | | 1.00 | | pH units | 06/06/15 | 06/06/15 | EPA 9045D |
| Carbon, Total Organic (9, 24, 306) | 22000 | 920 | 220 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | EPA 9060A |
| Alkalinity, Total (as CaCO ₃) (24) | 360 | 5.0 | 3.2 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO ₃) (24) | 360 | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO ₃) (24) | ND | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved (24) | 7660 | 10.0 | 1.00 | 1.00 | | mg/kg | 06/16/15 | 06/16/15 | SM 2540 C (M) |
| Phosphorus, Total (9, 24) | 3.5 | 0.92 | 0.22 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

06/06/15

Work Order:

15-06-0567

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|----------------------|-----------------------|----------------|
| EB01 | 15-06-0567-10 | 06/05/15 15:41 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|--------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 06/13/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 06/11/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 06/11/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 06/11/15 | SM 2320B |
| Solids, Total Dissolved (24) | ND | 0 | 0.870 | 1.00 | | mg/L | 06/11/15 | 06/11/15 | SM 2540 C |
| pH | 7.69 | 0.01 | | 1.00 | BV,BU | pH units | N/A | 06/06/15 | SM 4500 H+ B |
| Carbon, Total Organic (24) | 0.55 | 0.50 | 0.24 | 1.00 | | mg/L | 06/19/15 | 06/19/15 | SM 5310 B |

| Method Blank | N/A | Solid |
|--------------|-----|-------|
|--------------|-----|-------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|-------|---------------|---------------|-------------------|
| Moisture | ND | 0.100 | | 1.00 | | % | 06/08/15 | 06/08/15 | ASTM D-2216 (M) |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 06/13/15 | EPA 365.1 |
| Carbon, Total Organic (24) | ND | 500 | 120 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | EPA 9060A |
| Alkalinity, Total (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 06/11/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 06/11/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 06/11/15 | SM 2320B |
| Alkalinity, Total (as CaCO3) (24) | ND | 5.0 | 3.2 | 1.00 | | mg/kg | N/A | 06/15/15 | SM 2320B M |
| Bicarbonate (as CaCO3) (24) | ND | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Carbonate (as CaCO3) (24) | ND | 5.0 | 4.2 | 1.00 | | mg/kg | 06/15/15 | 06/15/15 | SM 2320B M |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.87 | 1.00 | | mg/L | 06/11/15 | 06/11/15 | SM 2540 C |
| Solids, Total Dissolved (24) | ND | 1.0 | 1.0 | 1.00 | | mg/kg | 06/16/15 | 06/16/15 | SM 2540 C (M) |
| Phosphorus, Total (24) | ND | 0.50 | 0.12 | 1.00 | | mg/kg | 06/10/15 | 06/11/15 | SM 4500 P B/E (M) |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 06/19/15 | 06/19/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-06-0570-2 | Sample | Aqueous | IC 7 | N/A | 06/06/15 12:37 | 150606S01 |
| 15-06-0570-2 | Matrix Spike | Aqueous | IC 7 | N/A | 06/06/15 13:43 | 150606S01 |
| 15-06-0570-2 | Matrix Spike Duplicate | Aqueous | IC 7 | N/A | 06/06/15 13:59 | 150606S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.6360 | 250.0 | 259.8 | 104 | 258.0 | 103 | 80-120 | 1 | 0-20 | |
| Chloride | 40.36 | 5000 | 4867 | 97 | 4867 | 97 | 80-120 | 0 | 0-20 | |
| Bromide | ND | 500.0 | 478.0 | 96 | 475.1 | 95 | 80-120 | 1 | 0-20 | |
| Nitrate (as N) | ND | 500.0 | 486.3 | 97 | 486.2 | 97 | 80-120 | 0 | 0-20 | |
| Sulfate | 89.02 | 5000 | 4968 | 98 | 4979 | 98 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|--------|------------|---------------|----------------|---------------------|
| 15-06-0536-1 | Sample | Solid | IC 10 | 06/09/15 | 06/09/15 14:26 | 150609S01P |
| 15-06-0536-1 | Matrix Spike | Solid | IC 10 | 06/09/15 | 06/09/15 16:37 | 150609S01P |
| 15-06-0536-1 | Matrix Spike Duplicate | Solid | IC 10 | 06/09/15 | 06/09/15 16:54 | 150609S01P |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 1.260 | 25.00 | 23.80 | 90 | 24.02 | 91 | 80-120 | 1 | 0-20 | |
| Chloride | ND | 500.0 | 539.0 | 108 | 520.6 | 104 | 80-120 | 3 | 0-20 | |
| Bromide | ND | 50.00 | 52.63 | 105 | 50.57 | 101 | 80-120 | 4 | 0-20 | |
| Nitrate (as N) | ND | 50.00 | 53.75 | 108 | 51.61 | 103 | 80-120 | 4 | 0-20 | |
| Sulfate | 205.6 | 500.0 | 788.5 | 117 | 755.4 | 110 | 80-120 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
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San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-06-0884-1 | Sample | Aqueous | ACA 1 | N/A | 06/13/15 14:38 | 150613S01 |
| 15-06-0884-1 | Matrix Spike | Aqueous | ACA 1 | N/A | 06/13/15 14:38 | 150613S01 |
| 15-06-0884-1 | Matrix Spike Duplicate | Aqueous | ACA 1 | N/A | 06/13/15 14:38 | 150613S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | ND | 0.2000 | 0.2211 | 111 | 0.2262 | 113 | 90-110 | 2 | 0-25 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

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Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 9060A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|--------|------------|---------------|----------------|---------------------|
| 15-06-0424-1 | Sample | Solid | TOC 5 | 06/15/15 | 06/15/15 16:48 | F0615TOCS1 |
| 15-06-0424-1 | Matrix Spike | Solid | TOC 5 | 06/15/15 | 06/15/15 16:48 | F0615TOCS1 |
| 15-06-0424-1 | Matrix Spike Duplicate | Solid | TOC 5 | 06/15/15 | 06/15/15 16:48 | F0615TOCS1 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Carbon, Total Organic | 600.0 | 30000 | 26100 | 85 | 24200 | 79 | 75-125 | 8 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
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San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 4500 P B/E (M)

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|--------|------------|---------------|----------------|---------------------|
| HSM320 | Sample | Solid | UV 7 | 06/10/15 | 06/11/15 16:26 | F0611TPS2 |
| HSM320 | Matrix Spike | Solid | UV 7 | 06/10/15 | 06/11/15 16:26 | F0611TPS2 |
| HSM320 | Matrix Spike Duplicate | Solid | UV 7 | 06/10/15 | 06/11/15 16:26 | F0611TPS2 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | 39.46 | 40.00 | 80.30 | 102 | 80.70 | 103 | 70-130 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

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San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-06-0884-1 | Sample | Aqueous | TOC 8 | 06/19/15 | 06/19/15 22:52 | F0619TOCS1 |
| 15-06-0884-1 | Matrix Spike | Aqueous | TOC 8 | 06/19/15 | 06/19/15 22:52 | F0619TOCS1 |
| 15-06-0884-1 | Matrix Spike Duplicate | Aqueous | TOC 8 | 06/19/15 | 06/19/15 22:52 | F0619TOCS1 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Carbon, Total Organic | 0.7770 | 10.00 | 9.700 | 89 | 9.420 | 86 | 31-145 | 3 | 0-23 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
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San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|------------------|---------------------|------------|---------------|-------------------|
| HSM310 | Sample | | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:13 | 150611S09A | | | |
| HSM310 | Matrix Spike | | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:15 | 150611S09A | | | |
| HSM310 | Matrix Spike Duplicate | | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:18 | 150611S09A | | | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Calcium | 43080 | 25.00 | 43440 | 4X | 45040 | 4X | 75-125 | 4X | 0-20 | Q |
| Magnesium | 1104 | 25.00 | 1152 | 4X | 1167 | 4X | 75-125 | 4X | 0-20 | Q |
| Potassium | 863.7 | 250.0 | 1192 | 131 | 1199 | 134 | 75-125 | 1 | 0-20 | 3 |
| Sodium | 95.79 | 250.0 | 421.6 | 130 | 426.0 | 132 | 75-125 | 1 | 0-20 | 3 |
| Strontium | 84.54 | 25.00 | 115.4 | 124 | 116.4 | 128 | 75-125 | 1 | 0-20 | 3 |
| Silicon | 915.1 | 25.00 | 942.5 | 4X | 957.0 | 4X | 75-125 | 4X | 0-20 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3010A Total
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-06-1010-11 | Sample | Aqueous | ICP 7300 | 06/15/15 | 06/16/15 14:06 | 150615SA4 |
| 15-06-1010-11 | Matrix Spike | Aqueous | ICP 7300 | 06/15/15 | 06/16/15 14:08 | 150615SA4 |
| 15-06-1010-11 | Matrix Spike Duplicate | Aqueous | ICP 7300 | 06/15/15 | 06/16/15 14:09 | 150615SA4 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Calcium | 36.00 | 0.5000 | 36.02 | 4X | 36.11 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 79.75 | 0.5000 | 79.89 | 4X | 79.35 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 0.5344 | 5.000 | 5.794 | 105 | 6.047 | 110 | 83-131 | 4 | 0-7 | |
| Sodium | 44.98 | 5.000 | 49.93 | 4X | 50.05 | 4X | 73-127 | 4X | 0-9 | Q |
| Strontium | 1.051 | 0.5000 | 1.577 | 105 | 1.604 | 111 | 81-123 | 2 | 0-6 | |
| Silicon | 13.42 | 0.5000 | 13.81 | 4X | 13.58 | 4X | 24-180 | 4X | 0-15 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-06-0965-1 | Sample | Solid | ICP/MS 03 | 06/12/15 | 06/12/15 17:47 | 150612S01 | | | | |
| 15-06-0965-1 | Matrix Spike | Solid | ICP/MS 03 | 06/12/15 | 06/12/15 17:32 | 150612S01 | | | | |
| 15-06-0965-1 | Matrix Spike Duplicate | Solid | ICP/MS 03 | 06/12/15 | 06/12/15 17:36 | 150612S01 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 25.00 | 9.499 | 38 | 9.348 | 37 | 1-97 | 2 | 0-39 | |
| Arsenic | 3.329 | 25.00 | 29.78 | 106 | 29.45 | 104 | 72-132 | 1 | 0-13 | |
| Barium | 192.6 | 25.00 | 237.7 | 4X | 235.3 | 4X | 50-152 | 4X | 0-41 | Q |
| Beryllium | ND | 25.00 | 28.23 | 113 | 27.82 | 111 | 61-121 | 1 | 0-13 | |
| Cadmium | ND | 25.00 | 28.73 | 115 | 28.04 | 112 | 85-121 | 2 | 0-12 | |
| Chromium | 37.93 | 25.00 | 66.38 | 114 | 65.17 | 109 | 20-182 | 2 | 0-15 | |
| Copper | 29.39 | 25.00 | 58.57 | 117 | 57.61 | 113 | 25-157 | 2 | 0-22 | |
| Lead | 9.600 | 25.00 | 36.59 | 108 | 36.60 | 108 | 62-134 | 0 | 0-23 | |
| Nickel | 25.79 | 25.00 | 52.61 | 107 | 50.85 | 100 | 46-154 | 3 | 0-15 | |
| Selenium | ND | 25.00 | 27.57 | 110 | 27.27 | 109 | 54-132 | 1 | 0-14 | |
| Silver | ND | 12.50 | 14.00 | 112 | 13.43 | 107 | 78-126 | 4 | 0-15 | |
| Thallium | ND | 25.00 | 24.81 | 99 | 24.97 | 100 | 79-115 | 1 | 0-11 | |
| Zinc | 96.60 | 25.00 | 131.5 | 140 | 131.0 | 138 | 23-173 | 0 | 0-18 | |
| Aluminum | 19620 | 25.00 | 22610 | 4X | 22010 | 4X | 80-120 | 4X | 0-20 | Q |
| Iron | 30610 | 25.00 | 32270 | 4X | 30950 | 4X | 80-120 | 4X | 0-20 | Q |
| Manganese | 312.0 | 25.00 | 320.6 | 4X | 313.9 | 4X | 80-120 | 4X | 0-20 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|---------------------------|------------------------|-------------|----------|------------|---------------|----------------|---------------------|-----|--------|------------|
| 15-06-0423-2 | Sample | | Aqueous | ICP/MS 04 | 06/08/15 | 06/09/15 17:26 | 150608SA2A | | | |
| 15-06-0423-2 | Matrix Spike | | Aqueous | ICP/MS 04 | 06/08/15 | 06/09/15 17:07 | 150608SA2A | | | |
| 15-06-0423-2 | Matrix Spike Duplicate | | Aqueous | ICP/MS 04 | 06/08/15 | 06/09/15 17:11 | 150608SA2A | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.09236 | 92 | 0.08643 | 86 | 85-133 | 7 | 0-11 | |
| Arsenic | ND | 0.1000 | 0.08479 | 85 | 0.08075 | 81 | 73-127 | 5 | 0-11 | |
| Barium | 0.05909 | 0.1000 | 0.1576 | 99 | 0.1571 | 98 | 74-128 | 0 | 0-10 | |
| Beryllium | ND | 0.1000 | 0.09956 | 100 | 0.09269 | 93 | 56-122 | 7 | 0-11 | |
| Cadmium | ND | 0.1000 | 0.08748 | 87 | 0.08354 | 84 | 84-114 | 5 | 0-8 | |
| Chromium | 0.001241 | 0.1000 | 0.1117 | 110 | 0.1002 | 99 | 73-133 | 11 | 0-11 | |
| Copper | 0.001283 | 0.1000 | 0.09526 | 94 | 0.09005 | 89 | 72-108 | 6 | 0-10 | |
| Lead | ND | 0.1000 | 0.1092 | 109 | 0.1029 | 103 | 79-121 | 6 | 0-10 | |
| Nickel | 0.006589 | 0.1000 | 0.09759 | 91 | 0.09413 | 88 | 68-122 | 4 | 0-10 | |
| Selenium | 0.001400 | 0.1000 | 0.07889 | 77 | 0.07473 | 73 | 59-125 | 5 | 0-12 | |
| Silver | ND | 0.05000 | 0.04526 | 91 | 0.04702 | 94 | 68-128 | 4 | 0-14 | |
| Thallium | ND | 0.1000 | 0.1080 | 108 | 0.1008 | 101 | 73-121 | 7 | 0-11 | |
| Zinc | ND | 0.1000 | 0.08150 | 82 | 0.07452 | 75 | 43-145 | 9 | 0-39 | |
| Aluminum | ND | 0.1000 | 0.1043 | 104 | 0.1024 | 102 | 47-161 | 2 | 0-24 | |
| Iron | 0.08723 | 5.100 | 5.337 | 103 | 5.329 | 103 | 27-201 | 0 | 0-24 | |
| Manganese | 0.03515 | 0.1000 | 0.1365 | 101 | 0.1322 | 97 | 72-126 | 3 | 0-42 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 7470A Total
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-06-0577-1 | Sample | Aqueous | Mercury 04 | 06/08/15 | 06/08/15 13:59 | 150608S01 |
| 15-06-0577-1 | Matrix Spike | Aqueous | Mercury 04 | 06/08/15 | 06/08/15 14:01 | 150608S01 |
| 15-06-0577-1 | Matrix Spike Duplicate | Aqueous | Mercury 04 | 06/08/15 | 06/08/15 14:03 | 150608S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | ND | 0.01000 | 0.01008 | 101 | 0.009980 | 100 | 57-141 | 1 | 0-10 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 7471A Total
Method: EPA 7471A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|--------|------------|---------------|----------------|---------------------|
| 15-06-1080-1 | Sample | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:12 | 150617S03 |
| 15-06-1080-1 | Matrix Spike | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:14 | 150617S03 |
| 15-06-1080-1 | Matrix Spike Duplicate | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:16 | 150617S03 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | 0.6237 | 0.8350 | 1.383 | 91 | 1.596 | 116 | 71-137 | 14 | 0-14 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|--------|------------|---------------|----------------|---------------------|
| 15-06-0778-2 | Sample | Solid | GC 41 | 06/10/15 | 06/11/15 19:49 | 150610S17 |
| 15-06-0778-2 | Matrix Spike | Solid | GC 41 | 06/10/15 | 06/11/15 19:02 | 150610S17 |
| 15-06-0778-2 | Matrix Spike Duplicate | Solid | GC 41 | 06/10/15 | 06/11/15 19:18 | 150610S17 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|--------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Aldrin | ND | 25.00 | 22.05 | 88 | 20.40 | 82 | 50-135 | 8 | 0-25 | |
| Alpha-BHC | ND | 25.00 | 20.34 | 81 | 19.27 | 77 | 50-135 | 5 | 0-25 | |
| Beta-BHC | ND | 25.00 | 20.16 | 81 | 20.56 | 82 | 50-135 | 2 | 0-25 | |
| 4,4'-DDD | 7.030 | 25.00 | 23.78 | 67 | 25.52 | 74 | 50-135 | 7 | 0-25 | |
| 4,4'-DDE | ND | 25.00 | 21.65 | 87 | 21.54 | 86 | 50-135 | 1 | 0-25 | |
| 4,4'-DDT | ND | 25.00 | 21.19 | 85 | 19.94 | 80 | 50-135 | 6 | 0-25 | |
| Delta-BHC | ND | 25.00 | 19.98 | 80 | 19.62 | 78 | 50-135 | 2 | 0-25 | |
| Dieldrin | ND | 25.00 | 21.49 | 86 | 20.07 | 80 | 50-135 | 7 | 0-25 | |
| Endosulfan I | ND | 25.00 | 18.67 | 75 | 17.22 | 69 | 50-135 | 8 | 0-25 | |
| Endosulfan II | ND | 25.00 | 19.02 | 76 | 17.54 | 70 | 50-135 | 8 | 0-25 | |
| Endosulfan Sulfate | ND | 25.00 | 18.09 | 72 | 17.33 | 69 | 50-135 | 4 | 0-25 | |
| Endrin | ND | 25.00 | 20.06 | 80 | 19.03 | 76 | 50-135 | 5 | 0-25 | |
| Endrin Aldehyde | ND | 25.00 | 16.33 | 65 | 16.32 | 65 | 50-135 | 0 | 0-25 | |
| Gamma-BHC | ND | 25.00 | 20.45 | 82 | 19.27 | 77 | 50-135 | 6 | 0-25 | |
| Heptachlor | ND | 25.00 | 19.92 | 80 | 18.70 | 75 | 50-135 | 6 | 0-25 | |
| Heptachlor Epoxide | ND | 25.00 | 18.65 | 75 | 17.54 | 70 | 50-135 | 6 | 0-25 | |
| Methoxychlor | ND | 25.00 | 18.88 | 76 | 18.53 | 74 | 50-135 | 2 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|--------|------------|---------------|----------------|---------------------|
| 15-06-0778-2 | Sample | Solid | GC 58 | 06/10/15 | 06/11/15 23:34 | 150610S18 |
| 15-06-0778-2 | Matrix Spike | Solid | GC 58 | 06/10/15 | 06/11/15 22:40 | 150610S18 |
| 15-06-0778-2 | Matrix Spike Duplicate | Solid | GC 58 | 06/10/15 | 06/11/15 22:58 | 150610S18 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|--------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Aroclor-1016 | ND | 100.0 | 116.1 | 116 | 128.6 | 129 | 50-135 | 10 | 0-20 | |
| Aroclor-1260 | ND | 100.0 | 98.04 | 98 | 106.2 | 106 | 50-135 | 8 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-05-2040-31 | Sample | Solid | GC 26 | 06/09/15 | 06/12/15 14:21 | 150609S20 | | | | |
| 15-05-2040-31 | Matrix Spike | Solid | GC 26 | 06/09/15 | 06/12/15 12:53 | 150609S20 | | | | |
| 15-05-2040-31 | Matrix Spike Duplicate | Solid | GC 26 | 06/09/15 | 06/12/15 13:37 | 150609S20 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | ND | 4.000 | 4.195 | 105 | 4.092 | 102 | 30-130 | 2 | 0-30 | |
| Bolstar | ND | 4.000 | 3.999 | 100 | 3.906 | 98 | 30-130 | 2 | 0-30 | |
| Chlorpyrifos | ND | 4.000 | 3.810 | 95 | 3.691 | 92 | 30-130 | 3 | 0-30 | |
| Coumaphos | ND | 4.000 | 4.344 | 109 | 4.281 | 107 | 30-130 | 1 | 0-30 | |
| Diazinon | ND | 4.000 | 3.873 | 97 | 3.883 | 97 | 30-130 | 0 | 0-30 | |
| Disulfoton | ND | 4.000 | 4.393 | 110 | 4.327 | 108 | 30-130 | 2 | 0-30 | |
| Ethoprop | ND | 4.000 | 4.096 | 102 | 4.008 | 100 | 30-130 | 2 | 0-30 | |
| Fensulfothion | ND | 4.000 | 4.015 | 100 | 4.000 | 100 | 30-130 | 0 | 0-30 | |
| Fenthion | ND | 4.000 | 3.881 | 97 | 3.760 | 94 | 30-130 | 3 | 0-30 | |
| Merphos | ND | 4.000 | 3.825 | 96 | 3.767 | 94 | 30-130 | 2 | 0-30 | |
| Methyl Parathion | ND | 4.000 | 3.784 | 95 | 3.655 | 91 | 30-130 | 3 | 0-30 | |
| Phorate | ND | 4.000 | 4.678 | 117 | 4.463 | 112 | 30-130 | 5 | 0-30 | |
| Ronnel | ND | 4.000 | 3.702 | 93 | 3.594 | 90 | 30-130 | 3 | 0-30 | |
| Stirophos | ND | 4.000 | 3.328 | 83 | 3.267 | 82 | 30-130 | 2 | 0-30 | |
| Tokuthion | ND | 4.000 | 3.249 | 81 | 3.139 | 78 | 30-130 | 3 | 0-30 | |
| Trichloronate | ND | 4.000 | 3.914 | 98 | 3.789 | 95 | 30-130 | 3 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|--------|------------|---------------|----------------|---------------------|
| 15-06-0559-4 | Sample | Solid | GC 40 | 06/10/15 | 06/11/15 21:39 | 150610S09 |
| 15-06-0559-4 | Matrix Spike | Solid | GC 40 | 06/10/15 | 06/11/15 19:44 | 150610S09 |
| 15-06-0559-4 | Matrix Spike Duplicate | Solid | GC 40 | 06/10/15 | 06/11/15 20:07 | 150610S09 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| 2,4-D | ND | 400.0 | 259.5 | 65 | 263.7 | 66 | 30-130 | 2 | 0-30 | |
| 2,4,5-T | ND | 40.00 | 26.20 | 66 | 26.50 | 66 | 30-130 | 1 | 0-30 | |
| 2,4-DB | ND | 400.0 | 308.8 | 77 | 284.7 | 71 | 30-130 | 8 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-06-0562-3 | Sample | Solid | GC/MS TT | 06/11/15 | 06/11/15 18:16 | 150611S02 | | | | |
| 15-06-0562-3 | Matrix Spike | Solid | GC/MS TT | 06/11/15 | 06/11/15 18:35 | 150611S02 | | | | |
| 15-06-0562-3 | Matrix Spike Duplicate | Solid | GC/MS TT | 06/11/15 | 06/12/15 12:07 | 150611S02 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | ND | 10.00 | 8.524 | 85 | 8.573 | 86 | 34-148 | 1 | 0-20 | |
| Acenaphthylene | ND | 10.00 | 8.345 | 83 | 8.398 | 84 | 53-120 | 1 | 0-20 | |
| Butyl Benzyl Phthalate | ND | 10.00 | 9.267 | 93 | 9.369 | 94 | 15-189 | 1 | 0-20 | |
| 4-Chloro-3-Methylphenol | ND | 10.00 | 8.396 | 84 | 9.066 | 91 | 32-120 | 8 | 0-20 | |
| 2-Chlorophenol | ND | 10.00 | 8.130 | 81 | 8.393 | 84 | 53-120 | 3 | 0-20 | |
| 1,4-Dichlorobenzene | ND | 10.00 | 6.024 | 60 | 6.091 | 61 | 43-120 | 1 | 0-26 | |
| Dimethyl Phthalate | ND | 10.00 | 7.777 | 78 | 8.138 | 81 | 44-122 | 5 | 0-20 | |
| 2,4-Dinitrotoluene | ND | 10.00 | 8.156 | 82 | 8.663 | 87 | 28-120 | 6 | 0-20 | |
| Fluorene | ND | 10.00 | 8.435 | 84 | 8.771 | 88 | 12-186 | 4 | 0-20 | |
| N-Nitroso-di-n-propylamine | ND | 10.00 | 6.369 | 64 | 6.891 | 69 | 38-140 | 8 | 0-20 | |
| Naphthalene | ND | 10.00 | 7.922 | 79 | 7.953 | 80 | 20-140 | 0 | 0-20 | |
| 4-Nitrophenol | ND | 10.00 | 8.077 | 81 | 8.997 | 90 | 14-128 | 11 | 0-59 | |
| Pentachlorophenol | ND | 10.00 | 7.978 | 80 | 8.826 | 88 | 10-124 | 10 | 0-20 | |
| Phenol | ND | 10.00 | 7.455 | 75 | 8.017 | 80 | 22-124 | 7 | 0-20 | |
| Pyrene | ND | 10.00 | 8.828 | 88 | 8.932 | 89 | 31-169 | 1 | 0-20 | |
| 1,2,4-Trichlorobenzene | ND | 10.00 | 7.838 | 78 | 7.923 | 79 | 56-120 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-06-1033-2 | Sample | Aqueous | GC/MS WW | 06/13/15 | 06/13/15 18:26 | 150613S016 | | | | |
| 15-06-1033-2 | Matrix Spike | Aqueous | GC/MS WW | 06/13/15 | 06/13/15 13:45 | 150613S016 | | | | |
| 15-06-1033-2 | Matrix Spike Duplicate | Aqueous | GC/MS WW | 06/13/15 | 06/13/15 14:17 | 150613S016 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 50.49 | 101 | 48.46 | 97 | 74-122 | 4 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 54.43 | 109 | 49.19 | 98 | 60-144 | 10 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 50.85 | 102 | 49.80 | 100 | 73-120 | 2 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 49.67 | 99 | 50.90 | 102 | 80-122 | 2 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 52.20 | 104 | 48.90 | 98 | 70-120 | 7 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 51.93 | 104 | 49.63 | 99 | 64-142 | 5 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 66.34 | 133 | 59.69 | 119 | 52-136 | 11 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 59.36 | 119 | 56.12 | 112 | 77-125 | 6 | 0-24 | |
| Toluene | ND | 50.00 | 54.80 | 110 | 51.31 | 103 | 72-126 | 7 | 0-23 | |
| Trichloroethene | 1.511 | 50.00 | 56.13 | 109 | 53.23 | 103 | 74-128 | 5 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 47.07 | 94 | 42.08 | 84 | 67-133 | 11 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 118.9 | 119 | 110.5 | 111 | 63-129 | 7 | 0-25 | |
| o-Xylene | ND | 50.00 | 57.91 | 116 | 55.08 | 110 | 62-128 | 5 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 62.38 | 125 | 61.11 | 122 | 68-134 | 2 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-06-0573-1 | Sample | Solid | GC/MS Q | 06/08/15 | 06/09/15 02:59 | 150608S026 | | | | |
| 15-06-0573-1 | Matrix Spike | Solid | GC/MS Q | 06/08/15 | 06/09/15 08:18 | 150608S026 | | | | |
| 15-06-0573-1 | Matrix Spike Duplicate | Solid | GC/MS Q | 06/08/15 | 06/09/15 08:44 | 150608S026 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 34.19 | 68 | 34.46 | 69 | 61-127 | 1 | 0-20 | |
| Carbon Tetrachloride | ND | 50.00 | 37.69 | 75 | 38.02 | 76 | 51-135 | 1 | 0-29 | |
| Chlorobenzene | ND | 50.00 | 29.06 | 58 | 29.64 | 59 | 57-123 | 2 | 0-20 | |
| 1,2-Dibromoethane | ND | 50.00 | 29.81 | 60 | 30.02 | 60 | 64-124 | 1 | 0-20 | 3 |
| 1,2-Dichlorobenzene | ND | 50.00 | 20.75 | 42 | 21.34 | 43 | 35-131 | 3 | 0-25 | |
| 1,2-Dichloroethane | ND | 50.00 | 28.37 | 57 | 28.63 | 57 | 80-120 | 1 | 0-20 | 3 |
| 1,1-Dichloroethene | ND | 50.00 | 35.68 | 71 | 36.06 | 72 | 47-143 | 1 | 0-25 | |
| Ethylbenzene | ND | 50.00 | 33.59 | 67 | 33.84 | 68 | 57-129 | 1 | 0-22 | |
| Toluene | ND | 50.00 | 34.11 | 68 | 33.84 | 68 | 63-123 | 1 | 0-20 | |
| Trichloroethene | ND | 50.00 | 37.73 | 75 | 37.43 | 75 | 44-158 | 1 | 0-20 | |
| Vinyl Chloride | ND | 50.00 | 40.85 | 82 | 41.57 | 83 | 49-139 | 2 | 0-47 | |
| p/m-Xylene | ND | 100.0 | 65.41 | 65 | 65.20 | 65 | 70-130 | 0 | 0-30 | 3 |
| o-Xylene | ND | 50.00 | 30.29 | 61 | 30.61 | 61 | 70-130 | 1 | 0-30 | 3 |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 30.36 | 61 | 30.62 | 61 | 57-123 | 1 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HSM310 | Sample | Solid | GC/MS BB | 06/07/15 | 06/09/15 01:29 | 150608S015 | | | | |
| HSM310 | Matrix Spike | Solid | GC/MS BB | 06/07/15 | 06/09/15 01:58 | 150608S015 | | | | |
| HSM310 | Matrix Spike Duplicate | Solid | GC/MS BB | 06/07/15 | 06/09/15 02:26 | 150608S015 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 34.19 | 68 | 29.52 | 59 | 61-127 | 15 | 0-20 | 3 |
| Carbon Tetrachloride | ND | 50.00 | 33.52 | 67 | 27.82 | 56 | 51-135 | 19 | 0-29 | |
| Chlorobenzene | ND | 50.00 | 29.91 | 60 | 25.61 | 51 | 57-123 | 15 | 0-20 | 3 |
| 1,2-Dibromoethane | ND | 50.00 | 32.34 | 65 | 29.26 | 59 | 64-124 | 10 | 0-20 | 3 |
| 1,2-Dichlorobenzene | ND | 50.00 | 23.31 | 47 | 20.32 | 41 | 35-131 | 14 | 0-25 | |
| 1,2-Dichloroethane | ND | 50.00 | 35.96 | 72 | 32.75 | 66 | 80-120 | 9 | 0-20 | 3 |
| 1,1-Dichloroethene | ND | 50.00 | 36.85 | 74 | 31.66 | 63 | 47-143 | 15 | 0-25 | |
| Ethylbenzene | ND | 50.00 | 33.64 | 67 | 28.28 | 57 | 57-129 | 17 | 0-22 | |
| Toluene | ND | 50.00 | 33.21 | 66 | 28.24 | 56 | 63-123 | 16 | 0-20 | 3 |
| Trichloroethene | ND | 50.00 | 37.33 | 75 | 32.56 | 65 | 44-158 | 14 | 0-20 | |
| Vinyl Chloride | ND | 50.00 | 42.51 | 85 | 37.14 | 74 | 49-139 | 13 | 0-47 | |
| p/m-Xylene | ND | 100.0 | 65.93 | 66 | 55.31 | 55 | 70-130 | 18 | 0-30 | 3 |
| o-Xylene | ND | 50.00 | 30.10 | 60 | 25.63 | 51 | 70-130 | 16 | 0-30 | 3 |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 34.27 | 69 | 32.08 | 64 | 57-123 | 7 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|--------------|-------------|------------|----------------|----------------|-----------------------|
| 15-06-0965-1 | Sample | Solid | ICP/MS 03 | 06/12/15 00:00 | 06/12/15 17:47 | 150612S01 |
| 15-06-0965-1 | PDS | Solid | ICP/MS 03 | 06/12/15 00:00 | 06/12/15 17:39 | 150612S01 |
| Parameter | Sample Conc. | Spike Added | PDS Conc. | PDS %Rec. | %Rec. CL | Qualifiers |
| Antimony | ND | 25.00 | 26.50 | 106 | 75-125 | |
| Arsenic | 3.329 | 25.00 | 29.85 | 106 | 75-125 | |
| Barium | 192.6 | 25.00 | 213.2 | 4X | 75-125 | Q |
| Beryllium | ND | 25.00 | 27.16 | 109 | 75-125 | |
| Cadmium | ND | 25.00 | 26.43 | 106 | 75-125 | |
| Chromium | 37.93 | 25.00 | 61.81 | 96 | 75-125 | |
| Copper | 29.39 | 25.00 | 53.68 | 97 | 75-125 | |
| Lead | 9.600 | 25.00 | 35.87 | 105 | 75-125 | |
| Nickel | 25.79 | 25.00 | 48.91 | 92 | 75-125 | |
| Selenium | ND | 25.00 | 27.85 | 111 | 75-125 | |
| Silver | ND | 12.50 | 12.32 | 99 | 75-125 | |
| Thallium | ND | 25.00 | 25.49 | 102 | 75-125 | |
| Zinc | 96.60 | 25.00 | 123.7 | 108 | 75-125 | |
| Aluminum | 19620 | 25.00 | 19940 | 4X | 75-125 | Q |
| Iron | 30610 | 25.00 | 30260 | 4X | 75-125 | Q |
| Manganese | 312.0 | 25.00 | 328.0 | 4X | 75-125 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|--------------|-------------|------------|----------------|----------------|-----------------------|
| 15-06-0423-2 | Sample | Aqueous | ICP/MS 04 | 06/08/15 00:00 | 06/09/15 17:26 | 150608SA2A |
| 15-06-0423-2 | PDS | Aqueous | ICP/MS 04 | 06/08/15 00:00 | 06/09/15 17:14 | 150608SA2A |
| Parameter | Sample Conc. | Spike Added | PDS Conc. | PDS %Rec. | %Rec. CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.09069 | 91 | 75-125 | |
| Arsenic | ND | 0.1000 | 0.08723 | 87 | 75-125 | |
| Barium | 0.05909 | 0.1000 | 0.1523 | 93 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.09423 | 94 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.08697 | 87 | 75-125 | |
| Chromium | 0.001241 | 0.1000 | 0.1023 | 101 | 75-125 | |
| Copper | 0.001283 | 0.1000 | 0.09096 | 90 | 75-125 | |
| Lead | ND | 0.1000 | 0.1013 | 101 | 75-125 | |
| Nickel | 0.006589 | 0.1000 | 0.09511 | 89 | 75-125 | |
| Selenium | 0.001400 | 0.1000 | 0.08100 | 80 | 75-125 | |
| Silver | ND | 0.05000 | 0.04173 | 83 | 75-125 | |
| Thallium | ND | 0.1000 | 0.09854 | 99 | 75-125 | |
| Zinc | ND | 0.1000 | 0.08036 | 80 | 75-125 | |
| Aluminum | ND | 0.1000 | 0.1023 | 102 | 75-125 | |
| Iron | 0.08723 | 5.100 | 4.732 | 91 | 75-125 | |
| Manganese | 0.03515 | 0.1000 | 0.1294 | 94 | 75-125 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: ASTM D-2216 (M)

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|---------------------|------------------|-----------------------|-----------------------|------------------------|
| HSM310 | Sample | Solid | N/A | 06/08/15 00:00 | 06/08/15 20:00 | F0608MOID6 |
| HSM310 | Sample Duplicate | Solid | N/A | 06/08/15 00:00 | 06/08/15 20:00 | F0608MOID6 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Moisture | | 71.30 | 72.00 | 1 | 0-10 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 9045D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|---------------------|------------------|-----------------------|-----------------------|------------------------|
| HSM310 | Sample | Solid | PH 4 | 06/06/15 00:00 | 06/06/15 14:23 | F0606PHD2 |
| HSM310 | Sample Duplicate | Solid | PH 4 | 06/06/15 00:00 | 06/06/15 14:23 | F0606PHD2 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| pH | | 7.710 | 7.730 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---|-------------------------|---------------------|------------------|---------------|-----------------------|------------------------|
| EB01 | Sample | Aqueous | PH1/BUR03 | N/A | 06/11/15 12:25 | F0611ALKD1 |
| EB01 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 06/11/15 12:25 | F0611ALKD1 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Alkalinity, Total (as CaCO ₃) | | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|-------------------------------------|-------------------------|---------------------|------------------|---------------|-----------------------|------------------------|
| EB01 | Sample | Aqueous | PH1/BUR03 | N/A | 06/11/15 12:25 | F0611HCOD1 |
| EB01 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 06/11/15 12:25 | F0611HCOD1 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Bicarbonate (as CaCO ₃) | | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|----------------|------------------|---------------|-----------------------|------------------------|
| EB01 | Sample | Aqueous | PH1/BUR03 | N/A | 06/11/15 12:25 | F0611CO3D1 |
| EB01 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 06/11/15 12:25 | F0611CO3D1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 2320B M

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|--------|------------|---------------|----------------|------------------------|
| 15-06-0770-3 | Sample | Solid | PH1/BUR03 | N/A | 06/15/15 16:35 | F0615ALKD2 |
| 15-06-0770-3 | Sample Duplicate | Solid | PH1/BUR03 | N/A | 06/15/15 16:35 | F0615ALKD2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|---|---------------------|------------------|------------|---------------|-------------------|
| Alkalinity, Total (as CaCO ₃) | 200.0 | 195.0 | 3 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 2320B M

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|--------|------------|----------------|----------------|------------------------|
| 15-06-0770-3 | Sample | Solid | PH1/BUR03 | 06/15/15 00:00 | 06/15/15 16:35 | F0615HCOD2 |
| 15-06-0770-3 | Sample Duplicate | Solid | PH1/BUR03 | 06/15/15 00:00 | 06/15/15 16:35 | F0615HCOD2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Bicarbonate (as CaCO ₃) | 200.0 | 195.0 | 3 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 2320B M

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|--------|------------|----------------|----------------|------------------------|
| 15-06-0770-3 | Sample | Solid | PH1/BUR03 | 06/15/15 00:00 | 06/15/15 16:35 | F0615CO3D2 |
| 15-06-0770-3 | Sample Duplicate | Solid | PH1/BUR03 | 06/15/15 00:00 | 06/15/15 16:35 | F0615CO3D2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-06-0863-1 | Sample | Aqueous | SC 2 | 06/11/15 00:00 | 06/11/15 18:00 | F0611TDSD1 |
| 15-06-0863-1 | Sample Duplicate | Aqueous | SC 2 | 06/11/15 00:00 | 06/11/15 18:00 | F0611TDSD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Dissolved | 7900 | 7830 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 2540 C (M)

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|---------------------|------------------|-----------------------|-----------------------|------------------------|
| HSM310 | Sample | Solid | SC 2 | 06/16/15 00:00 | 06/16/15 19:00 | F0616TDSD1 |
| HSM310 | Sample Duplicate | Solid | SC 2 | 06/16/15 00:00 | 06/16/15 19:00 | F0616TDSD1 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Solids, Total Dissolved | | 1395 | 1429 | 2 | 0-10 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-06-0467-1 | Sample | Aqueous | PH 1 | N/A | 06/06/15 12:30 | F0606PHD1 |
| 15-06-0467-1 | Sample Duplicate | Aqueous | PH 1 | N/A | 06/06/15 12:30 | F0606PHD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|------------------|---------------------|------------------|------------|---------------|-------------------|
| pH | 7.480 | 7.460 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|--------------------|------------------------|------------------|-----------------|-------------------|
| 099-12-906-5827 | LCS | Aqueous | IC 7 | N/A | 06/06/15 10:42 | 150606L01 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Fluoride | | 2.500 | 2.475 | 99 | 90-110 | |
| Chloride | | 50.00 | 47.67 | 95 | 90-110 | |
| Bromide | | 5.000 | 4.707 | 94 | 90-110 | |
| Nitrate (as N) | | 5.000 | 4.810 | 96 | 90-110 | |
| Sulfate | | 50.00 | 48.67 | 97 | 90-110 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|--------|------------|---------------|----------------|------------------|
| 099-12-922-582 | LCS | Solid | IC 10 | 06/09/15 | 06/09/15 11:56 | 150609L01P |

| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
|----------------|-------------|-----------------|-----------|----------|------------|
| Fluoride | 25.00 | 23.64 | 95 | 90-110 | |
| Chloride | 500.0 | 513.1 | 103 | 90-110 | |
| Bromide | 50.00 | 50.56 | 101 | 90-110 | |
| Nitrate (as N) | 50.00 | 51.62 | 103 | 90-110 | |
| Sulfate | 500.0 | 516.2 | 103 | 90-110 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-739-107 | LCS | Aqueous | ACA 1 | N/A | 06/13/15 14:38 | 150613L01 | | | |
| 099-12-739-107 | LCSD | Aqueous | ACA 1 | N/A | 06/13/15 14:38 | 150613L01 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Phosphorus, Total | 0.2000 | 0.1957 | 98 | 0.1980 | 99 | 90-110 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: EPA 9060A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-06-013-1275 | LCS | Solid | TOC 5 | 06/15/15 | 06/15/15 16:48 | F0615TOCL1 | | | |
| 099-06-013-1275 | LCSD | Solid | TOC 5 | 06/15/15 | 06/15/15 16:48 | F0615TOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 6000 | 5946 | 99 | 5901 | 98 | 80-120 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-981-104 | LCS | Aqueous | PH1/BUR03 | N/A | 06/11/15 12:25 | F0611ALKB1 | | | |
| 099-15-981-104 | LCSD | Aqueous | PH1/BUR03 | N/A | 06/11/15 12:25 | F0611ALKB1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 10.00 | 10.00 | 100 | 10.00 | 100 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4619 | LCS | Aqueous | SC 2 | 06/11/15 | 06/11/15 18:00 | F0611TDSL1 | | | |
| 099-12-180-4619 | LCSD | Aqueous | SC 2 | 06/11/15 | 06/11/15 18:00 | F0611TDSL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 90.00 | 90 | 95.00 | 95 | 80-120 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 4500 P B/E (M)

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------------|------------|---------------|-------------------|
| 099-05-001-5419 | LCS | Solid | UV 7 | 06/10/15 | 06/11/15 16:26 | F0611TPL2 | | | |
| 099-05-001-5419 | LCSD | Solid | UV 7 | 06/10/15 | 06/11/15 16:26 | F0611TPL2 | | | |
| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Phosphorus, Total | 2.000 | 1.940 | 97 | 1.945 | 97 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-097-5675 | LCS | Aqueous | TOC 8 | 06/19/15 | 06/19/15 22:52 | F0619TOCL1 | | | |
| 099-05-097-5675 | LCSD | Aqueous | TOC 8 | 06/19/15 | 06/19/15 22:52 | F0619TOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 10.00 | 8.700 | 87 | 9.320 | 93 | 80-120 | 7 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
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San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 097-01-002-21301 | LCS | Solid | ICP 7300 | 06/11/15 | 06/24/15 17:49 | 150611L09A |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Calcium | | 25.00 | 23.84 | 95 | 80-120 | |
| Magnesium | | 25.00 | 23.97 | 96 | 80-120 | |
| Potassium | | 250.0 | 222.2 | 89 | 80-120 | |
| Sodium | | 250.0 | 222.7 | 89 | 80-120 | |
| Strontium | | 25.00 | 24.92 | 100 | 80-120 | |
| Silicon | | 25.00 | 20.35 | 81 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

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San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-15-683-1299 | LCS | Aqueous | ICP 7300 | 06/15/15 | 06/16/15 11:29 | 150615LA4F |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Calcium | | 0.5000 | 0.5004 | 100 | 80-120 | |
| Magnesium | | 0.5000 | 0.5464 | 109 | 80-120 | |
| Potassium | | 5.000 | 5.072 | 101 | 80-120 | |
| Sodium | | 5.000 | 5.025 | 101 | 80-120 | |
| Strontium | | 0.5000 | 0.5134 | 103 | 80-120 | |
| Silicon | | 0.5000 | 0.4580 | 92 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3050B
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-621-955 | LCS | Solid | ICP/MS 03 | 06/12/15 | 06/12/15 17:29 | 150612L01 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 25.00 | 25.38 | 102 | 80-120 | 73-127 | |
| Arsenic | 25.00 | 27.13 | 109 | 80-120 | 73-127 | |
| Barium | 25.00 | 24.89 | 100 | 80-120 | 73-127 | |
| Beryllium | 25.00 | 26.54 | 106 | 80-120 | 73-127 | |
| Cadmium | 25.00 | 25.76 | 103 | 80-120 | 73-127 | |
| Chromium | 25.00 | 27.45 | 110 | 80-120 | 73-127 | |
| Copper | 25.00 | 27.35 | 109 | 80-120 | 73-127 | |
| Lead | 25.00 | 25.99 | 104 | 80-120 | 73-127 | |
| Nickel | 25.00 | 26.73 | 107 | 80-120 | 73-127 | |
| Selenium | 25.00 | 26.60 | 106 | 80-120 | 73-127 | |
| Silver | 12.50 | 12.30 | 98 | 80-120 | 73-127 | |
| Thallium | 25.00 | 24.90 | 100 | 80-120 | 73-127 | |
| Zinc | 25.00 | 26.62 | 106 | 80-120 | 73-127 | |
| Aluminum | 25.00 | 21.91 | 88 | 80-120 | 73-127 | |
| Iron | 25.00 | 23.45 | 94 | 80-120 | 73-127 | |
| Manganese | 25.00 | 27.37 | 109 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-842 | LCS | Aqueous | ICP/MS 04 | 06/08/15 | 06/09/15 17:03 | 150608LA2F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.1040 | 104 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.1063 | 106 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.1029 | 103 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.1028 | 103 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.1028 | 103 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.1071 | 107 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.1069 | 107 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.1030 | 103 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.1039 | 104 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.1021 | 102 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.04912 | 98 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.1001 | 100 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.1084 | 108 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.1064 | 106 | 80-120 | 73-127 | |
| Iron | 5.100 | 5.346 | 105 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.1049 | 105 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| 099-15-763-567 | LCS | Aqueous | Mercury 04 | 06/08/15 | 06/08/15 13:56 | 150608L01F |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.01000 | 0.009903 | 99 | 85-121 | |


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Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 7471A Total
Method: EPA 7471A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------|-------------------|-----------------|-----------------------|------------------|
| 099-16-272-1355 | LCS | Solid | Mercury 05 | 06/17/15 | 06/17/15 20:05 | 150617L03 |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.8350 | 1.004 | 120 | 85-121 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-12-537-2135 | LCS | Solid | GC 41 | 06/10/15 | 06/11/15 20:20 | 150610L17 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Aldrin | 25.00 | 20.45 | 82 | 50-135 | 36-149 | |
| Alpha-BHC | 25.00 | 20.79 | 83 | 50-135 | 36-149 | |
| Beta-BHC | 25.00 | 27.23 | 109 | 50-135 | 36-149 | |
| 4,4'-DDD | 25.00 | 22.65 | 91 | 50-135 | 36-149 | |
| 4,4'-DDE | 25.00 | 23.00 | 92 | 50-135 | 36-149 | |
| 4,4'-DDT | 25.00 | 22.73 | 91 | 50-135 | 36-149 | |
| Delta-BHC | 25.00 | 22.29 | 89 | 50-135 | 36-149 | |
| Dieldrin | 25.00 | 24.68 | 99 | 50-135 | 36-149 | |
| Endosulfan I | 25.00 | 22.91 | 92 | 50-135 | 36-149 | |
| Endosulfan II | 25.00 | 23.81 | 95 | 50-135 | 36-149 | |
| Endosulfan Sulfate | 25.00 | 22.64 | 91 | 50-135 | 36-149 | |
| Endrin | 25.00 | 23.83 | 95 | 50-135 | 36-149 | |
| Endrin Aldehyde | 25.00 | 19.96 | 80 | 50-135 | 36-149 | |
| Gamma-BHC | 25.00 | 21.53 | 86 | 50-135 | 36-149 | |
| Heptachlor | 25.00 | 23.29 | 93 | 50-135 | 36-149 | |
| Heptachlor Epoxide | 25.00 | 23.29 | 93 | 50-135 | 36-149 | |
| Methoxychlor | 25.00 | 23.76 | 95 | 50-135 | 36-149 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-529-813 | LCS | Aqueous | | GC 44 | 06/08/15 | 06/10/15 13:57 | 150608L02 | | | |
| 099-12-529-813 | LCSD | Aqueous | | GC 44 | 06/08/15 | 06/10/15 14:12 | 150608L02 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Alpha-BHC | 0.5000 | 0.4410 | 88 | 0.4620 | 92 | 50-135 | 36-149 | 5 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.4504 | 90 | 0.4709 | 94 | 50-135 | 36-149 | 4 | 0-25 | |
| Beta-BHC | 0.5000 | 0.4414 | 88 | 0.4599 | 92 | 50-135 | 36-149 | 4 | 0-25 | |
| Heptachlor | 0.5000 | 0.4442 | 89 | 0.4606 | 92 | 50-135 | 36-149 | 4 | 0-25 | |
| Delta-BHC | 0.5000 | 0.4371 | 87 | 0.4555 | 91 | 50-135 | 36-149 | 4 | 0-25 | |
| Aldrin | 0.5000 | 0.4354 | 87 | 0.4617 | 92 | 50-135 | 36-149 | 6 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.4343 | 87 | 0.4542 | 91 | 50-135 | 36-149 | 4 | 0-25 | |
| Endosulfan I | 0.5000 | 0.4494 | 90 | 0.4778 | 96 | 50-135 | 36-149 | 6 | 0-25 | |
| Dieldrin | 0.5000 | 0.4510 | 90 | 0.4718 | 94 | 50-135 | 36-149 | 4 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.4065 | 81 | 0.4165 | 83 | 50-135 | 36-149 | 2 | 0-25 | |
| Endrin | 0.5000 | 0.4627 | 93 | 0.4719 | 94 | 50-135 | 36-149 | 2 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.3976 | 80 | 0.4285 | 86 | 50-135 | 36-149 | 7 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.4460 | 89 | 0.4669 | 93 | 50-135 | 36-149 | 5 | 0-25 | |
| Endosulfan II | 0.5000 | 0.4405 | 88 | 0.4605 | 92 | 50-135 | 36-149 | 4 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.3572 | 71 | 0.3586 | 72 | 50-135 | 36-149 | 0 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.4403 | 88 | 0.4498 | 90 | 50-135 | 36-149 | 2 | 0-25 | |
| Methoxychlor | 0.5000 | 0.3351 | 67 | 0.3408 | 68 | 50-135 | 36-149 | 2 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 099-12-535-3266 | LCS | Solid | GC 58 | 06/10/15 | 06/11/15 22:03 | 150610L18 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Aroclor-1016 | | 100.0 | 74.42 | 74 | 50-135 | |
| Aroclor-1260 | | 100.0 | 56.80 | 57 | 50-135 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-533-1050 | LCS | Aqueous | GC 31 | 06/08/15 | 06/09/15 18:42 | 150608L03 | | | |
| 099-12-533-1050 | LCSD | Aqueous | GC 31 | 06/08/15 | 06/09/15 19:01 | 150608L03 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Aroclor-1016 | 2.000 | 2.265 | 113 | 2.014 | 101 | 50-135 | 12 | 0-25 | |
| Aroclor-1260 | 2.000 | 1.990 | 100 | 1.716 | 86 | 50-135 | 15 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

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Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-963-97 | LCS | Aqueous | | GC 26 | 06/08/15 | 06/10/15 12:52 | 150608L04 | | | |
| 099-15-963-97 | LCSD | Aqueous | | GC 26 | 06/08/15 | 06/10/15 13:37 | 150608L04 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.04300 | 108 | 0.04589 | 115 | 30-130 | 13-147 | 7 | 0-30 | |
| Bolstar | 0.04000 | 0.03955 | 99 | 0.04185 | 105 | 30-130 | 13-147 | 6 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.03756 | 94 | 0.04046 | 101 | 30-130 | 13-147 | 7 | 0-30 | |
| Coumaphos | 0.04000 | 0.04345 | 109 | 0.04521 | 113 | 30-130 | 13-147 | 4 | 0-30 | |
| Diazinon | 0.04000 | 0.04000 | 100 | 0.04434 | 111 | 30-130 | 13-147 | 10 | 0-30 | |
| Disulfoton | 0.04000 | 0.04213 | 105 | 0.04673 | 117 | 30-130 | 13-147 | 10 | 0-30 | |
| Ethoprop | 0.04000 | 0.04083 | 102 | 0.04688 | 117 | 30-130 | 13-147 | 14 | 0-30 | |
| Fensulfothion | 0.04000 | 0.04101 | 103 | 0.04297 | 107 | 30-130 | 13-147 | 5 | 0-30 | |
| Fenthion | 0.04000 | 0.03866 | 97 | 0.04328 | 108 | 30-130 | 13-147 | 11 | 0-30 | |
| Merphos | 0.04000 | 0.03773 | 94 | 0.04108 | 103 | 30-130 | 13-147 | 9 | 0-30 | |
| Methyl Parathion | 0.04000 | 0.03595 | 90 | 0.04168 | 104 | 30-130 | 13-147 | 15 | 0-30 | |
| Phorate | 0.04000 | 0.04692 | 117 | 0.04556 | 114 | 30-130 | 13-147 | 3 | 0-30 | |
| Ronnel | 0.04000 | 0.03567 | 89 | 0.04195 | 105 | 30-130 | 13-147 | 16 | 0-30 | |
| Stirophos | 0.04000 | 0.03312 | 83 | 0.03577 | 89 | 30-130 | 13-147 | 8 | 0-30 | |
| Tokuthion | 0.04000 | 0.04239 | 106 | 0.04464 | 112 | 30-130 | 13-147 | 5 | 0-30 | |
| Trichloronate | 0.04000 | 0.03861 | 97 | 0.04175 | 104 | 30-130 | 13-147 | 8 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-973-188 | LCS | Solid | GC 26 | 06/09/15 | 06/12/15 12:08 | 150609L20 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Azinphos Methyl | 4.000 | 4.285 | 107 | 30-130 | 13-147 | |
| Bolstar | 4.000 | 4.166 | 104 | 30-130 | 13-147 | |
| Chlorpyrifos | 4.000 | 3.975 | 99 | 30-130 | 13-147 | |
| Coumaphos | 4.000 | 4.409 | 110 | 30-130 | 13-147 | |
| Diazinon | 4.000 | 4.067 | 102 | 30-130 | 13-147 | |
| Disulfoton | 4.000 | 4.556 | 114 | 30-130 | 13-147 | |
| Ethoprop | 4.000 | 4.265 | 107 | 30-130 | 13-147 | |
| Fensulfothion | 4.000 | 4.241 | 106 | 30-130 | 13-147 | |
| Fenthion | 4.000 | 4.046 | 101 | 30-130 | 13-147 | |
| Merphos | 4.000 | 4.057 | 101 | 30-130 | 13-147 | |
| Methyl Parathion | 4.000 | 3.836 | 96 | 30-130 | 13-147 | |
| Phorate | 4.000 | 5.086 | 127 | 30-130 | 13-147 | |
| Ronnel | 4.000 | 3.789 | 95 | 30-130 | 13-147 | |
| Stirophos | 4.000 | 3.444 | 86 | 30-130 | 13-147 | |
| Tokuthion | 4.000 | 4.172 | 104 | 30-130 | 13-147 | |
| Trichloronate | 4.000 | 4.033 | 101 | 30-130 | 13-147 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 095-01-033-1294 | LCS | Solid | GC 40 | 06/10/15 | 06/11/15 18:57 | 150610L09 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| 2,4-D | | 400.0 | 284.1 | 71 | 30-130 | |
| 2,4,5-T | | 40.00 | 28.70 | 72 | 30-130 | |
| 2,4-DB | | 400.0 | 291.7 | 73 | 30-130 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-034-647 | LCS | Aqueous | GC 40 | 06/08/15 | 06/12/15 03:48 | 150608L06 | | | |
| 095-01-034-647 | LCSD | Aqueous | GC 40 | 06/08/15 | 06/12/15 04:11 | 150608L06 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| 2,4-D | 20.00 | 20.57 | 103 | 16.46 | 82 | 30-130 | 22 | 0-30 | |
| 2,4,5-T | 2.000 | 2.010 | 100 | 1.665 | 83 | 30-130 | 19 | 0-30 | |
| 2,4-DB | 20.00 | 22.49 | 112 | 18.35 | 92 | 30-130 | 20 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-4051 | LCS | Aqueous | | GC/MS TT | 06/08/15 | 06/08/15 16:14 | 150608L01 | | | |
| 095-01-003-4051 | LCSD | Aqueous | | GC/MS TT | 06/08/15 | 06/08/15 16:33 | 150608L01 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 200.0 | 158.8 | 79 | 171.2 | 86 | 61-120 | 51-130 | 8 | 0-20 | |
| Acenaphthylene | 200.0 | 155.7 | 78 | 167.0 | 84 | 55-120 | 44-131 | 7 | 0-20 | |
| Butyl Benzyl Phthalate | 200.0 | 161.0 | 80 | 171.0 | 86 | 56-122 | 45-133 | 6 | 0-20 | |
| 4-Chloro-3-Methylphenol | 200.0 | 151.4 | 76 | 160.9 | 80 | 52-120 | 41-131 | 6 | 0-20 | |
| 2-Chlorophenol | 200.0 | 151.9 | 76 | 159.7 | 80 | 47-120 | 35-132 | 5 | 0-20 | |
| 1,4-Dichlorobenzene | 200.0 | 147.8 | 74 | 155.7 | 78 | 36-120 | 22-134 | 5 | 0-20 | |
| Dimethyl Phthalate | 200.0 | 154.9 | 77 | 163.9 | 82 | 60-120 | 50-130 | 6 | 0-20 | |
| 2,4-Dinitrotoluene | 200.0 | 153.9 | 77 | 164.7 | 82 | 61-121 | 51-131 | 7 | 0-20 | |
| Fluorene | 200.0 | 157.3 | 79 | 169.1 | 85 | 67-120 | 58-129 | 7 | 0-20 | |
| N-Nitroso-di-n-propylamine | 200.0 | 139.6 | 70 | 143.6 | 72 | 39-123 | 25-137 | 3 | 0-20 | |
| Naphthalene | 200.0 | 156.0 | 78 | 166.9 | 83 | 54-120 | 43-131 | 7 | 0-20 | |
| 4-Nitrophenol | 200.0 | 66.80 | 33 | 69.92 | 35 | 14-120 | 0-138 | 5 | 0-20 | |
| Pentachlorophenol | 200.0 | 157.2 | 79 | 165.5 | 83 | 31-127 | 15-143 | 5 | 0-20 | |
| Phenol | 200.0 | 72.85 | 36 | 74.14 | 37 | 17-120 | 0-137 | 2 | 0-20 | |
| Pyrene | 200.0 | 162.4 | 81 | 172.5 | 86 | 58-124 | 47-135 | 6 | 0-20 | |
| 1,2,4-Trichlorobenzene | 200.0 | 164.0 | 82 | 172.5 | 86 | 49-120 | 37-132 | 5 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 3545
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-12-549-3303 | LCS | Solid | GC/MS TT | 06/11/15 | 06/11/15 16:35 | 150611L02 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Acenaphthene | 10.00 | 8.342 | 83 | 51-123 | 39-135 | |
| Acenaphthylene | 10.00 | 8.197 | 82 | 52-120 | 41-131 | |
| Butyl Benzyl Phthalate | 10.00 | 9.211 | 92 | 43-139 | 27-155 | |
| 4-Chloro-3-Methylphenol | 10.00 | 8.287 | 83 | 55-121 | 44-132 | |
| 2-Chlorophenol | 10.00 | 8.255 | 83 | 58-124 | 47-135 | |
| 1,4-Dichlorobenzene | 10.00 | 7.641 | 76 | 42-132 | 27-147 | |
| Dimethyl Phthalate | 10.00 | 7.933 | 79 | 51-123 | 39-135 | |
| 2,4-Dinitrotoluene | 10.00 | 8.347 | 83 | 51-129 | 38-142 | |
| Fluorene | 10.00 | 8.266 | 83 | 54-126 | 42-138 | |
| N-Nitroso-di-n-propylamine | 10.00 | 6.744 | 67 | 40-136 | 24-152 | |
| Naphthalene | 10.00 | 8.350 | 84 | 32-146 | 13-165 | |
| 4-Nitrophenol | 10.00 | 7.644 | 76 | 24-126 | 7-143 | |
| Pentachlorophenol | 10.00 | 8.449 | 84 | 23-131 | 5-149 | |
| Phenol | 10.00 | 7.455 | 75 | 40-130 | 25-145 | |
| Pyrene | 10.00 | 8.656 | 87 | 47-143 | 31-159 | |
| 1,2,4-Trichlorobenzene | 10.00 | 8.784 | 88 | 45-129 | 31-143 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-17386 | LCS | Aqueous | GC/MS WW | 06/13/15 | 06/13/15 13:06 | 150613L015 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 45.74 | 91 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 48.33 | 97 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 47.32 | 95 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 49.00 | 98 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 48.18 | 96 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 50.72 | 101 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 58.69 | 117 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 54.35 | 109 | 80-123 | 73-130 | |
| Toluene | 50.00 | 49.61 | 99 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 49.95 | 100 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 40.85 | 82 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 110.6 | 111 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 53.42 | 107 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 59.59 | 119 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-12-796-9818 | LCS | Solid | GC/MS Q | 06/08/15 | 06/09/15 01:13 | 150608L054 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 44.86 | 90 | 78-120 | 71-127 | |
| Carbon Tetrachloride | 50.00 | 41.59 | 83 | 49-139 | 34-154 | |
| Chlorobenzene | 50.00 | 43.99 | 88 | 79-120 | 72-127 | |
| 1,2-Dibromoethane | 50.00 | 49.41 | 99 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 44.19 | 88 | 75-120 | 68-128 | |
| 1,2-Dichloroethane | 50.00 | 43.77 | 88 | 80-120 | 73-127 | |
| 1,1-Dichloroethene | 50.00 | 39.78 | 80 | 74-122 | 66-130 | |
| Ethylbenzene | 50.00 | 45.21 | 90 | 76-120 | 69-127 | |
| Toluene | 50.00 | 44.54 | 89 | 77-120 | 70-127 | |
| Trichloroethene | 50.00 | 46.24 | 92 | 80-120 | 73-127 | |
| Vinyl Chloride | 50.00 | 46.42 | 93 | 68-122 | 59-131 | |
| p/m-Xylene | 100.0 | 89.56 | 90 | 75-125 | 67-133 | |
| o-Xylene | 50.00 | 43.84 | 88 | 75-125 | 67-133 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 45.12 | 90 | 77-120 | 70-127 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/06/15
Work Order: 15-06-0567
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-12-796-9792 | LCS | Solid | GC/MS BB | 06/08/15 | 06/08/15 23:35 | 150608L032 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 47.92 | 96 | 78-120 | 71-127 | |
| Carbon Tetrachloride | 50.00 | 48.50 | 97 | 49-139 | 34-154 | |
| Chlorobenzene | 50.00 | 46.87 | 94 | 79-120 | 72-127 | |
| 1,2-Dibromoethane | 50.00 | 52.05 | 104 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 47.61 | 95 | 75-120 | 68-128 | |
| 1,2-Dichloroethane | 50.00 | 52.94 | 106 | 80-120 | 73-127 | |
| 1,1-Dichloroethene | 50.00 | 45.84 | 92 | 74-122 | 66-130 | |
| Ethylbenzene | 50.00 | 50.03 | 100 | 76-120 | 69-127 | |
| Toluene | 50.00 | 47.50 | 95 | 77-120 | 70-127 | |
| Trichloroethene | 50.00 | 49.38 | 99 | 80-120 | 73-127 | |
| Vinyl Chloride | 50.00 | 50.26 | 101 | 68-122 | 59-131 | |
| p/m-Xylene | 100.0 | 99.65 | 100 | 75-125 | 67-133 | |
| o-Xylene | 50.00 | 46.85 | 94 | 75-125 | 67-133 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 49.85 | 100 | 77-120 | 70-127 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

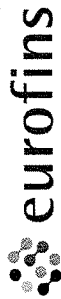
RPD: Relative Percent Difference. CL: Control Limits

Glossary of Terms and Qualifiers

Work Order: 15-06-0567

Page 1 of 1

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|---|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| CI | See case narrative. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |
| | Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time. |
| | A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations. |



Calscience

7440 Lincoln Way, Garden Grove, CA 92841-1427 • (714) 895-5494
For courier service / sample drop off information, contact us26_sales@eurofins.com or call us.

CHAIN OF CUSTODY RECORD

DATE: 06/05/15
PAGE: 1 OF 1

WO# / LAB USE ONLY
15-06-0567

| | | | | | |
|---|--|---|--|---|--|
| LABORATORY CLIENT: SWCA Environmental Consultants | | CLIENT PROJECT NAME / NUMBER: EAA 27122 | | P.O. NO.: | |
| ADDRESS: 6200 UTSA Blvd. Suite 102 | | PROJECT CONTACT: Philip Pearce | | SAMPLER(S) (PRINT): Philip Pearce/Jennifer Moreland | |
| CITY: San Antonio | | STATE: TX | | ZIP: 78249-1618 | |
| TEL: 210.877.2847 | | E-MAIL: P Pearce@swca.com | | | |

REQUESTED ANALYSES

| | | | | | | | | |
|---|-----------|---------------|------|--------|--------------|-------------|-----------|----------------|
| TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD"): | | LOG CODE: | | | | | | |
| <input type="checkbox"/> SAME DAY <input type="checkbox"/> 24 HR <input type="checkbox"/> 48 HR <input type="checkbox"/> 72 HR <input checked="" type="checkbox"/> 10 Days (standard) | | | | | | | | |
| <input type="checkbox"/> COELT EDF | | | | | | | | |
| GLOBAL ID: | | | | | | | | |
| SPECIAL INSTRUCTIONS: | | | | | | | | |
| Please analyze each sample for pH EPA 6020 Metals: Zn, Ti, Se, Sb, Pb, Ni, Cu, Cr, Cd, Be, Ba, As, Al, Ag, Fe, Mn | | | | | | | | |
| LAB USE ONLY | SAMPLE ID | SAMPLING DATE | TIME | MATRIX | NO. OF CONT. | Unpreserved | Preserved | Field Filtered |
| 1 | HSM310 | 6/5/2015 | 1003 | Solid | 2 | | | |
| 2 | HSM320 | 6/5/2015 | 1048 | Solid | 2 | | | |
| 3 | HSM330 | 6/5/2015 | 1103 | Solid | 2 | | | |
| 4 | HSM340 | 6/5/2015 | 1121 | Solid | 2 | | | |
| 5 | HSM350 | 6/5/2015 | 1224 | Solid | 2 | | | |
| 6 | HSM360 | 6/5/2015 | 1313 | Solid | 2 | | | |
| 7 | HSM370 | 6/5/2015 | 1351 | Solid | 2 | | | |
| 8 | FDHSM370 | 6/5/2015 | 1351 | Solid | 2 | | | |
| 9 | TB08 | 6/5/2015 | NA | NPW | 1 | | | |
| 10 | EB01 | 6/5/2015 | 1541 | NPW | 9 | | | |

| | | | | | |
|------------------------------|--|--------------------------------------|--|--------------|------------|
| Relinquished by: (Signature) | | Received by: (Signature/Affiliation) | | Date: 6/5/15 | Time: 1700 |
| Relinquished by: (Signature) | | Received by: (Signature/Affiliation) | | Date: 6/6/15 | Time: 0930 |
| Relinquished by: (Signature) | | Received by: (Signature/Affiliation) | | Date: | Time: |



0567

ORIGIN ID:SATA (210) 877-2847
SWCA INC

6200 UTSA BLVD STE 102

SAN ANTONIO, TX 782491618
UNITED STATES US

SHIP DATE: 05JUN15
ACTWGT: 45.0 LB MAN
CAD: /OFFC1601
DIMS: 23x14x14 IN

BILL SENDER

ORIGIN ID:SATA (210) 877-2847
SWCA INC

6200 UTSA BLVD STE 102

SAN ANTONIO, TX 782491618
UNITED STATES US

SHIP DATE: 05JUN15
ACTWGT: 58.0 LB MAN
CAD: /OFFC1601
DIMS: 23x14x14 IN

BILL SENDER

TO

CALSCIENCE
7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895-5494

REF:

YNU:

PO:

DEPT:

CALSCIENCE
7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895-5494

REF:

YNU:

PO:

DEPT:

FedEx
Express



J151215022301uv

FedEx
Express



AV100220912191J

2 of 2

MPS# 7807 7591 1587
0681

Mstr# 8062 7398 5461

0215

XO APVA

SATURDAY 12:00P
PRIORITY OVERNIGHT

DSR

92841

CA-US SNA

1 of 2

TRK# 8062 7398 5461
0215

MASTER

XO APVA

SATURDAY 12:00P
PRIORITY OVERNIGHT

DSR

92841

CA-US SNA

Return to Contents

SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 2

CLIENT: SWCA

DATE: 06/06/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 2.1 °C (w/ CF): 2.8 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEAL:

Cooler ☒ Present and Intact

☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 802

Sample(s) ☐ Present and Intact

☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 965

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE: 3

(Trip Blank Lot Number: 150508B)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2901 core ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☒ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☒ 202 CGJ

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ **Other Matrix** (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 965

s = H₂SO₄, u = ultra-pure, z₂na = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 802

SAMPLE RECEIPT CHECKLIST

COOLER 2 OF 2

CLIENT: SWCA

DATE: 06 / 06 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 3.0 °C (w/ CF): 2.7 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEAL:

Cooler ☒ Present and Intact

☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 802

Sample(s) ☐ Present and Intact

☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 802

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|--------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☐ VOA_h ☐ VOA_{na2} ☐ 100PJ ☐ 100PJ_{na2} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_z_{na} ☐ 250AGB ☐ 250CGB ☐ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s
☐ 500PB ☐ 1AGB ☐ 1AGB_{na2} ☐ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☐ _____ ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☒ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☒ 1ozCGJ

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ **Other Matrix** (_____) ☐ _____ ☐ _____

Container: **A** = Amber, **B** = Bottle, **C** = Clear, **E** = Envelope, **G** = Glass, **J** = Jar, **P** = Plastic, and **Z** = Ziploc/Resealable Bag

Preservative: **b** = buffered, **f** = filtered, **h** = HCl, **n** = HNO₃, **na** = NaOH, **na₂** = Na₂S₂O₃, **p** = H₃PO₄, Labeled/Checked by: 965

s = H₂SO₄, **u** = ultra-pure, **z**_{na} = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 802



Calscience



WORK ORDER NUMBER: 15-06-0884

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: SWCA Environmental Consultants

Client Project Name: EAA 27122

Attention: Philip Pearce
6200 UTSA Blvd.
Suite 102
San Antonio, TX 78249-1618

H. Burley FOR

Approved for release on 06/25/2015 by:
Don Burley
Project Manager

ResultLink ▶

Email your PM ▶



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Client Project Name: EAA 27122
Work Order Number: 15-06-0884

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Work Order Narrative

Work Order: 15-06-0884

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Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 06/11/15. They were assigned to Work Order 15-06-0884.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.



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Sample Summary

| | | | |
|---------|--------------------------------|-----------------------|----------------|
| Client: | SWCA Environmental Consultants | Work Order: | 15-06-0884 |
| | 6200 UTSA Blvd., Suite 102 | Project Name: | EAA 27122 |
| | San Antonio, TX 78249-1618 | PO Number: | |
| | | Date/Time Received: | 06/11/15 09:35 |
| | | Number of Containers: | 9 |

Attn: Philip Pearce

| Sample Identification | Lab Number | Collection Date and Time | Number of Containers | Matrix |
|-----------------------|--------------|--------------------------|----------------------|---------|
| EB02 | 15-06-0884-1 | 06/10/15 13:26 | 9 | Aqueous |


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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB02 | 15-06-0884-1-I | 06/10/15 13:26 | Aqueous | IC 10 | N/A | 06/11/15 15:40 | 150611L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | ND | 0.10 | 0.025 | 1.00 | |
| Chloride | ND | 1.0 | 0.12 | 1.00 | |
| Bromide | ND | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | ND | 0.10 | 0.025 | 1.00 | |
| Sulfate | ND | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|--------------|-----------------|-----|---------|-------|-----|----------------|-----------|
| Method Blank | 099-12-906-5828 | N/A | Aqueous | IC 10 | N/A | 06/11/15 10:08 | 150611L01 |
|--------------|-----------------|-----|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | ND | 0.10 | 0.025 | 1.00 | |
| Chloride | ND | 1.0 | 0.12 | 1.00 | |
| Bromide | ND | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | ND | 0.10 | 0.025 | 1.00 | |
| Sulfate | ND | 1.0 | 0.19 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB02 | 15-06-0884-1-F | 06/10/15 13:26 | Aqueous | ICP 7300 | 06/15/15 | 06/23/15 15:05 | 150615LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | ND | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 0.0131 | 0.100 | 0.00336 | 1.00 | B,J |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|-----------------|-----|---------|----------|----------|----------------|------------|
| Method Blank | 099-15-683-1299 | N/A | Aqueous | ICP 7300 | 06/15/15 | 06/16/15 11:27 | 150615LA4F |
|--------------|-----------------|-----|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | ND | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 0.0372 | 0.100 | 0.00336 | 1.00 | J |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 1 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB02 | 15-06-0884-1-F | 06/10/15 13:26 | Aqueous | ICP/MS 03 | 06/12/15 | 06/16/15 00:45 | 150612LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.000130 | 0.00100 | 0.0000986 | 1.00 | J |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00155 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.000425 | 0.00100 | 0.000132 | 1.00 | J |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00129 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.000253 | 0.00100 | 0.000139 | 1.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 2 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-844 | N/A | Aqueous | ICP/MS 03 | 06/12/15 | 06/15/15 18:56 | 150612LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB02 | 15-06-0884-1-F | 06/10/15 13:26 | Aqueous | Mercury 04 | 06/15/15 | 06/15/15 19:00 | 150615L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|-----|---------|------------|----------|----------------|------------|
| Method Blank | 099-15-763-571 | N/A | Aqueous | Mercury 04 | 06/15/15 | 06/15/15 18:08 | 150615L03F |
|--------------|----------------|-----|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

Page 1 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB02 | 15-06-0884-1-I | 06/10/15 13:26 | Aqueous | GC 41 | 06/15/15 | 06/17/15 18:36 | 150615L01B |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 96 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 103 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

Page 2 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-817 | N/A | Aqueous | GC 41 | 06/15/15 | 06/17/15 17:35 | 150615L01B |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 100 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 78 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB02 | 15-06-0884-1-I | 06/10/15 13:26 | Aqueous | GC 41 | 06/15/15 | 06/15/15 23:54 | 150615L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 98 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 111 | 50-135 | |

| Method Blank | 099-12-533-1052 | N/A | Aqueous | GC 41 | 06/15/15 | 06/15/15 20:52 | 150615L01 |
|--------------|-----------------|-----|---------|-------|----------|----------------|-----------|
|--------------|-----------------|-----|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 94 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 1 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB02 | 15-06-0884-1-I | 06/10/15 13:26 | Aqueous | GC 26 | 06/12/15 | 06/16/15 22:46 | 150612L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 116 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 2 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-99 | N/A | Aqueous | GC 26 | 06/12/15 | 06/16/15 20:33 | 150612L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 116 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB02 | 15-06-0884-1-I | 06/10/15 13:26 | Aqueous | GC 40 | 06/12/15 | 06/16/15 19:56 | 150612L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 65 | 0-123 | |

| Method Blank | 095-01-034-648 | N/A | Aqueous | GC 40 | 06/12/15 | 06/16/15 18:23 | 150612L12 |
|--------------|----------------|-----|---------|-------|----------|----------------|-----------|
|--------------|----------------|-----|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 86 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 1 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB02 | 15-06-0884-1-E | 06/10/15 13:26 | Aqueous | GC/MS SS | 06/12/15 | 06/15/15 13:30 | 150612L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 83 | 33-120 | |
| 2-Fluorophenol | 56 | 24-120 | |
| Nitrobenzene-d5 | 79 | 38-120 | |
| p-Terphenyl-d14 | 81 | 41-137 | |
| Phenol-d6 | 35 | 16-120 | |
| 2,4,6-Tribromophenol | 96 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 4 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-4053 | N/A | Aqueous | GC/MS SS | 06/12/15 | 06/13/15 16:19 | 150612L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 5 of 6

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 86 | 33-120 | |
| 2-Fluorophenol | 62 | 24-120 | |
| Nitrobenzene-d5 | 81 | 38-120 | |
| p-Terphenyl-d14 | 87 | 41-137 | |
| Phenol-d6 | 41 | 16-120 | |
| 2,4,6-Tribromophenol | 106 | 27-159 | |


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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 1 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| EB02 | 15-06-0884-1-A | 06/10/15 13:26 | Aqueous | GC/MS Z | 06/15/15 | 06/16/15 03:09 | 150615L021 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | 26 | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 2 of 4

| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | 0.45 | 1.0 | 0.24 | 1.00 | J |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 104 | 80-120 | |
| Dibromofluoromethane | 99 | 78-126 | |
| 1,2-Dichloroethane-d4 | 95 | 75-135 | |
| Toluene-d8 | 108 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 3 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-17406 | N/A | Aqueous | GC/MS Z | 06/15/15 | 06/16/15 00:37 | 150615L021 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 106 | 80-120 | |
| Dibromofluoromethane | 99 | 78-126 | |
| 1,2-Dichloroethane-d4 | 94 | 75-135 | |
| Toluene-d8 | 108 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

06/11/15

Work Order:

15-06-0884

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| EB02 | 15-06-0884-1 | 06/10/15 13:26 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|--------------|
| Phosphorus, Total (24) | 0.020 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 06/13/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 06/16/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 06/16/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 06/16/15 | SM 2320B |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.87 | 1.00 | | mg/L | 06/17/15 | 06/17/15 | SM 2540 C |
| pH | 6.85 | 0.01 | | 1.00 | BV,BU | pH units | N/A | 06/11/15 | SM 4500 H+ B |
| Carbon, Total Organic (24) | 0.78 | 0.50 | 0.24 | 1.00 | | mg/L | 06/19/15 | 06/19/15 | SM 5310 B |

| Method Blank | N/A | Aqueous |
|--------------|-----|---------|
|--------------|-----|---------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|-------|---------------|---------------|-----------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 06/13/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 06/16/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 06/16/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 06/16/15 | SM 2320B |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.87 | 1.00 | | mg/L | 06/17/15 | 06/17/15 | SM 2540 C |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 06/19/15 | 06/19/15 | SM 5310 B |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-06-0862-10 | Sample | Aqueous | IC 10 | N/A | 06/11/15 13:25 | 150611S01 |
| 15-06-0862-10 | Matrix Spike | Aqueous | IC 10 | N/A | 06/11/15 14:35 | 150611S01 |
| 15-06-0862-10 | Matrix Spike Duplicate | Aqueous | IC 10 | N/A | 06/11/15 14:51 | 150611S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.2620 | 250.0 | 257.3 | 103 | 265.5 | 106 | 80-120 | 3 | 0-20 | |
| Chloride | 950.6 | 5000 | 6428 | 110 | 6425 | 109 | 80-120 | 0 | 0-20 | |
| Bromide | 2.124 | 500.0 | 535.0 | 107 | 531.9 | 106 | 80-120 | 1 | 0-20 | |
| Nitrate (as N) | 5.228 | 500.0 | 551.3 | 109 | 548.8 | 109 | 80-120 | 0 | 0-20 | |
| Sulfate | 48.92 | 5000 | 5533 | 110 | 5549 | 110 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

Page 2 of 7

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| EB02 | Sample | Aqueous | ACA 1 | N/A | 06/13/15 14:38 | 150613S01 |
| EB02 | Matrix Spike | Aqueous | ACA 1 | N/A | 06/13/15 14:38 | 150613S01 |
| EB02 | Matrix Spike Duplicate | Aqueous | ACA 1 | N/A | 06/13/15 14:38 | 150613S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | ND | 0.2000 | 0.2211 | 111 | 0.2262 | 113 | 90-110 | 2 | 0-25 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| EB02 | Sample | Aqueous | TOC 8 | 06/19/15 | 06/19/15 22:52 | F0619TOCS1 |
| EB02 | Matrix Spike | Aqueous | TOC 8 | 06/19/15 | 06/19/15 22:52 | F0619TOCS1 |
| EB02 | Matrix Spike Duplicate | Aqueous | TOC 8 | 06/19/15 | 06/19/15 22:52 | F0619TOCS1 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Carbon, Total Organic | 0.7770 | 10.00 | 9.700 | 89 | 9.420 | 86 | 31-145 | 3 | 0-23 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3010A Total
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-06-1010-11 | Sample | Aqueous | ICP 7300 | 06/15/15 | 06/16/15 14:06 | 150615SA4 |
| 15-06-1010-11 | Matrix Spike | Aqueous | ICP 7300 | 06/15/15 | 06/16/15 14:08 | 150615SA4 |
| 15-06-1010-11 | Matrix Spike Duplicate | Aqueous | ICP 7300 | 06/15/15 | 06/16/15 14:09 | 150615SA4 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Calcium | 36.00 | 0.5000 | 36.02 | 4X | 36.11 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 79.75 | 0.5000 | 79.89 | 4X | 79.35 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 0.5344 | 5.000 | 5.794 | 105 | 6.047 | 110 | 83-131 | 4 | 0-7 | |
| Sodium | 44.98 | 5.000 | 49.93 | 4X | 50.05 | 4X | 73-127 | 4X | 0-9 | Q |
| Strontium | 1.051 | 0.5000 | 1.577 | 105 | 1.604 | 111 | 81-123 | 2 | 0-6 | |
| Silicon | 13.42 | 0.5000 | 13.81 | 4X | 13.58 | 4X | 24-180 | 4X | 0-15 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3020A Total
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|---------------------------|------------------------|-------------|----------|------------|---------------|----------------|---------------------|-----|--------|------------|
| 15-06-0784-4 | Sample | | Aqueous | ICP/MS 03 | 06/12/15 | 06/15/15 19:21 | 150612SA3 | | | |
| 15-06-0784-4 | Matrix Spike | | Aqueous | ICP/MS 03 | 06/12/15 | 06/15/15 19:04 | 150612SA3 | | | |
| 15-06-0784-4 | Matrix Spike Duplicate | | Aqueous | ICP/MS 03 | 06/12/15 | 06/15/15 19:07 | 150612SA3 | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.09791 | 98 | 0.1023 | 102 | 80-120 | 4 | 0-20 | |
| Arsenic | ND | 0.1000 | 0.1015 | 102 | 0.1079 | 108 | 80-120 | 6 | 0-20 | |
| Barium | 0.09502 | 0.1000 | 0.1934 | 98 | 0.2042 | 109 | 80-120 | 5 | 0-20 | |
| Beryllium | ND | 0.1000 | 0.09246 | 92 | 0.09985 | 100 | 80-120 | 8 | 0-20 | |
| Cadmium | ND | 0.1000 | 0.09231 | 92 | 0.09672 | 97 | 80-120 | 5 | 0-20 | |
| Chromium | ND | 0.1000 | 0.1110 | 111 | 0.1144 | 114 | 80-120 | 3 | 0-20 | |
| Copper | 0.001257 | 0.1000 | 0.1076 | 106 | 0.1147 | 113 | 80-120 | 6 | 0-20 | |
| Lead | ND | 0.1000 | 0.1036 | 104 | 0.1092 | 109 | 80-120 | 5 | 0-20 | |
| Nickel | 0.001860 | 0.1000 | 0.1066 | 105 | 0.1148 | 113 | 80-120 | 7 | 0-20 | |
| Selenium | ND | 0.1000 | 0.08518 | 85 | 0.08642 | 86 | 80-120 | 1 | 0-20 | |
| Silver | ND | 0.05000 | 0.04771 | 95 | 0.04936 | 99 | 80-120 | 3 | 0-20 | |
| Thallium | ND | 0.1000 | 0.1010 | 101 | 0.1067 | 107 | 80-120 | 6 | 0-20 | |
| Zinc | 0.008488 | 0.1000 | 0.09729 | 89 | 0.1027 | 94 | 80-120 | 5 | 0-20 | |
| Aluminum | ND | 0.1000 | 0.1012 | 101 | 0.1071 | 107 | 80-120 | 6 | 0-20 | |
| Iron | 0.5884 | 5.100 | 6.237 | 111 | 6.665 | 119 | 80-120 | 7 | 0-20 | |
| Manganese | 0.01884 | 0.1000 | 0.1267 | 108 | 0.1361 | 117 | 80-120 | 7 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 7470A Total
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-06-1127-2 | Sample | Aqueous | Mercury 04 | 06/15/15 | 06/15/15 18:35 | 150615S03 |
| 15-06-1127-2 | Matrix Spike | Aqueous | Mercury 04 | 06/15/15 | 06/15/15 18:24 | 150615S03 |
| 15-06-1127-2 | Matrix Spike Duplicate | Aqueous | Mercury 04 | 06/15/15 | 06/15/15 18:26 | 150615S03 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | ND | 0.01000 | 0.008127 | 81 | 0.008788 | 88 | 57-141 | 8 | 0-10 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-06-1097-2 | Sample | Aqueous | GC/MS Z | 06/15/15 | 06/16/15 01:08 | 150615S010 | | | | |
| 15-06-1097-2 | Matrix Spike | Aqueous | GC/MS Z | 06/15/15 | 06/16/15 01:38 | 150615S010 | | | | |
| 15-06-1097-2 | Matrix Spike Duplicate | Aqueous | GC/MS Z | 06/15/15 | 06/16/15 02:08 | 150615S010 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 54.01 | 108 | 55.06 | 110 | 74-122 | 2 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 45.83 | 92 | 47.12 | 94 | 60-144 | 3 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 49.76 | 100 | 50.50 | 101 | 73-120 | 1 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 52.42 | 105 | 52.91 | 106 | 80-122 | 1 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 50.27 | 101 | 51.27 | 103 | 70-120 | 2 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 53.50 | 107 | 54.02 | 108 | 64-142 | 1 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 56.58 | 113 | 59.43 | 119 | 52-136 | 5 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 51.96 | 104 | 52.49 | 105 | 77-125 | 1 | 0-24 | |
| Toluene | ND | 50.00 | 53.36 | 107 | 54.30 | 109 | 72-126 | 2 | 0-23 | |
| Trichloroethene | ND | 50.00 | 55.61 | 111 | 56.00 | 112 | 74-128 | 1 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 46.16 | 92 | 46.06 | 92 | 67-133 | 0 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 93.65 | 94 | 95.11 | 95 | 63-129 | 2 | 0-25 | |
| o-Xylene | ND | 50.00 | 46.76 | 94 | 47.88 | 96 | 62-128 | 2 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 52.11 | 104 | 52.96 | 106 | 68-134 | 2 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3020A Total
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|--------------|-------------|------------|----------------|----------------|-----------------------|
| 15-06-0784-4 | Sample | Aqueous | ICP/MS 03 | 06/12/15 00:00 | 06/15/15 19:21 | 150612SA3 |
| 15-06-0784-4 | PDS | Aqueous | ICP/MS 03 | 06/12/15 00:00 | 06/15/15 19:11 | 150612SA3 |
| Parameter | Sample Conc. | Spike Added | PDS Conc. | PDS %Rec. | %Rec. CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.1008 | 101 | 75-125 | |
| Arsenic | ND | 0.1000 | 0.1011 | 101 | 75-125 | |
| Barium | 0.09502 | 0.1000 | 0.1967 | 102 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.09512 | 95 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.09426 | 94 | 75-125 | |
| Chromium | ND | 0.1000 | 0.1082 | 108 | 75-125 | |
| Copper | 0.001257 | 0.1000 | 0.1063 | 105 | 75-125 | |
| Lead | ND | 0.1000 | 0.1043 | 104 | 75-125 | |
| Nickel | 0.001860 | 0.1000 | 0.1051 | 103 | 75-125 | |
| Selenium | ND | 0.1000 | 0.08213 | 82 | 75-125 | |
| Silver | ND | 0.05000 | 0.04435 | 89 | 75-125 | |
| Thallium | ND | 0.1000 | 0.1023 | 102 | 75-125 | |
| Zinc | 0.008488 | 0.1000 | 0.09970 | 91 | 75-125 | |
| Aluminum | ND | 0.1000 | 0.1024 | 102 | 75-125 | |
| Iron | 0.5884 | 5.100 | 6.014 | 106 | 75-125 | |
| Manganese | 0.01884 | 0.1000 | 0.1234 | 105 | 75-125 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-06-0773-7 | Sample | Aqueous | PH1/BUR03 | N/A | 06/16/15 17:35 | F0616ALKD2 |
| 15-06-0773-7 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 06/16/15 17:35 | F0616ALKD2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|---|---------------------|------------------|------------|---------------|-------------------|
| Alkalinity, Total (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-06-0773-7 | Sample | Aqueous | PH1/BUR03 | N/A | 06/16/15 17:35 | F0616HCOD2 |
| 15-06-0773-7 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 06/16/15 17:35 | F0616HCOD2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Bicarbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-06-0773-7 | Sample | Aqueous | PH1/BUR03 | N/A | 06/16/15 17:35 | F0616CO3D2 |
| 15-06-0773-7 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 06/16/15 17:35 | F0616CO3D2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-06-1111-8 | Sample | Aqueous | SC 2 | 06/17/15 00:00 | 06/17/15 19:00 | F0617TDSD1 |
| 15-06-1111-8 | Sample Duplicate | Aqueous | SC 2 | 06/17/15 00:00 | 06/17/15 19:00 | F0617TDSD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Dissolved | 3245 | 3270 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-06-0986-1 | Sample | Aqueous | PH 1 | N/A | 06/11/15 20:15 | F0611PHD1 |
| 15-06-0986-1 | Sample Duplicate | Aqueous | PH 1 | N/A | 06/11/15 20:15 | F0611PHD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|------------------|---------------------|------------------|------------|---------------|-------------------|
| pH | 7.170 | 7.160 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-12-906-5828 | LCS | Aqueous | IC 10 | N/A | 06/11/15 10:25 | 150611L01 |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Fluoride | | 2.500 | 2.419 | 97 | 90-110 | |
| Chloride | | 50.00 | 52.65 | 105 | 90-110 | |
| Bromide | | 5.000 | 5.106 | 102 | 90-110 | |
| Nitrate (as N) | | 5.000 | 5.257 | 105 | 90-110 | |
| Sulfate | | 50.00 | 52.70 | 105 | 90-110 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-739-107 | LCS | Aqueous | ACA 1 | N/A | 06/13/15 14:38 | 150613L01 | | | |
| 099-12-739-107 | LCSD | Aqueous | ACA 1 | N/A | 06/13/15 14:38 | 150613L01 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Phosphorus, Total | 0.2000 | 0.1957 | 98 | 0.1980 | 99 | 90-110 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-981-106 | LCS | Aqueous | PH1/BUR03 | N/A | 06/16/15 17:35 | F0616ALKB2 | | | |
| 099-15-981-106 | LCSD | Aqueous | PH1/BUR03 | N/A | 06/16/15 17:35 | F0616ALKB2 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 10.00 | 10.00 | 100 | 10.00 | 100 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4633 | LCS | Aqueous | SC 2 | 06/17/15 | 06/17/15 19:00 | F0617TDSL1 | | | |
| 099-12-180-4633 | LCSD | Aqueous | SC 2 | 06/17/15 | 06/17/15 19:00 | F0617TDSL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 85.00 | 85 | 90.00 | 90 | 80-120 | 6 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-097-5675 | LCS | Aqueous | TOC 8 | 06/19/15 | 06/19/15 22:52 | F0619TOCL1 | | | |
| 099-05-097-5675 | LCSD | Aqueous | TOC 8 | 06/19/15 | 06/19/15 22:52 | F0619TOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 10.00 | 8.700 | 87 | 9.320 | 93 | 80-120 | 7 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-15-683-1299 | LCS | Aqueous | ICP 7300 | 06/15/15 | 06/16/15 11:29 | 150615LA4F |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Calcium | | 0.5000 | 0.5004 | 100 | 80-120 | |
| Magnesium | | 0.5000 | 0.5464 | 109 | 80-120 | |
| Potassium | | 5.000 | 5.072 | 101 | 80-120 | |
| Sodium | | 5.000 | 5.025 | 101 | 80-120 | |
| Strontium | | 0.5000 | 0.5134 | 103 | 80-120 | |
| Silicon | | 0.5000 | 0.4580 | 92 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-844 | LCS | Aqueous | ICP/MS 03 | 06/12/15 | 06/15/15 19:00 | 150612LA3F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.09679 | 97 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.1032 | 103 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.09541 | 95 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.1007 | 101 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.09663 | 97 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.1054 | 105 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.1056 | 106 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.09719 | 97 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.1039 | 104 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.09973 | 100 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.04529 | 91 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.09394 | 94 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.1074 | 107 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.1033 | 103 | 80-120 | 73-127 | |
| Iron | 5.100 | 5.389 | 106 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.1063 | 106 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|------------------|
| 099-15-763-571 | LCS | Aqueous | Mercury 04 | 06/15/15 | 06/15/15 18:13 | 150615L03F |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.01000 | 0.01023 | 102 | 85-121 | |

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Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 099-12-529-817 | LCS | Aqueous | GC 41 | 06/15/15 | 06/17/15 18:06 | 150615L01B |
| 099-12-529-817 | LCSD | Aqueous | GC 41 | 06/15/15 | 06/17/15 18:21 | 150615L01B |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
|--------------------|-------------|-----------|-----------|------------|------------|----------|--------|-----|--------|------------|
| Alpha-BHC | 0.5000 | 0.5132 | 103 | 0.5162 | 103 | 50-135 | 36-149 | 1 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.5555 | 111 | 0.5549 | 111 | 50-135 | 36-149 | 0 | 0-25 | |
| Beta-BHC | 0.5000 | 0.4180 | 84 | 0.4220 | 84 | 50-135 | 36-149 | 1 | 0-25 | |
| Heptachlor | 0.5000 | 0.4800 | 96 | 0.4578 | 92 | 50-135 | 36-149 | 5 | 0-25 | |
| Delta-BHC | 0.5000 | 0.5110 | 102 | 0.4912 | 98 | 50-135 | 36-149 | 4 | 0-25 | |
| Aldrin | 0.5000 | 0.4622 | 92 | 0.4394 | 88 | 50-135 | 36-149 | 5 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.4966 | 99 | 0.5009 | 100 | 50-135 | 36-149 | 1 | 0-25 | |
| Endosulfan I | 0.5000 | 0.4422 | 88 | 0.4503 | 90 | 50-135 | 36-149 | 2 | 0-25 | |
| Dieldrin | 0.5000 | 0.4442 | 89 | 0.4501 | 90 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.5264 | 105 | 0.5205 | 104 | 50-135 | 36-149 | 1 | 0-25 | |
| Endrin | 0.5000 | 0.5963 | 119 | 0.5986 | 120 | 50-135 | 36-149 | 0 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.4765 | 95 | 0.4809 | 96 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.5831 | 117 | 0.6197 | 124 | 50-135 | 36-149 | 6 | 0-25 | |
| Endosulfan II | 0.5000 | 0.4348 | 87 | 0.4520 | 90 | 50-135 | 36-149 | 4 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.2994 | 60 | 0.3084 | 62 | 50-135 | 36-149 | 3 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.4985 | 100 | 0.5081 | 102 | 50-135 | 36-149 | 2 | 0-25 | |
| Methoxychlor | 0.5000 | 0.2688 | 54 | 0.2801 | 56 | 50-135 | 36-149 | 4 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------------|------------|---------------|-------------------|
| 099-12-533-1052 | LCS | Aqueous | GC 41 | 06/15/15 | 06/16/15 00:24 | 150615L01 | | | |
| 099-12-533-1052 | LCSD | Aqueous | GC 41 | 06/15/15 | 06/16/15 00:39 | 150615L01 | | | |
| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Aroclor-1016 | 2.000 | 1.837 | 92 | 1.844 | 92 | 50-135 | 0 | 0-25 | |
| Aroclor-1260 | 2.000 | 2.291 | 115 | 1.974 | 99 | 50-135 | 15 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-963-99 | LCS | Aqueous | | GC 26 | 06/12/15 | 06/16/15 21:17 | 150612L11 | | | |
| 099-15-963-99 | LCSD | Aqueous | | GC 26 | 06/12/15 | 06/16/15 22:01 | 150612L11 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.04122 | 103 | 0.04086 | 102 | 30-130 | 13-147 | 1 | 0-30 | |
| Bolstar | 0.04000 | 0.04824 | 121 | 0.04466 | 112 | 30-130 | 13-147 | 8 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.04567 | 114 | 0.04151 | 104 | 30-130 | 13-147 | 10 | 0-30 | |
| Coumaphos | 0.04000 | 0.04288 | 107 | 0.04187 | 105 | 30-130 | 13-147 | 2 | 0-30 | |
| Diazinon | 0.04000 | 0.04904 | 123 | 0.04466 | 112 | 30-130 | 13-147 | 9 | 0-30 | |
| Disulfoton | 0.04000 | 0.05134 | 128 | 0.05013 | 125 | 30-130 | 13-147 | 2 | 0-30 | |
| Ethoprop | 0.04000 | 0.05132 | 128 | 0.04613 | 115 | 30-130 | 13-147 | 11 | 0-30 | |
| Fensulfothion | 0.04000 | 0.04716 | 118 | 0.04446 | 111 | 30-130 | 13-147 | 6 | 0-30 | |
| Fenthion | 0.04000 | 0.04909 | 123 | 0.04464 | 112 | 30-130 | 13-147 | 9 | 0-30 | |
| Merphos | 0.04000 | 0.04695 | 117 | 0.04267 | 107 | 30-130 | 13-147 | 10 | 0-30 | |
| Methyl Parathion | 0.04000 | 0.04531 | 113 | 0.04105 | 103 | 30-130 | 13-147 | 10 | 0-30 | |
| Phorate | 0.04000 | 0.04349 | 109 | 0.04377 | 109 | 30-130 | 13-147 | 1 | 0-30 | |
| Ronnel | 0.04000 | 0.04424 | 111 | 0.04011 | 100 | 30-130 | 13-147 | 10 | 0-30 | |
| Stirophos | 0.04000 | 0.03888 | 97 | 0.03526 | 88 | 30-130 | 13-147 | 10 | 0-30 | |
| Tokuthion | 0.04000 | 0.05031 | 126 | 0.04818 | 120 | 30-130 | 13-147 | 4 | 0-30 | |
| Trichloronate | 0.04000 | 0.04643 | 116 | 0.04206 | 105 | 30-130 | 13-147 | 10 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-034-648 | LCS | Aqueous | GC 40 | 06/12/15 | 06/16/15 17:37 | 150612L12 | | | |
| 095-01-034-648 | LCSD | Aqueous | GC 40 | 06/12/15 | 06/16/15 18:00 | 150612L12 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| 2,4-D | 20.00 | 15.30 | 76 | 16.88 | 84 | 30-130 | 10 | 0-30 | |
| 2,4,5-T | 2.000 | 1.545 | 77 | 1.685 | 84 | 30-130 | 9 | 0-30 | |
| 2,4-DB | 20.00 | 15.28 | 76 | 16.26 | 81 | 30-130 | 6 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-4053 | LCS | Aqueous | | GC/MS SS | 06/12/15 | 06/13/15 16:38 | 150612L03 | | | |
| 095-01-003-4053 | LCSD | Aqueous | | GC/MS SS | 06/12/15 | 06/13/15 16:58 | 150612L03 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 100.0 | 100.4 | 100 | 100.7 | 101 | 61-120 | 51-130 | 0 | 0-20 | |
| Acenaphthylene | 100.0 | 97.82 | 98 | 97.85 | 98 | 55-120 | 44-131 | 0 | 0-20 | |
| Butyl Benzyl Phthalate | 100.0 | 97.05 | 97 | 96.07 | 96 | 56-122 | 45-133 | 1 | 0-20 | |
| 4-Chloro-3-Methylphenol | 100.0 | 90.94 | 91 | 88.20 | 88 | 52-120 | 41-131 | 3 | 0-20 | |
| 2-Chlorophenol | 100.0 | 89.50 | 90 | 87.19 | 87 | 47-120 | 35-132 | 3 | 0-20 | |
| 1,4-Dichlorobenzene | 100.0 | 64.79 | 65 | 60.05 | 60 | 36-120 | 22-134 | 8 | 0-20 | |
| Dimethyl Phthalate | 100.0 | 101.4 | 101 | 101.3 | 101 | 60-120 | 50-130 | 0 | 0-20 | |
| 2,4-Dinitrotoluene | 100.0 | 121.3 | 121 | 121.4 | 121 | 61-121 | 51-131 | 0 | 0-20 | |
| Fluorene | 100.0 | 105.3 | 105 | 103.3 | 103 | 67-120 | 58-129 | 2 | 0-20 | |
| N-Nitroso-di-n-propylamine | 100.0 | 81.74 | 82 | 81.19 | 81 | 39-123 | 25-137 | 1 | 0-20 | |
| Naphthalene | 100.0 | 81.91 | 82 | 78.06 | 78 | 54-120 | 43-131 | 5 | 0-20 | |
| 4-Nitrophenol | 100.0 | 49.71 | 50 | 47.00 | 47 | 14-120 | 0-138 | 6 | 0-20 | |
| Pentachlorophenol | 100.0 | 84.07 | 84 | 84.13 | 84 | 31-127 | 15-143 | 0 | 0-20 | |
| Phenol | 100.0 | 41.40 | 41 | 40.75 | 41 | 17-120 | 0-137 | 2 | 0-20 | |
| Pyrene | 100.0 | 96.01 | 96 | 94.06 | 94 | 58-124 | 47-135 | 2 | 0-20 | |
| 1,2,4-Trichlorobenzene | 100.0 | 76.35 | 76 | 70.36 | 70 | 49-120 | 37-132 | 8 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 06/11/15
Work Order: 15-06-0884
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-17406 | LCS | Aqueous | GC/MS Z | 06/15/15 | 06/15/15 23:37 | 150615L021 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 54.96 | 110 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 46.32 | 93 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 51.07 | 102 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 54.12 | 108 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 51.92 | 104 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 54.23 | 108 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 57.78 | 116 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 53.10 | 106 | 80-123 | 73-130 | |
| Toluene | 50.00 | 54.81 | 110 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 57.32 | 115 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 45.62 | 91 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 97.35 | 97 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 48.32 | 97 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 56.94 | 114 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits

Glossary of Terms and Qualifiers

Work Order: 15-06-0884

Page 1 of 1

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|---|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| CI | See case narrative. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |
| | Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time. |
| | A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations. |

00002

00281

FedEx Package
Express **US Airbill**FedEx
Tracking
Number

8062 7398 5440

0215

Recipient

1 From

Date

Sender's
Name

Phil Pearce

Phone

210 877-2847

Company

SWCA INC

Address

6200 UTSA BLVD STE 102

Dept./Floor/Suite/Room

City

SAN ANTONIO

State

TX

ZIP

78249-1618

2 Your Internal Billing Reference

27122.02.05

3 To

Recipient's
Name

Phone

714 895-5494

Company

Address

We cannot deliver to P.O. boxes or P.O. ZIP codes.

Dept./Floor/Suite/Room

Address

Use this line for the HOLD location address or for continuation of your shipping address.

City

Garden Grove

State

CA

ZIP

92841

0115649666



8062 7398 5440

THU - 11 JUN AA

4 Express Package Service * To most locations.
NOTE: Service order has changed. Please select carefully.

Next Business Day

- ☐ **FedEx First Overnight**
Earliest next business morning delivery to select locations. Friday shipments will be delivered on Monday unless SATURDAY Delivery is selected.
- ☐ **FedEx Priority Overnight**
Next business morning. Friday shipments will be delivered on Monday unless SATURDAY Delivery is selected.
- ☒ **FedEx Standard Overnight**
Next business afternoon. Saturday Delivery NOT available.

2 or 3 Business Days

- ☐ **FedEx 2Day A.M.**
Second business morning. Saturday Delivery NOT available.
- ☐ **FedEx 2Day**
Second business afternoon. Thursday shipments will be delivered on Monday unless SATURDAY Delivery is selected.
- ☐ **FedEx Express Saver**
Third business day. Saturday Delivery NOT available.

5 Packaging * Declared value limit \$500.

- ☐ FedEx Envelope* ☐ FedEx Pak* ☐ FedEx Box ☐ FedEx Tube

6 Special Handling and Delivery Signature Options

- ☐ **SATURDAY Delivery**
NOT available for FedEx Standard Overnight, FedEx 2Day A.M., or FedEx Express Saver.

- ☒ **No Signature Required**
Package may be left without obtaining a signature for delivery.
- ☐ **Direct Signature**
Someone at recipient's address may sign for delivery. Fee applies.
- ☐ **Indirect**
If no one is at address, address is resident's.

Does this shipment contain dangerous goods?

- ☒ **No** ☐ **Yes** ☐ **Yes**
As per attached Shipper's Declaration. Shipper's Declaration not required.
- ☐ **Dry Ice**
Dry Ice, 9, UN 1845
- ☐ **Cargo Aircraft**

7 Payment Bill to:

- ☒ **Sender** ☐ **Recipient** ☐ **Third Party** ☐ **Credit Card**

Enter FedEx Acct. No. or Credit Card No. below:

Total Packages Total Weight

Total Packages 1 Total Weight 00 lbs.

Credit Card Auth.

Your liability is limited to US\$100 unless you declare a higher value. See the current FedEx Service Guide for details.

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SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 1

CLIENT: SWCA Env'l. Consultants

DATE: 06 / 11 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 2-7 °C (w/ CF): 2-4 °C; ☐ Blank ☒ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 836

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 836

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 836

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznnna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 g Cube ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 836

s = H₂SO₄, u = ultra-pure, znnna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 836

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 210 Lead
Sample Number: AA92419
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 00:36
Receipt Date/Time: 1/22/15 10:41

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time was exceeded, per customer proceed with analyses.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92419-A | E. coli | √ | SM 9223B-2004 | 2200 | MPN/100 mL | H | 1 | 41460 | 1/22/15 | 13:20 | RSC |
| AA92419-A | E. Coli Holding Time - IDEXX Colilert | | NA | 12.73 | hours | | 0.00 | 41459 | 1/22/15 | 13:20 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 24, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41460

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/24/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 240 Lead
Sample Number: AA92420
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 00:51
Receipt Date/Time: 1/22/15 10:41

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time was exceeded, per customer proceed with analyses.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92420-A | E. coli | √ | SM 9223B-2004 | 910 | MPN/100 mL | H | 1 | 41460 | 1/22/15 | 13:20 | RSC |
| AA92420-A | E. Coli Holding Time - IDEXX Colilert | | NA | 12.48 | hours | | 0.00 | 41459 | 1/22/15 | 13:20 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 24, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41460

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/24/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 24, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 250 Lead
Sample Number: AA92421
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 00:35
Receipt Date/Time: 1/22/15 10:41

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time was exceeded, per customer proceed with analyses.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92421-A | E. coli | √ | SM 9223B-2004 | 630 | MPN/100 mL | H | 1 | 41460 | 1/22/15 | 13:20 | RSC |
| AA92421-A | E. Coli Holding Time - IDEXX Colilert | | NA | 12.75 | hours | | 0.00 | 41459 | 1/22/15 | 13:20 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 24, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41460

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



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1/24/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 24, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 260 Lead
Sample Number: AA92422
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 01:10
Receipt Date/Time: 1/22/15 10:41

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time was exceeded, per customer proceed with analyses.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92422-A | E. coli | √ | SM 9223B-2004 | 370 | MPN/100 mL | H | 1 | 41460 | 1/22/15 | 13:20 | RSC |
| AA92422-A | E. Coli Holding Time - IDEXX Colilert | | NA | 12.17 | hours | | 0.00 | 41459 | 1/22/15 | 13:20 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 24, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41460

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



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1/24/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 270 Lead
Sample Number: AA92423
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 00:54
Receipt Date/Time: 1/22/15 10:41

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time was exceeded, per customer proceed with analyses.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92423-A | E. coli | √ | SM 9223B-2004 | 1300 | MPN/100 mL | H | 1 | 41460 | 1/22/15 | 13:20 | RSC |
| AA92423-A | E. Coli Holding Time - IDEXX Colilert | | NA | 12.43 | hours | | 0.00 | 41459 | 1/22/15 | 13:20 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41460

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/24/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 210 Peak
Sample Number: AA92424
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 08:14
Receipt Date/Time: 1/22/15 10:42

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Analysis Comments: AA92424-A E. Coli Holding Time - IDEXX Colilert
Not analyzed per customer request.
AA92424-A E. coli
Not analyzed per customer request.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92424-A | E. coli | √ | SM 9223B-2004 | Scratched | MPN/100 mL | * | | | 1/23/15 | 11:08 | JLM |
| AA92424-A | E. Coli Holding Time - IDEXX Colilert | | NA | Scratched | hours | * | | | 1/23/15 | 11:08 | JLM |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 240 Peak
Sample Number: AA92425
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 08:31
Receipt Date/Time: 1/22/15 10:42

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Analysis Comments: AA92425-A E. Coli Holding Time - IDEXX Colilert
Not analyzed per customer request.
AA92425-A E. coli
Not analyzed per customer request.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92425-A | E. coli | √ | SM 9223B-2004 | Scratched | MPN/100 mL | * | | | 1/23/15 | 11:08 | JLM |
| AA92425-A | E. Coli Holding Time - IDEXX Colilert | | NA | Scratched | hours | * | | | 1/23/15 | 11:08 | JLM |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 250 Peak
Sample Number: AA92426
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 08:10
Receipt Date/Time: 1/22/15 10:42

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Analysis Comments: AA92426-A E. Coli Holding Time - IDEXX Colilert
Not analyzed per customer request.
AA92426-A E. coli
Not analyzed per customer request.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92426-A | E. coli | √ | SM 9223B-2004 | Scratched | MPN/100 mL | * | | | 1/23/15 | 11:08 | JLM |
| AA92426-A | E. Coli Holding Time - IDEXX Colilert | | NA | Scratched | hours | * | | | 1/23/15 | 11:08 | JLM |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 260 Peak
Sample Number: AA92427
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 08:43
Receipt Date/Time: 1/22/15 10:42

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Analysis Comments: AA92427-A E. Coli Holding Time - IDEXX Colilert
Not analyzed per customer request.
AA92427-A E. coli
Not analyzed per customer request.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92427-A | E. coli | √ | SM 9223B-2004 | Scratched | MPN/100 mL | * | | | 1/23/15 | 11:08 | JLM |
| AA92427-A | E. Coli Holding Time - IDEXX Colilert | | NA | Scratched | hours | * | | | 1/23/15 | 11:08 | JLM |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 270 Peak
Sample Number: AA92428
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 08:30
Receipt Date/Time: 1/22/15 10:42

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Analysis Comments: AA92428-A E. Coli Holding Time - IDEXX Colilert
Not analyzed per customer request.
AA92428-A E. coli
Not analyzed per customer request.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92428-A | E. coli | √ | SM 9223B-2004 | Scratched | MPN/100 mL | * | | | 1/23/15 | 11:08 | JLM |
| AA92428-A | E. Coli Holding Time - IDEXX Colilert | | NA | Scratched | hours | * | | | 1/23/15 | 11:08 | JLM |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HCS 210 Peak
Sample Number: AA92461
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 16:28
Receipt Date/Time: 1/23/15 09:00

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments:

Hold time was exceeded, per customer proceed with analyses. A discrepancy was noted on the COC at sample validation. A time gap exists between a time of relinquish and a time of receipt. This occurred prior to sample submission to the laboratory.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92461-A | E. coli | √ | SM 9223B-2004 | 1600 | MPN/100 mL | H | 1 | 41481 | 1/23/15 | 15:03 | RSC/MAG |
| AA92461-A | E. Coli Holding Time - IDEXX Colilert | | NA | 22.58 | hours | | 0.00 | 41479 | 1/23/15 | 15:03 | RSC/MAG |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41481

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 1 of 2

Client:

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6200 UTSA Blvd Ste 102
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Sample Location: HCS 240 Peak
Sample Number: AA92462
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 16:51
Receipt Date/Time: 1/23/15 09:00

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments:

Hold time was exceeded, per customer proceed with analyses. A discrepancy was noted on the COC at sample validation. A time gap exists between a time of relinquish and a time of receipt. This occurred prior to sample submission to the laboratory.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92462-A | E. coli | √ | SM 9223B-2004 | 550 | MPN/100 mL | H | 1 | 41481 | 1/23/15 | 15:03 | RSC/MAG |
| AA92462-A | E. Coli Holding Time - IDEXX Colilert | | NA | 22.20 | hours | | 0.00 | 41479 | 1/23/15 | 15:03 | RSC/MAG |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41481

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 250 Peak
Sample Number: AA92463
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 16:40
Receipt Date/Time: 1/23/15 09:00

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments:

Hold time was exceeded, per customer proceed with analyses. A discrepancy was noted on the COC at sample validation. A time gap exists between a time of relinquish and a time of receipt. This occurred prior to sample submission to the laboratory.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92463-A | E. coli | √ | SM 9223B-2004 | 1800 | MPN/100 mL | H | 1 | 41481 | 1/23/15 | 15:03 | RSC/MAG |
| AA92463-A | E. Coli Holding Time - IDEXX Colilert | | NA | 22.38 | hours | | 0.00 | 41479 | 1/23/15 | 15:03 | RSC/MAG |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41481

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 260 Peak
Sample Number: AA92464
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 17:03
Receipt Date/Time: 1/23/15 09:00

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments:

Hold time was exceeded, per customer proceed with analyses. A discrepancy was noted on the COC at sample validation. A time gap exists between a time of relinquish and a time of receipt. This occurred prior to sample submission to the laboratory.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92464-A | E. coli | √ | SM 9223B-2004 | 10000 | MPN/100 mL | H | 1 | 41481 | 1/23/15 | 15:03 | RSC/MAG |
| AA92464-A | E. Coli Holding Time - IDEXX Colilert | | NA | 22.00 | hours | | 0.00 | 41479 | 1/23/15 | 15:03 | RSC/MAG |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41481

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 270 Peak
Sample Number: AA92465
Sample Matrix: Non Potable Water

Collection Date/Time: 1/22/15 16:55
Receipt Date/Time: 1/23/15 09:00

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments:

Hold time was exceeded, per customer proceed with analyses. A discrepancy was noted on the COC at sample validation. A time gap exists between a time of relinquish and a time of receipt. This occurred prior to sample submission to the laboratory.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92465-A | E. coli | √ | SM 9223B-2004 | 5500 | MPN/100 mL | H | 1 | 41481 | 1/23/15 | 15:03 | RSC/MAG |
| AA92465-A | E. Coli Holding Time - IDEXX Colilert | | NA | 22.13 | hours | | 0.00 | 41479 | 1/23/15 | 15:03 | RSC/MAG |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41481

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
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Sample Location: HCS 210 Trail
Sample Number: AA92466
Sample Matrix: Non Potable Water

Collection Date/Time: 1/23/15 03:01
Receipt Date/Time: 1/23/15 09:00

CASE NARRATIVE

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Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Hold time was exceeded, per customer proceed with analyses.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92466-A | E. coli | √ | SM 9223B-2004 | 880 | MPN/100 mL | H | 1 | 41481 | 1/23/15 | 15:03 | RSC/MAG |
| AA92466-A | E. Coli Holding Time - IDEXX Colilert | | NA | 12.03 | hours | | 0.00 | 41479 | 1/23/15 | 15:03 | RSC/MAG |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41481

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 240 Trail
Sample Number: AA92467
Sample Matrix: Non Potable Water

Collection Date/Time: 1/23/15 03:21
Receipt Date/Time: 1/23/15 09:00

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time was exceeded, per customer proceed with analyses.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92467-A | E. coli | √ | SM 9223B-2004 | 110 | MPN/100 mL | H | 1 | 41481 | 1/23/15 | 15:03 | RSC/MAG |
| AA92467-A | E. Coli Holding Time - IDEXX Colilert | | NA | 11.70 | hours | | 0.00 | 41479 | 1/23/15 | 15:03 | RSC/MAG |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41481

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 250 Trail
Sample Number: AA92468
Sample Matrix: Non Potable Water

Collection Date/Time: 1/23/15 03:03
Receipt Date/Time: 1/23/15 09:00

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time was exceeded, per customer proceed with analyses.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92468-A | E. coli | √ | SM 9223B-2004 | 340 | MPN/100 mL | H | 1 | 41482 | 1/23/15 | 15:03 | RSC/MAG |
| AA92468-A | E. Coli Holding Time - IDEXX Colilert | | NA | 12.00 | hours | | 0.00 | 41480 | 1/23/15 | 15:03 | RSC/MAG |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41482

QC Analyte Name

Initial Blank for E. coli

Log Range for E. coli

Result

Absent

0.1246

Units

Qualifier

H

H

Lower

0.0

Acceptance Criteria

Target

Absent

Upper

0.5



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 1 of 2

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Sample Location: HCS 260 Trail
Sample Number: AA92469
Sample Matrix: Non Potable Water

Collection Date/Time: 1/23/15 03:36
Receipt Date/Time: 1/23/15 09:00

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time was exceeded, per customer proceed with analyses.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92469-A | E. coli | √ | SM 9223B-2004 | 2300 | MPN/100 mL | H | 1 | 41482 | 1/23/15 | 15:03 | RSC/MAG |
| AA92469-A | E. Coli Holding Time - IDEXX Colilert | | NA | 11.45 | hours | | 0.00 | 41480 | 1/23/15 | 15:03 | RSC/MAG |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41482

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

H

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS 270 Trail
Sample Number: AA92470
Sample Matrix: Non Potable Water

Collection Date/Time: 1/23/15 03:30
Receipt Date/Time: 1/23/15 09:00

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time was exceeded, per customer proceed with analyses.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92470-A | E. coli | √ | SM 9223B-2004 | 4400 | MPN/100 mL | H | 1 | 41482 | 1/23/15 | 15:03 | RSC/MAG |
| AA92470-A | E. Coli Holding Time - IDEXX Colilert | | NA | 11.55 | hours | | 0.00 | 41480 | 1/23/15 | 15:03 | RSC/MAG |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41482

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

H

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: FD HCS 260 Trail
Sample Number: AA92471
Sample Matrix: Non Potable Water

Collection Date/Time: 1/23/15 03:36
Receipt Date/Time: 1/23/15 09:00

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time was exceeded, per customer proceed with analyses.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92471-A | E. coli | √ | SM 9223B-2004 | 3400 | MPN/100 mL | H | 1 | 41482 | 1/23/15 | 15:03 | RSC/MAG |
| AA92471-A | E. Coli Holding Time - IDEXX Colilert | | NA | 11.45 | hours | | 0.00 | 41480 | 1/23/15 | 15:03 | RSC/MAG |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41482

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

H

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: FD HCS 270 Trail
Sample Number: AA92472
Sample Matrix: Non Potable Water

Collection Date/Time: 1/23/15 03:30
Receipt Date/Time: 1/23/15 09:00

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time was exceeded, per customer proceed with analyses.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA92472-A | E. coli | √ | SM 9223B-2004 | 2500 | MPN/100 mL | H | 1 | 41482 | 1/23/15 | 15:03 | RSC/MAG |
| AA92472-A | E. Coli Holding Time - IDEXX Colilert | | NA | 11.55 | hours | | 0.00 | 41480 | 1/23/15 | 15:03 | RSC/MAG |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



January 27, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-41482

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

H

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

1/27/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable



Calscience



WORK ORDER NUMBER: 15-01-1404

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: SWCA Environmental Consultants

Client Project Name: EAA 27122

Attention: Philip Pearce
6200 UTSA Blvd.
Suite 102
San Antonio, TX 78249-1618

Approved for release on 02/09/2015 by:
Don Burley
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

Contents

Client Project Name: EAA 27122
Work Order Number: 15-01-1404

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Work Order Narrative

Work Order: 15-01-1404

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Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 01/23/15. They were assigned to Work Order 15-01-1404.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

EPA 6010B Silicon LCS recovery was below the control limits. Silicon was detected in each of the samples. The results may be biased low.

EPA 8141A Merphos LCS and LCSD recoveries were above the control limits. Merphos was not detected in the samples. The results are reported without further clarification.



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Sample Summary

| | | |
|--|-----------------------|----------------|
| Client: SWCA Environmental Consultants | Work Order: | 15-01-1404 |
| 6200 UTSA Blvd., Suite 102 | Project Name: | EAA 27122 |
| San Antonio, TX 78249-1618 | PO Number: | |
| | Date/Time Received: | 01/23/15 10:00 |
| | Number of Containers: | 45 |

Attn: Philip Pearce

| Sample Identification | Lab Number | Collection Date and Time | Number of Containers | Matrix |
|-----------------------|--------------|--------------------------|----------------------|---------|
| HCS 210 Lead | 15-01-1404-1 | 01/22/15 00:36 | 9 | Aqueous |
| HCS 240 Lead | 15-01-1404-2 | 01/22/15 00:51 | 9 | Aqueous |
| HCS 250 Lead | 15-01-1404-3 | 01/22/15 00:35 | 9 | Aqueous |
| HCS 260 Lead | 15-01-1404-4 | 01/22/15 01:10 | 9 | Aqueous |
| HCS 270 Lead | 15-01-1404-5 | 01/22/15 00:54 | 9 | Aqueous |

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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

Page 1 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-01-1404-1-F | 01/22/15 00:36 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:24 | 150126L6FF |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 15.4 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 1.05 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 2.02 | 0.500 | 0.103 | 1.00 | |
| Sodium | 0.624 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.0466 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 0.842 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS 240 Lead | 15-01-1404-2-F | 01/22/15 00:51 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:31 | 150126L6FF |
|--------------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 73.9 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 14.4 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.77 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.626 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.47 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS 250 Lead | 15-01-1404-3-F | 01/22/15 00:35 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:32 | 150126L6FF |
|--------------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 73.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 14.2 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.58 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.2 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.622 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.36 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

Page 2 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-01-1404-4-F | 01/22/15 01:10 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:34 | 150126L6FF |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 75.1 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 14.5 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.53 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.6 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.637 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.33 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS 270 Lead | 15-01-1404-5-F | 01/22/15 00:54 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:36 | 150126L6FF |
|--------------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 67.3 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 13.2 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.76 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.7 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.571 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.11 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|-----------------|-----|---------|----------|----------|----------------|------------|
| Method Blank | 099-15-683-1120 | N/A | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 17:52 | 150126L6FF |
|--------------|-----------------|-----|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | ND | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 0.0522 | 0.100 | 0.00336 | 1.00 | J |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 1 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-01-1404-1-F | 01/22/15 00:36 | Aqueous | ICP/MS 03 | 01/26/15 | 01/27/15 13:20 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000288 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.00845 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00148 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.000670 | 0.00100 | 0.000132 | 1.00 | J |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00431 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | 0.0115 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0168 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00315 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 2 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Lead | 15-01-1404-2-F | 01/22/15 00:51 | Aqueous | ICP/MS 03 | 01/26/15 | 01/27/15 13:23 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000113 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0567 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000650 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00161 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000312 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00549 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0160 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0564 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00186 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 3 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-01-1404-3-F | 01/22/15 00:35 | Aqueous | ICP/MS 03 | 01/26/15 | 01/27/15 13:26 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000113 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0539 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000819 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00155 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000333 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00891 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00613 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0536 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00205 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-01-1404-4-F | 01/22/15 01:10 | Aqueous | ICP/MS 03 | 01/26/15 | 01/27/15 13:30 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0546 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000651 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00168 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000335 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00195 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | 0.00950 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0440 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00260 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-01-1404-5-F | 01/22/15 00:54 | Aqueous | ICP/MS 03 | 01/26/15 | 01/27/15 13:33 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|-----------|---------|-----------|------|------------|
| Antimony | 0.000194 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0494 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000789 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | 0.0000928 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00145 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000350 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0146 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00488 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0249 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00226 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 6 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-719 | N/A | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 08:18 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-01-1404-1-F | 01/22/15 00:36 | Aqueous | Mercury 04 | 01/29/15 | 01/29/15 17:28 | 150129L06F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 240 Lead | 15-01-1404-2-F | 01/22/15 00:51 | Aqueous | Mercury 04 | 01/29/15 | 01/29/15 17:30 | 150129L06F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 250 Lead | 15-01-1404-3-F | 01/22/15 00:35 | Aqueous | Mercury 04 | 01/29/15 | 01/29/15 17:32 | 150129L06F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 260 Lead | 15-01-1404-4-F | 01/22/15 01:10 | Aqueous | Mercury 04 | 01/29/15 | 01/29/15 17:35 | 150129L06F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 270 Lead | 15-01-1404-5-F | 01/22/15 00:54 | Aqueous | Mercury 04 | 01/29/15 | 01/29/15 17:37 | 150129L06F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|-----|---------|------------|----------|----------------|------------|
| Method Blank | 099-15-763-484 | N/A | Aqueous | Mercury 04 | 01/29/15 | 01/29/15 17:01 | 150129L06F |
|--------------|----------------|-----|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

Page 1 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|-----------------------|-----------------------|-----------------------|----------------|-------------------|-----------------------|-------------------|
| HCS 210 Lead | 15-01-1404-1-I | 01/22/15 00:36 | Aqueous | LC/TQ 2 | 01/29/15 | 02/02/15 18:49 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | <u>RL</u> | | <u>DF</u> | | <u>Qualifiers</u> |
| Caffeine | | 480 | 100 | | 10.0 | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | <u>Control Limits</u> | | <u>Qualifiers</u> | | |
| Caffeine-C13 (Surrogate) | | 45 | 31-200 | | | | |
| HCS 240 Lead | 15-01-1404-2-I | 01/22/15 00:51 | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 13:16 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | <u>RL</u> | | <u>DF</u> | | <u>Qualifiers</u> |
| Caffeine | | 47 | 10 | | 1.00 | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | <u>Control Limits</u> | | <u>Qualifiers</u> | | |
| Caffeine-C13 (Surrogate) | | 54 | 31-200 | | | | |
| HCS 250 Lead | 15-01-1404-3-I | 01/22/15 00:35 | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 13:25 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | <u>RL</u> | | <u>DF</u> | | <u>Qualifiers</u> |
| Caffeine | | 55 | 10 | | 1.00 | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | <u>Control Limits</u> | | <u>Qualifiers</u> | | |
| Caffeine-C13 (Surrogate) | | 46 | 31-200 | | | | |
| HCS 260 Lead | 15-01-1404-4-I | 01/22/15 01:10 | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 13:33 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | <u>RL</u> | | <u>DF</u> | | <u>Qualifiers</u> |
| Caffeine | | 57 | 10 | | 1.00 | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | <u>Control Limits</u> | | <u>Qualifiers</u> | | |
| Caffeine-C13 (Surrogate) | | 48 | 31-200 | | | | |
| HCS 270 Lead | 15-01-1404-5-I | 01/22/15 00:54 | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 17:44 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | <u>RL</u> | | <u>DF</u> | | <u>Qualifiers</u> |
| Caffeine | | 95 | 10 | | 1.00 | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | <u>Control Limits</u> | | <u>Qualifiers</u> | | |
| Caffeine-C13 (Surrogate) | | 47 | 31-200 | | | | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|-------------------|---------------------|---------|-----------------------|---------------|--------------------|-------------------|
| Method Blank | 099-16-376-10 | N/A | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 17:06 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | <u>DF</u> | | <u>Qualifiers</u> |
| Caffeine | | ND | | 10 | 1.00 | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 49 | | 31-200 | | | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

Page 1 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-01-1404-1-I | 01/22/15 00:36 | Aqueous | GC 51 | 01/24/15 | 02/04/15 07:57 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.099 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.099 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.099 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.099 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.099 | 0.028 | 1.00 | |
| Aldrin | ND | 0.099 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.099 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.099 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.099 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.099 | 0.026 | 1.00 | |
| Endrin | ND | 0.099 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.099 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.099 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.099 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.099 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.099 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.099 | 0.025 | 1.00 | |
| Chlordane | ND | 0.99 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.099 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 99 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 98 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

Page 2 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Lead | 15-01-1404-2-I | 01/22/15 00:51 | Aqueous | GC 51 | 01/24/15 | 02/04/15 08:12 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.096 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.096 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.096 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.096 | 0.025 | 1.00 | |
| Delta-BHC | ND | 0.096 | 0.027 | 1.00 | |
| Aldrin | ND | 0.096 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.096 | 0.024 | 1.00 | |
| Endosulfan I | ND | 0.096 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.096 | 0.027 | 1.00 | |
| 4,4'-DDE | ND | 0.096 | 0.026 | 1.00 | |
| Endrin | ND | 0.096 | 0.029 | 1.00 | |
| Endrin Aldehyde | ND | 0.096 | 0.025 | 1.00 | |
| 4,4'-DDD | ND | 0.096 | 0.026 | 1.00 | |
| Endosulfan II | ND | 0.096 | 0.026 | 1.00 | |
| 4,4'-DDT | ND | 0.096 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.096 | 0.028 | 1.00 | |
| Methoxychlor | ND | 0.096 | 0.024 | 1.00 | |
| Chlordane | ND | 0.96 | 0.32 | 1.00 | |
| Toxaphene | ND | 1.9 | 0.57 | 1.00 | |
| Endrin Ketone | ND | 0.096 | 0.023 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 105 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 97 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-01-1404-3-I | 01/22/15 00:35 | Aqueous | GC 51 | 01/24/15 | 02/04/15 08:26 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.095 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.095 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.095 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.095 | 0.025 | 1.00 | |
| Delta-BHC | ND | 0.095 | 0.027 | 1.00 | |
| Aldrin | ND | 0.095 | 0.025 | 1.00 | |
| Heptachlor Epoxide | ND | 0.095 | 0.024 | 1.00 | |
| Endosulfan I | ND | 0.095 | 0.026 | 1.00 | |
| Dieldrin | ND | 0.095 | 0.027 | 1.00 | |
| 4,4'-DDE | ND | 0.095 | 0.025 | 1.00 | |
| Endrin | ND | 0.095 | 0.029 | 1.00 | |
| Endrin Aldehyde | ND | 0.095 | 0.025 | 1.00 | |
| 4,4'-DDD | ND | 0.095 | 0.026 | 1.00 | |
| Endosulfan II | ND | 0.095 | 0.026 | 1.00 | |
| 4,4'-DDT | ND | 0.095 | 0.025 | 1.00 | |
| Endosulfan Sulfate | ND | 0.095 | 0.028 | 1.00 | |
| Methoxychlor | ND | 0.095 | 0.024 | 1.00 | |
| Chlordane | ND | 0.95 | 0.31 | 1.00 | |
| Toxaphene | ND | 1.9 | 0.56 | 1.00 | |
| Endrin Ketone | ND | 0.095 | 0.023 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 91 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-01-1404-4-I | 01/22/15 01:10 | Aqueous | GC 51 | 01/24/15 | 02/04/15 08:40 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.098 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.098 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.098 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.098 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.098 | 0.028 | 1.00 | |
| Aldrin | ND | 0.098 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.098 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.098 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.098 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.098 | 0.026 | 1.00 | |
| Endrin | ND | 0.098 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.098 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.098 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.098 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.098 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.098 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.098 | 0.025 | 1.00 | |
| Chlordane | ND | 0.98 | 0.32 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.58 | 1.00 | |
| Endrin Ketone | ND | 0.098 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 101 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 96 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

Page 5 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-01-1404-5-I | 01/22/15 00:54 | Aqueous | GC 51 | 01/24/15 | 02/04/15 08:55 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.099 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.099 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.099 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.099 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.099 | 0.028 | 1.00 | |
| Aldrin | ND | 0.099 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.099 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.099 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.099 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.099 | 0.026 | 1.00 | |
| Endrin | ND | 0.099 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.099 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.099 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.099 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.099 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.099 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.099 | 0.025 | 1.00 | |
| Chlordane | ND | 0.99 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.099 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 95 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 92 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-778 | N/A | Aqueous | GC 51 | 01/24/15 | 02/04/15 04:50 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 97 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 90 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

Page 1 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HCS 210 Lead | 15-01-1404-1-I | 01/22/15 00:36 | Aqueous | GC 31 | 01/24/15 | 02/04/15 11:41 | 150124L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 0.99 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.99 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 0.99 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 0.99 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 0.99 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.99 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.99 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.99 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 73 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 84 | 50-135 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HCS 240 Lead | 15-01-1404-2-I | 01/22/15 00:51 | Aqueous | GC 31 | 01/24/15 | 02/04/15 12:00 | 150124L09 |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 0.96 | 0.28 | 1.00 | |
| Aroclor-1221 | ND | 0.96 | 0.27 | 1.00 | |
| Aroclor-1232 | ND | 0.96 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.96 | 0.17 | 1.00 | |
| Aroclor-1248 | ND | 0.96 | 0.19 | 1.00 | |
| Aroclor-1254 | ND | 0.96 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.96 | 0.25 | 1.00 | |
| Aroclor-1262 | ND | 0.96 | 0.25 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 79 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 88 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

Page 2 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-01-1404-3-I | 01/22/15 00:35 | Aqueous | GC 31 | 01/24/15 | 02/04/15 12:19 | 150124L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|------|------|------|------------|
| Aroclor-1016 | ND | 0.95 | 0.28 | 1.00 | |
| Aroclor-1221 | ND | 0.95 | 0.27 | 1.00 | |
| Aroclor-1232 | ND | 0.95 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.95 | 0.17 | 1.00 | |
| Aroclor-1248 | ND | 0.95 | 0.19 | 1.00 | |
| Aroclor-1254 | ND | 0.95 | 0.21 | 1.00 | |
| Aroclor-1260 | ND | 0.95 | 0.25 | 1.00 | |
| Aroclor-1262 | ND | 0.95 | 0.25 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 72 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 79 | 50-135 | |

| HCS 260 Lead | 15-01-1404-4-I | 01/22/15 01:10 | Aqueous | GC 31 | 01/24/15 | 02/04/15 12:38 | 150124L09 |
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|------|------|------|------------|
| Aroclor-1016 | ND | 0.98 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.98 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 0.98 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.98 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 0.98 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.98 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.98 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.98 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 70 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 78 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

Page 3 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-01-1404-5-I | 01/22/15 00:54 | Aqueous | GC 31 | 01/24/15 | 02/04/15 12:57 | 150124L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|------|------|------|------------|
| Aroclor-1016 | ND | 0.99 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.99 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 0.99 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 0.99 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 0.99 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.99 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.99 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.99 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 70 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 78 | 50-135 | |

| Method Blank | 099-12-533-1000 | N/A | Aqueous | GC 31 | 01/24/15 | 02/04/15 11:03 | 150124L09 |
|--------------|-----------------|-----|---------|-------|----------|----------------|-----------|
|--------------|-----------------|-----|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 71 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 79 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 1 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-01-1404-1-I | 01/22/15 00:36 | Aqueous | GC 26 | 01/24/15 | 02/06/15 20:53 | 150124L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 124 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 2 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Lead | 15-01-1404-2-I | 01/22/15 00:51 | Aqueous | GC 26 | 01/24/15 | 02/06/15 21:37 | 150124L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 126 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 3 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-01-1404-3-I | 01/22/15 00:35 | Aqueous | GC 26 | 01/24/15 | 02/06/15 22:21 | 150124L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 87 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-01-1404-4-I | 01/22/15 01:10 | Aqueous | GC 26 | 01/24/15 | 02/06/15 23:05 | 150124L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 122 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 5 of 6

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-01-1404-5-I | 01/22/15 00:54 | Aqueous | GC 26 | 01/24/15 | 02/06/15 23:49 | 150124L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 122 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-82 | N/A | Aqueous | GC 26 | 01/24/15 | 02/06/15 18:41 | 150124L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 115 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

Page 1 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-01-1404-1-I | 01/22/15 00:36 | Aqueous | GC 40 | 01/26/15 | 01/29/15 20:17 | 150126L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 66 | 0-123 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|
| HCS 240 Lead | 15-01-1404-2-I | 01/22/15 00:51 | Aqueous | GC 40 | 01/26/15 | 01/29/15 20:40 | 150126L04 |
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 70 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

Page 2 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-01-1404-3-I | 01/22/15 00:35 | Aqueous | GC 40 | 01/26/15 | 01/29/15 21:02 | 150126L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 64 | 0-123 | |

| HCS 260 Lead | 15-01-1404-4-I | 01/22/15 01:10 | Aqueous | GC 40 | 01/26/15 | 01/29/15 21:26 | 150126L04 |
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 75 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-01-1404-5-I | 01/22/15 00:54 | Aqueous | GC 40 | 01/26/15 | 01/29/15 21:49 | 150126L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPD | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 63 | 0-123 | |

| Method Blank | 095-01-034-634 | N/A | Aqueous | GC 40 | 01/26/15 | 01/29/15 19:53 | 150126L04 |
|--------------|----------------|-----|---------|-------|----------|----------------|-----------|
|--------------|----------------|-----|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPD | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 60 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-01-1404-1-H | 01/22/15 00:36 | Aqueous | GC/MS TT | 01/24/15 | 02/02/15 21:03 | 150124L02B |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.8 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 9.8 | 2.8 | 1.00 | |
| Aniline | ND | 9.8 | 1.5 | 1.00 | |
| Anthracene | ND | 9.8 | 3.0 | 1.00 | |
| Azobenzene | ND | 9.8 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.4 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.8 | 4.6 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.8 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.8 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.8 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.8 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.8 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.8 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.8 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.8 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.8 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.8 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.8 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.8 | 2.6 | 1.00 | |
| Chrysene | ND | 9.8 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.8 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.8 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.8 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.8 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.8 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.8 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.8 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.8 | 3.0 | 1.00 | |
| Fluorene | ND | 9.8 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.8 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.8 | 1.00 | |
| Hexachloroethane | ND | 9.8 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.8 | 2.1 | 1.00 | |
| Isophorone | ND | 9.8 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.8 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.8 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.8 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.8 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 9.8 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 9.8 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.8 | 4.6 | 1.00 | |
| Phenanthrene | ND | 9.8 | 2.9 | 1.00 | |
| Phenol | ND | 9.8 | 2.0 | 1.00 | |
| Pyrene | ND | 9.8 | 2.9 | 1.00 | |
| Pyridine | ND | 9.8 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 80 | 33-120 | |
| 2-Fluorophenol | 55 | 24-120 | |
| Nitrobenzene-d5 | 74 | 38-120 | |
| p-Terphenyl-d14 | 78 | 41-137 | |
| Phenol-d6 | 35 | 16-120 | |
| 2,4,6-Tribromophenol | 105 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 4 of 18

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Lead | 15-01-1404-2-H | 01/22/15 00:51 | Aqueous | GC/MS TT | 01/24/15 | 02/02/15 21:22 | 150124L02B |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.8 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 9.8 | 2.8 | 1.00 | |
| Aniline | ND | 9.8 | 1.5 | 1.00 | |
| Anthracene | ND | 9.8 | 3.0 | 1.00 | |
| Azobenzene | ND | 9.8 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.4 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.8 | 4.6 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.8 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.8 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.8 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.8 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.8 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.8 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.8 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.8 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.8 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.8 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.8 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.8 | 2.6 | 1.00 | |
| Chrysene | ND | 9.8 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.8 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.8 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.8 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.8 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.8 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.8 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.8 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.8 | 3.0 | 1.00 | |
| Fluorene | ND | 9.8 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.8 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.8 | 1.00 | |
| Hexachloroethane | ND | 9.8 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.8 | 2.1 | 1.00 | |
| Isophorone | ND | 9.8 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.8 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.8 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.8 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.8 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 9.8 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 9.8 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.8 | 4.6 | 1.00 | |
| Phenanthrene | ND | 9.8 | 2.9 | 1.00 | |
| Phenol | ND | 9.8 | 2.0 | 1.00 | |
| Pyrene | ND | 9.8 | 2.9 | 1.00 | |
| Pyridine | ND | 9.8 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 71 | 33-120 | |
| 2-Fluorophenol | 54 | 24-120 | |
| Nitrobenzene-d5 | 69 | 38-120 | |
| p-Terphenyl-d14 | 69 | 41-137 | |
| Phenol-d6 | 33 | 16-120 | |
| 2,4,6-Tribromophenol | 90 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-01-1404-3-H | 01/22/15 00:35 | Aqueous | GC/MS TT | 01/24/15 | 02/02/15 21:40 | 150124L02B |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 76 | 33-120 | |
| 2-Fluorophenol | 54 | 24-120 | |
| Nitrobenzene-d5 | 73 | 38-120 | |
| p-Terphenyl-d14 | 75 | 41-137 | |
| Phenol-d6 | 34 | 16-120 | |
| 2,4,6-Tribromophenol | 99 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-01-1404-4-H | 01/22/15 01:10 | Aqueous | GC/MS TT | 01/24/15 | 02/02/15 21:59 | 150124L02B |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.8 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 9.8 | 2.8 | 1.00 | |
| Aniline | ND | 9.8 | 1.5 | 1.00 | |
| Anthracene | ND | 9.8 | 3.0 | 1.00 | |
| Azobenzene | ND | 9.8 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.4 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.8 | 4.6 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.8 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.8 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.8 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.8 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.8 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.8 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.8 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.8 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.8 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.8 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.8 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.8 | 2.6 | 1.00 | |
| Chrysene | ND | 9.8 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.8 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.8 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.8 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.8 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.8 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.8 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.8 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.8 | 3.0 | 1.00 | |
| Fluorene | ND | 9.8 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.8 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.8 | 1.00 | |
| Hexachloroethane | ND | 9.8 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.8 | 2.1 | 1.00 | |
| Isophorone | ND | 9.8 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.8 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.8 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.8 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.8 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 9.8 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 9.8 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.8 | 4.6 | 1.00 | |
| Phenanthrene | ND | 9.8 | 2.9 | 1.00 | |
| Phenol | ND | 9.8 | 2.0 | 1.00 | |
| Pyrene | ND | 9.8 | 2.9 | 1.00 | |
| Pyridine | ND | 9.8 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 67 | 33-120 | |
| 2-Fluorophenol | 51 | 24-120 | |
| Nitrobenzene-d5 | 64 | 38-120 | |
| p-Terphenyl-d14 | 67 | 41-137 | |
| Phenol-d6 | 32 | 16-120 | |
| 2,4,6-Tribromophenol | 86 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 13 of 18

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-01-1404-5-H | 01/22/15 00:54 | Aqueous | GC/MS TT | 01/24/15 | 02/02/15 22:18 | 150124L02B |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 75 | 33-120 | |
| 2-Fluorophenol | 54 | 24-120 | |
| Nitrobenzene-d5 | 72 | 38-120 | |
| p-Terphenyl-d14 | 75 | 41-137 | |
| Phenol-d6 | 33 | 16-120 | |
| 2,4,6-Tribromophenol | 100 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-3990 | N/A | Aqueous | GC/MS TT | 01/24/15 | 01/28/15 12:09 | 150124L02B |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 69 | 33-120 | |
| 2-Fluorophenol | 58 | 24-120 | |
| Nitrobenzene-d5 | 75 | 38-120 | |
| p-Terphenyl-d14 | 76 | 41-137 | |
| Phenol-d6 | 36 | 16-120 | |
| 2,4,6-Tribromophenol | 101 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 1 of 14

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-01-1404-1-A | 01/22/15 00:36 | Aqueous | GC/MS LL | 01/23/15 | 01/23/15 17:05 | 150123L019 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | 11 | 20 | 10 | 1.00 | J |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | 3.7 | 10 | 2.2 | 1.00 | J |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 102 | 78-126 | |
| 1,2-Dichloroethane-d4 | 111 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Lead | 15-01-1404-2-B | 01/22/15 00:51 | Aqueous | GC/MS JJ | 01/26/15 | 01/26/15 18:19 | 150126L019 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | 0.41 | 10 | 0.41 | 1.00 | J |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 99 | 80-120 | |
| Dibromofluoromethane | 94 | 78-126 | |
| 1,2-Dichloroethane-d4 | 101 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 5 of 14

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-01-1404-3-A | 01/22/15 00:35 | Aqueous | GC/MS LL | 01/23/15 | 01/23/15 20:50 | 150123L019 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 6 of 14

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 103 | 78-126 | |
| 1,2-Dichloroethane-d4 | 114 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 7 of 14

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-01-1404-4-A | 01/22/15 01:10 | Aqueous | GC/MS LL | 01/23/15 | 01/23/15 21:18 | 150123L019 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 99 | 80-120 | |
| Dibromofluoromethane | 103 | 78-126 | |
| 1,2-Dichloroethane-d4 | 114 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 9 of 14

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-01-1404-5-A | 01/22/15 00:54 | Aqueous | GC/MS LL | 01/23/15 | 01/23/15 21:46 | 150123L019 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 10 of 14

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 100 | 78-126 | |
| 1,2-Dichloroethane-d4 | 113 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 11 of 14

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-16235 | N/A | Aqueous | GC/MS LL | 01/23/15 | 01/23/15 16:37 | 150123L019 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 99 | 78-126 | |
| 1,2-Dichloroethane-d4 | 106 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-16243 | N/A | Aqueous | GC/MS JJ | 01/26/15 | 01/26/15 15:05 | 150126L019 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|----------|----------------|------------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| 1,4-Bromofluorobenzene | 100 | 80-120 | | | |
| Dibromofluoromethane | 94 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 98 | 75-135 | | | |
| Toluene-d8 | 100 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

01/23/15

Work Order:

15-01-1404

Page 1 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 210 Lead | 15-01-1404-1 | 01/22/15 00:36 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.059 | 0.10 | 0.025 | 1.00 | J | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Chloride (24) | 0.92 | 1.0 | 0.12 | 1.00 | J | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Nitrate (as N) (24) | 0.55 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Sulfate (24) | 3.1 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.10 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 41.0 | 1.00 | 0.848 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 41.0 | 1.00 | 0.848 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 65 | 1.0 | 0.82 | 1.00 | | mg/L | 01/26/15 | 01/26/15 | SM 2540 C |
| Solids, Total Suspended (24) | 20 | 1.0 | 0.95 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 D |
| pH (24) | 7.83 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/23/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.3 | 0.50 | 0.28 | 1.00 | | mg/L | 02/03/15 | 02/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 26 | 2.5 | 1.2 | 5.00 | | mg/L | 01/29/15 | 01/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 4.7 | 0.50 | 0.24 | 1.00 | | mg/L | 01/23/15 | 01/24/15 | SM 5310 B |

| | | | |
|---------------------|---------------------|-----------------------|----------------|
| HCS 240 Lead | 15-01-1404-2 | 01/22/15 00:51 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.20 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Chloride (24) | 16 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.7 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Sulfate (24) | 28 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.032 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 216 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 216 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 275 | 1.00 | 0.820 | 1.00 | | mg/L | 01/26/15 | 01/26/15 | SM 2540 C |
| Solids, Total Suspended (24) | 3.6 | 1.0 | 0.95 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 D |
| pH (24) | 7.68 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/23/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 02/03/15 | 02/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 11 | 2.5 | 1.2 | 5.00 | | mg/L | 01/29/15 | 01/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 4.9 | 0.50 | 0.24 | 1.00 | | mg/L | 01/23/15 | 01/24/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

01/23/15

Work Order:

15-01-1404

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 250 Lead | 15-01-1404-3 | 01/22/15 00:35 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.20 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Chloride (24) | 16 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Bromide (24) | 0.098 | 0.10 | 0.037 | 1.00 | J | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.6 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Sulfate (24) | 27 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.024 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 211 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 211 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 265 | 1.00 | 0.820 | 1.00 | | mg/L | 01/26/15 | 01/26/15 | SM 2540 C |
| Solids, Total Suspended (24) | 8.1 | 1.0 | 0.95 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 D |
| pH (24) | 7.79 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/23/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 02/03/15 | 02/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 8.6 | 2.5 | 1.2 | 5.00 | | mg/L | 01/29/15 | 01/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 7.2 | 0.50 | 0.24 | 1.00 | | mg/L | 01/23/15 | 01/24/15 | SM 5310 B |

| | | | |
|---------------------|---------------------|-----------------------|----------------|
| HCS 260 Lead | 15-01-1404-4 | 01/22/15 01:10 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.21 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Chloride (24) | 17 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Bromide (24) | 0.11 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.7 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Sulfate (24) | 33 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.037 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 222 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 222 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 275 | 1.00 | 0.820 | 1.00 | | mg/L | 01/26/15 | 01/26/15 | SM 2540 C |
| Solids, Total Suspended (24) | 11 | 1.0 | 0.95 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 D |
| pH (24) | 7.80 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/23/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.70 | 0.50 | 0.28 | 1.00 | | mg/L | 02/03/15 | 02/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.6 | 2.5 | 1.2 | 5.00 | | mg/L | 01/29/15 | 01/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 4.7 | 0.50 | 0.24 | 1.00 | | mg/L | 01/23/15 | 01/24/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

01/23/15

Work Order:

15-01-1404

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 270 Lead | 15-01-1404-5 | 01/22/15 00:54 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.21 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Chloride (24) | 16 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.6 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Sulfate (24) | 29 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.031 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 205 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 205 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | 265 | 1.00 | 0.820 | 1.00 | | mg/L | 01/26/15 | 01/26/15 | SM 2540 C |
| Solids, Total Suspended (24) | 13 | 1.0 | 0.95 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 D |
| pH (24) | 7.94 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/23/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.63 | 0.50 | 0.28 | 1.00 | | mg/L | 02/03/15 | 02/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 4.5 | 2.5 | 1.2 | 5.00 | | mg/L | 01/29/15 | 01/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 3.1 | 0.50 | 0.24 | 1.00 | | mg/L | 01/23/15 | 01/24/15 | SM 5310 B |

| | | |
|---------------------|------------|----------------|
| Method Blank | N/A | Aqueous |
|---------------------|------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|-------|---------------|---------------|-----------------|
| Fluoride (24) | ND | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Chloride (24) | ND | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Nitrate (as N) (24) | ND | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Sulfate (24) | ND | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/23/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/29/15 | SM 2320B |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.82 | 1.00 | | mg/L | 01/26/15 | 01/26/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.95 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 D |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 02/03/15 | 02/03/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 01/29/15 | 01/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 01/23/15 | 01/24/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-01-1370-4 | Sample | Aqueous | IC 15 | N/A | 01/23/15 16:52 | 150123S01 |
| 15-01-1370-4 | Matrix Spike | Aqueous | IC 15 | N/A | 01/23/15 18:17 | 150123S01 |
| 15-01-1370-4 | Matrix Spike Duplicate | Aqueous | IC 15 | N/A | 01/23/15 18:33 | 150123S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.1518 | 250.0 | 251.7 | 101 | 248.8 | 99 | 80-120 | 1 | 0-20 | |
| Chloride | 119.4 | 5000 | 5285 | 103 | 5293 | 103 | 80-120 | 0 | 0-20 | |
| Bromide | 0.2959 | 500.0 | 519.6 | 104 | 519.6 | 104 | 80-120 | 0 | 0-20 | |
| Nitrate (as N) | ND | 500.0 | 517.8 | 104 | 519.6 | 104 | 80-120 | 0 | 0-20 | |
| Sulfate | ND | 5000 | 5140 | 103 | 5149 | 103 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HCS 210 Lead | Sample | Aqueous | ACA 1 | N/A | 02/02/15 14:36 | 150202SO1 |
| HCS 210 Lead | Matrix Spike | Aqueous | ACA 1 | N/A | 02/02/15 14:36 | 150202SO1 |
| HCS 210 Lead | Matrix Spike Duplicate | Aqueous | ACA 1 | N/A | 02/02/15 14:36 | 150202SO1 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | 0.1013 | 0.2000 | 0.3235 | 111 | 0.3258 | 112 | 90-110 | 1 | 0-25 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HCS 210 Lead | Sample | Aqueous | TOC 8 | 01/29/15 | 01/30/15 08:20 | F0129TOCS3 |
| HCS 210 Lead | Matrix Spike | Aqueous | TOC 8 | 01/29/15 | 01/30/15 08:20 | F0129TOCS3 |
| HCS 210 Lead | Matrix Spike Duplicate | Aqueous | TOC 8 | 01/29/15 | 01/30/15 08:20 | F0129TOCS3 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Carbon, Total Organic | 25.90 | 50.00 | 59.00 | 66 | 61.00 | 70 | 31-145 | 3 | 0-23 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-01-1355-4 | Sample | Aqueous | TOC 8 | 01/23/15 | 01/24/15 06:25 | F0123DOCS2 |
| 15-01-1355-4 | Matrix Spike | Aqueous | TOC 8 | 01/23/15 | 01/24/15 06:25 | F0123DOCS2 |
| 15-01-1355-4 | Matrix Spike Duplicate | Aqueous | TOC 8 | 01/23/15 | 01/24/15 06:25 | F0123DOCS2 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|---------------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Carbon, Dissolved Organic | 29.40 | 50.00 | 68.00 | 77 | 68.00 | 77 | 31-145 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-01-1511-1 | Sample | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 17:55 | 150126SA6 |
| 15-01-1511-1 | Matrix Spike | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 17:57 | 150126SA6 |
| 15-01-1511-1 | Matrix Spike Duplicate | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 17:58 | 150126SA6 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Calcium | 19.19 | 0.5000 | 20.14 | 4X | 20.20 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 2.172 | 0.5000 | 2.657 | 4X | 2.674 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 1.560 | 5.000 | 6.465 | 98 | 6.337 | 96 | 83-131 | 2 | 0-7 | |
| Sodium | 1.042 | 5.000 | 5.732 | 94 | 5.907 | 97 | 73-127 | 3 | 0-9 | |
| Strontium | 0.09780 | 0.5000 | 0.6090 | 102 | 0.6223 | 105 | 81-123 | 2 | 0-6 | |
| Silicon | 1.065 | 0.5000 | 1.494 | 86 | 1.491 | 85 | 24-180 | 0 | 0-15 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
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San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|---------------------------|------------------------|-------------|----------|------------|---------------|----------------|---------------------|-----|--------|------------|
| 15-01-1511-2 | Sample | | Aqueous | ICP/MS 04 | 01/26/15 | 01/28/15 11:59 | 150126S03 | | | |
| 15-01-1511-2 | Matrix Spike | | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 08:25 | 150126S03 | | | |
| 15-01-1511-2 | Matrix Spike Duplicate | | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 08:28 | 150126S03 | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.09735 | 97 | 0.09348 | 93 | 85-133 | 4 | 0-11 | |
| Arsenic | ND | 0.1000 | 0.09587 | 96 | 0.09385 | 94 | 73-127 | 2 | 0-11 | |
| Barium | 0.05252 | 0.1000 | 0.1547 | 102 | 0.1527 | 100 | 74-128 | 1 | 0-10 | |
| Beryllium | ND | 0.1000 | 0.09530 | 95 | 0.09080 | 91 | 56-122 | 5 | 0-11 | |
| Cadmium | ND | 0.1000 | 0.09322 | 93 | 0.08931 | 89 | 84-114 | 4 | 0-8 | |
| Chromium | ND | 0.1000 | 0.1030 | 103 | 0.1001 | 100 | 73-133 | 3 | 0-11 | |
| Copper | 0.001256 | 0.1000 | 0.09476 | 94 | 0.09231 | 91 | 72-108 | 3 | 0-10 | |
| Lead | ND | 0.1000 | 0.1001 | 100 | 0.09724 | 97 | 79-121 | 3 | 0-10 | |
| Nickel | 0.001371 | 0.1000 | 0.09097 | 90 | 0.08788 | 87 | 68-122 | 3 | 0-10 | |
| Selenium | ND | 0.1000 | 0.09183 | 92 | 0.08674 | 87 | 59-125 | 6 | 0-12 | |
| Silver | ND | 0.05000 | 0.04271 | 85 | 0.03964 | 79 | 68-128 | 7 | 0-14 | |
| Thallium | ND | 0.1000 | 0.09514 | 95 | 0.09216 | 92 | 73-121 | 3 | 0-11 | |
| Zinc | 0.01878 | 0.1000 | 0.1226 | 104 | 0.09249 | 74 | 43-145 | 28 | 0-39 | |
| Aluminum | ND | 0.1000 | 0.1196 | 120 | 0.1191 | 119 | 47-161 | 0 | 0-24 | |
| Iron | ND | 5.100 | 5.315 | 104 | 5.347 | 105 | 27-201 | 1 | 0-24 | |
| Manganese | 0.001311 | 0.1000 | 0.09499 | 94 | 0.09083 | 90 | 72-126 | 4 | 0-42 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 7470A Total
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-01-1366-3 | Sample | Aqueous | Mercury 04 | 01/29/15 | 01/29/15 17:06 | 150129S06 |
| 15-01-1366-3 | Matrix Spike | Aqueous | Mercury 04 | 01/29/15 | 01/29/15 17:08 | 150129S06 |
| 15-01-1366-3 | Matrix Spike Duplicate | Aqueous | Mercury 04 | 01/29/15 | 01/29/15 17:10 | 150129S06 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | ND | 0.01000 | 0.01060 | 106 | 0.01080 | 108 | 66-126 | 2 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HCS 260 Lead | Sample | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 13:33 | 150129S19 |
| HCS 260 Lead | Matrix Spike | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 13:42 | 150129S19 |
| HCS 260 Lead | Matrix Spike Duplicate | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 17:15 | 150129S19 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Caffeine | 57.25 | 100.0 | 176.7 | 119 | 171.2 | 114 | 70-130 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|-----------------------------|------------------------|-------------|----------|------------|---------------|----------------|---------------------|-----|--------|------------|
| HCS 240 Lead | Sample | | Aqueous | GC/MS JJ | 01/26/15 | 01/26/15 18:19 | 150126S016 | | | |
| HCS 240 Lead | Matrix Spike | | Aqueous | GC/MS JJ | 01/26/15 | 01/26/15 18:47 | 150126S016 | | | |
| HCS 240 Lead | Matrix Spike Duplicate | | Aqueous | GC/MS JJ | 01/26/15 | 01/26/15 19:14 | 150126S016 | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 45.57 | 91 | 47.86 | 96 | 74-122 | 5 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 45.00 | 90 | 48.21 | 96 | 60-144 | 7 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 44.04 | 88 | 46.31 | 93 | 73-120 | 5 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 47.48 | 95 | 49.24 | 98 | 80-122 | 4 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 44.96 | 90 | 47.46 | 95 | 70-120 | 5 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 49.58 | 99 | 51.27 | 103 | 64-142 | 3 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 42.41 | 85 | 43.52 | 87 | 52-136 | 3 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 46.50 | 93 | 48.43 | 97 | 77-125 | 4 | 0-24 | |
| Toluene | ND | 50.00 | 45.90 | 92 | 48.23 | 96 | 72-126 | 5 | 0-23 | |
| Trichloroethene | ND | 50.00 | 46.78 | 94 | 48.76 | 98 | 74-128 | 4 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 33.56 | 67 | 35.26 | 71 | 67-133 | 5 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 92.06 | 92 | 96.04 | 96 | 63-129 | 4 | 0-25 | |
| o-Xylene | ND | 50.00 | 45.76 | 92 | 47.70 | 95 | 62-128 | 4 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 39.91 | 80 | 42.64 | 85 | 68-134 | 7 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|-----------------------------|------------------------|-------------|----------|------------|---------------|----------------|---------------------|-----|--------|------------|
| HCS 210 Lead | Sample | | Aqueous | GC/MS LL | 01/23/15 | 01/23/15 17:05 | 150123S015 | | | |
| HCS 210 Lead | Matrix Spike | | Aqueous | GC/MS LL | 01/23/15 | 01/23/15 17:33 | 150123S015 | | | |
| HCS 210 Lead | Matrix Spike Duplicate | | Aqueous | GC/MS LL | 01/23/15 | 01/23/15 18:01 | 150123S015 | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 52.69 | 105 | 58.35 | 117 | 74-122 | 10 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 50.79 | 102 | 58.11 | 116 | 60-144 | 13 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 50.97 | 102 | 56.51 | 113 | 73-120 | 10 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 50.85 | 102 | 57.73 | 115 | 80-122 | 13 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 50.55 | 101 | 56.50 | 113 | 70-120 | 11 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 54.87 | 110 | 60.90 | 122 | 64-142 | 10 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 50.92 | 102 | 55.97 | 112 | 52-136 | 9 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 52.12 | 104 | 58.45 | 117 | 77-125 | 11 | 0-24 | |
| Toluene | ND | 50.00 | 52.75 | 105 | 58.53 | 117 | 72-126 | 10 | 0-23 | |
| Trichloroethene | ND | 50.00 | 50.51 | 101 | 56.74 | 113 | 74-128 | 12 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 54.49 | 109 | 56.26 | 113 | 67-133 | 3 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 105.6 | 106 | 116.7 | 117 | 63-129 | 10 | 0-25 | |
| o-Xylene | ND | 50.00 | 52.20 | 104 | 57.76 | 116 | 62-128 | 10 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 49.98 | 100 | 55.50 | 111 | 68-134 | 10 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|--------------|-------------|------------|----------------|----------------|-----------------------|
| 15-01-1511-2 | Sample | Aqueous | ICP/MS 04 | 01/26/15 00:00 | 01/28/15 11:59 | 150126S03 |
| 15-01-1511-2 | PDS | Aqueous | ICP/MS 04 | 01/26/15 00:00 | 01/27/15 08:31 | 150126S03 |
| Parameter | Sample Conc. | Spike Added | PDS Conc. | PDS %Rec. | %Rec. CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.09422 | 94 | 75-125 | |
| Arsenic | ND | 0.1000 | 0.09373 | 94 | 75-125 | |
| Barium | 0.05252 | 0.1000 | 0.1449 | 92 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.09330 | 93 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.08992 | 90 | 75-125 | |
| Chromium | ND | 0.1000 | 0.1011 | 101 | 75-125 | |
| Copper | 0.001256 | 0.1000 | 0.09289 | 92 | 75-125 | |
| Lead | ND | 0.1000 | 0.09506 | 95 | 75-125 | |
| Nickel | 0.001371 | 0.1000 | 0.08826 | 87 | 75-125 | |
| Selenium | ND | 0.1000 | 0.08392 | 84 | 75-125 | |
| Silver | ND | 0.05000 | 0.04708 | 94 | 75-125 | |
| Thallium | ND | 0.1000 | 0.09009 | 90 | 75-125 | |
| Zinc | 0.01878 | 0.1000 | 0.1007 | 82 | 75-125 | |
| Aluminum | ND | 0.1000 | 0.1196 | 120 | 75-125 | |
| Iron | ND | 5.100 | 4.713 | 92 | 75-125 | |
| Manganese | 0.001311 | 0.1000 | 0.08988 | 89 | 75-125 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-01-1355-3 | Sample | Aqueous | PH1/BUR03 | N/A | 01/29/15 16:10 | F0129ALKD1 |
| 15-01-1355-3 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 01/29/15 16:10 | F0129ALKD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|---|---------------------|------------------|------------|---------------|-------------------|
| Alkalinity, Total (as CaCO ₃) | 440.0 | 438.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-01-1355-3 | Sample | Aqueous | PH1/BUR03 | N/A | 01/29/15 16:10 | F0129HCOD1 |
| 15-01-1355-3 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 01/29/15 16:10 | F0129HCOD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Bicarbonate (as CaCO ₃) | 440.0 | 438.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-01-1355-3 | Sample | Aqueous | PH1/BUR03 | N/A | 01/29/15 16:10 | F0129CO3D1 |
| 15-01-1355-3 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 01/29/15 16:10 | F0129CO3D1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-01-1447-2 | Sample | Aqueous | SC 5 | 01/26/15 00:00 | 01/26/15 16:00 | F0126TDSD1 |
| 15-01-1447-2 | Sample Duplicate | Aqueous | SC 5 | 01/26/15 00:00 | 01/26/15 16:00 | F0126TDSD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Dissolved | 630.0 | 670.0 | 6 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-01-1544-1 | Sample | Aqueous | N/A | 01/27/15 00:00 | 01/27/15 14:00 | F0127TSSD1 |
| 15-01-1544-1 | Sample Duplicate | Aqueous | N/A | 01/27/15 00:00 | 01/27/15 14:00 | F0127TSSD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Suspended | 26.80 | 26.00 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-01-1394-1 | Sample | Aqueous | PH 1 | N/A | 01/23/15 18:45 | F0123PHD1 |
| 15-01-1394-1 | Sample Duplicate | Aqueous | PH 1 | N/A | 01/23/15 18:45 | F0123PHD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|------------------|---------------------|------------------|------------|---------------|-------------------|
| pH | 6.670 | 6.680 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 4500 N Org B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-01-1729-2 | Sample | Aqueous | BUR05 | 02/03/15 00:00 | 02/03/15 17:24 | F0203TKND1 |
| 15-01-1729-2 | Sample Duplicate | Aqueous | BUR05 | 02/03/15 00:00 | 02/03/15 17:24 | F0203TKND1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Total Kjeldahl Nitrogen | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-12-906-5371 | LCS | Aqueous | IC 15 | N/A | 01/23/15 11:43 | 150123L01 |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Fluoride | | 2.500 | 2.463 | 99 | 90-110 | |
| Chloride | | 50.00 | 49.50 | 99 | 90-110 | |
| Bromide | | 5.000 | 5.048 | 101 | 90-110 | |
| Nitrate (as N) | | 5.000 | 5.108 | 102 | 90-110 | |
| Sulfate | | 50.00 | 50.06 | 100 | 90-110 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------------|------------|---------------|-------------------|
| 099-12-739-97 | LCS | Aqueous | ACA 1 | N/A | 02/02/15 14:36 | 150202LO1 | | | |
| 099-12-739-97 | LCSD | Aqueous | ACA 1 | N/A | 02/02/15 14:36 | 150202LO1 | | | |
| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Phosphorus, Total | 0.2000 | 0.1852 | 93 | 0.1862 | 93 | 90-110 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-859-575 | LCS | Aqueous | PH1/BUR03 | N/A | 01/29/15 16:10 | F0129ALKB1 | | | |
| 099-15-859-575 | LCSD | Aqueous | PH1/BUR03 | N/A | 01/29/15 16:10 | F0129ALKB1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 100.0 | 97.00 | 97 | 97.00 | 97 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4400 | LCS | Aqueous | SC 5 | 01/26/15 | 01/26/15 16:00 | F0126TDSL1 | | | |
| 099-12-180-4400 | LCSD | Aqueous | SC 5 | 01/26/15 | 01/26/15 16:00 | F0126TDSL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 85.00 | 85 | 90.00 | 90 | 80-120 | 6 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-09-010-7018 | LCS | Aqueous | N/A | 01/27/15 | 01/27/15 14:00 | F0127TSSL1 | | | |
| 099-09-010-7018 | LCSD | Aqueous | N/A | 01/27/15 | 01/27/15 14:00 | F0127TSSL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Suspended | 100.0 | 85.00 | 85 | 91.00 | 91 | 80-120 | 7 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 099-05-097-5503 | LCS | Aqueous | TOC 8 | 01/29/15 | 01/30/15 08:20 | F0129TOCL3 |
| 099-05-097-5503 | LCSD | Aqueous | TOC 8 | 01/29/15 | 01/30/15 08:20 | F0129TOCL3 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| Carbon, Total Organic | 10.00 | 10.50 | 105 | 10.80 | 108 | 80-120 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-115-1397 | LCS | Aqueous | TOC 8 | 01/23/15 | 01/24/15 07:33 | F0123DOCL2 | | | |
| 099-05-115-1397 | LCSD | Aqueous | TOC 8 | 01/23/15 | 01/24/15 07:33 | F0123DOCL2 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Dissolved Organic | 10.00 | 10.40 | 104 | 10.70 | 107 | 80-120 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|--------------------|------------------------|------------------|-----------------|-------------------|
| 099-15-683-1120 | LCS | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 17:53 | 150126L6FF |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Calcium | | 0.5000 | 0.4279 | 86 | 80-120 | |
| Magnesium | | 0.5000 | 0.4097 | 82 | 80-120 | |
| Potassium | | 5.000 | 4.428 | 89 | 80-120 | |
| Sodium | | 5.000 | 4.045 | 81 | 80-120 | |
| Strontium | | 0.5000 | 0.4624 | 92 | 80-120 | |
| Silicon | | 0.5000 | 0.2593 | 52 | 80-120 | X |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-719 | LCS | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 15:23 | 150126L03F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.1029 | 103 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.1052 | 105 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.09971 | 100 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.1018 | 102 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.1012 | 101 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.1053 | 105 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.09728 | 97 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.1023 | 102 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.1020 | 102 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.1022 | 102 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.05273 | 105 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.09673 | 97 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.1070 | 107 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.1175 | 118 | 80-120 | 73-127 | |
| Iron | 5.100 | 5.029 | 99 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.1022 | 102 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|------------------|
| 099-15-763-484 | LCS | Aqueous | Mercury 04 | 01/29/15 | 01/29/15 17:03 | 150129L06F |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.01000 | 0.01066 | 107 | 85-121 | |

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Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 099-16-376-10 | LCS | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 12:55 | 150129L19 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Caffeine | | 100.0 | 103.8 | 104 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-529-778 | LCS | Aqueous | | GC 51 | 01/24/15 | 02/04/15 04:22 | 150124L08A | | | |
| 099-12-529-778 | LCSD | Aqueous | | GC 51 | 01/24/15 | 02/04/15 04:36 | 150124L08A | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Alpha-BHC | 0.5000 | 0.5324 | 106 | 0.5460 | 109 | 50-135 | 36-149 | 3 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.5466 | 109 | 0.5590 | 112 | 50-135 | 36-149 | 2 | 0-25 | |
| Beta-BHC | 0.5000 | 0.4827 | 97 | 0.4746 | 95 | 50-135 | 36-149 | 2 | 0-25 | |
| Heptachlor | 0.5000 | 0.5339 | 107 | 0.5402 | 108 | 50-135 | 36-149 | 1 | 0-25 | |
| Delta-BHC | 0.5000 | 0.6051 | 121 | 0.6324 | 126 | 50-135 | 36-149 | 4 | 0-25 | |
| Aldrin | 0.5000 | 0.4714 | 94 | 0.4712 | 94 | 50-135 | 36-149 | 0 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.5246 | 105 | 0.5276 | 106 | 50-135 | 36-149 | 1 | 0-25 | |
| Endosulfan I | 0.5000 | 0.5515 | 110 | 0.5555 | 111 | 50-135 | 36-149 | 1 | 0-25 | |
| Dieldrin | 0.5000 | 0.5378 | 108 | 0.5415 | 108 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.4811 | 96 | 0.4794 | 96 | 50-135 | 36-149 | 0 | 0-25 | |
| Endrin | 0.5000 | 0.5335 | 107 | 0.5371 | 107 | 50-135 | 36-149 | 1 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.6733 | 135 | 0.6103 | 122 | 50-135 | 36-149 | 10 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.5057 | 101 | 0.5046 | 101 | 50-135 | 36-149 | 0 | 0-25 | |
| Endosulfan II | 0.5000 | 0.5349 | 107 | 0.5386 | 108 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.5675 | 114 | 0.5887 | 118 | 50-135 | 36-149 | 4 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.5345 | 107 | 0.5392 | 108 | 50-135 | 36-149 | 1 | 0-25 | |
| Methoxychlor | 0.5000 | 0.5191 | 104 | 0.5246 | 105 | 50-135 | 36-149 | 1 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------------|------------|---------------|-------------------|
| 099-12-533-1000 | LCS | Aqueous | GC 31 | 01/24/15 | 02/04/15 10:25 | 150124L09 | | | |
| 099-12-533-1000 | LCSD | Aqueous | GC 31 | 01/24/15 | 02/04/15 10:44 | 150124L09 | | | |
| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Aroclor-1016 | 2.000 | 1.741 | 87 | 1.763 | 88 | 50-135 | 1 | 0-25 | |
| Aroclor-1260 | 2.000 | 1.654 | 83 | 1.659 | 83 | 50-135 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-963-82 | LCS | Aqueous | | GC 26 | 01/24/15 | 02/06/15 19:25 | 150124L03 | | | |
| 099-15-963-82 | LCSD | Aqueous | | GC 26 | 01/24/15 | 02/06/15 20:09 | 150124L03 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.04413 | 110 | 0.04395 | 110 | 30-130 | 13-147 | 0 | 0-30 | |
| Bolstar | 0.04000 | 0.04712 | 118 | 0.04669 | 117 | 30-130 | 13-147 | 1 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.04763 | 119 | 0.04691 | 117 | 30-130 | 13-147 | 2 | 0-30 | |
| Coumaphos | 0.04000 | 0.04735 | 118 | 0.04833 | 121 | 30-130 | 13-147 | 2 | 0-30 | |
| Diazinon | 0.04000 | 0.04995 | 125 | 0.04985 | 125 | 30-130 | 13-147 | 0 | 0-30 | |
| Disulfoton | 0.04000 | 0.04472 | 112 | 0.04335 | 108 | 30-130 | 13-147 | 3 | 0-30 | |
| Ethoprop | 0.04000 | 0.04985 | 125 | 0.04905 | 123 | 30-130 | 13-147 | 2 | 0-30 | |
| Fensulfothion | 0.04000 | 0.04732 | 118 | 0.04644 | 116 | 30-130 | 13-147 | 2 | 0-30 | |
| Fenthion | 0.04000 | 0.04798 | 120 | 0.04733 | 118 | 30-130 | 13-147 | 1 | 0-30 | |
| Merphos | 0.04000 | 0.06099 | 152 | 0.05996 | 150 | 30-130 | 13-147 | 2 | 0-30 | X |
| Methyl Parathion | 0.04000 | 0.04832 | 121 | 0.04758 | 119 | 30-130 | 13-147 | 2 | 0-30 | |
| Phorate | 0.04000 | 0.04887 | 122 | 0.04843 | 121 | 30-130 | 13-147 | 1 | 0-30 | |
| Ronnel | 0.04000 | 0.04912 | 123 | 0.04823 | 121 | 30-130 | 13-147 | 2 | 0-30 | |
| Stirophos | 0.04000 | 0.04117 | 103 | 0.04067 | 102 | 30-130 | 13-147 | 1 | 0-30 | |
| Tokuthion | 0.04000 | 0.04646 | 116 | 0.04588 | 115 | 30-130 | 13-147 | 1 | 0-30 | |
| Trichloronate | 0.04000 | 0.04793 | 120 | 0.04745 | 119 | 30-130 | 13-147 | 1 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

Page 15 of 18

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 095-01-034-634 | LCS | Aqueous | GC 40 | 01/26/15 | 01/29/15 19:07 | 150126L04 |
| 095-01-034-634 | LCSD | Aqueous | GC 40 | 01/26/15 | 01/29/15 19:30 | 150126L04 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| 2,4-D | 20.00 | 13.92 | 70 | 12.66 | 63 | 30-130 | 10 | 0-30 | |
| 2,4,5-T | 2.000 | 1.485 | 74 | 1.365 | 68 | 30-130 | 8 | 0-30 | |
| 2,4-DB | 20.00 | 13.54 | 68 | 12.15 | 61 | 30-130 | 11 | 0-30 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-3990 | LCS | Aqueous | | GC/MS TT | 01/24/15 | 01/28/15 13:24 | 150124L02B | | | |
| 095-01-003-3990 | LCSD | Aqueous | | GC/MS TT | 01/24/15 | 01/28/15 20:49 | 150124L02B | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 200.0 | 155.6 | 78 | 156.2 | 78 | 61-120 | 51-130 | 0 | 0-20 | |
| Acenaphthylene | 200.0 | 149.3 | 75 | 150.2 | 75 | 55-120 | 44-131 | 1 | 0-20 | |
| Butyl Benzyl Phthalate | 200.0 | 150.3 | 75 | 149.1 | 75 | 56-122 | 45-133 | 1 | 0-20 | |
| 4-Chloro-3-Methylphenol | 200.0 | 137.6 | 69 | 140.8 | 70 | 52-120 | 41-131 | 2 | 0-20 | |
| 2-Chlorophenol | 200.0 | 147.0 | 74 | 145.8 | 73 | 47-120 | 35-132 | 1 | 0-20 | |
| 1,4-Dichlorobenzene | 200.0 | 148.8 | 74 | 149.0 | 75 | 36-120 | 22-134 | 0 | 0-20 | |
| Dimethyl Phthalate | 200.0 | 149.3 | 75 | 153.3 | 77 | 60-120 | 50-130 | 3 | 0-20 | |
| 2,4-Dinitrotoluene | 200.0 | 152.6 | 76 | 158.6 | 79 | 61-121 | 51-131 | 4 | 0-20 | |
| Fluorene | 200.0 | 158.6 | 79 | 163.1 | 82 | 67-120 | 58-129 | 3 | 0-20 | |
| N-Nitroso-di-n-propylamine | 200.0 | 131.5 | 66 | 130.6 | 65 | 39-123 | 25-137 | 1 | 0-20 | |
| Naphthalene | 200.0 | 151.2 | 76 | 152.8 | 76 | 54-120 | 43-131 | 1 | 0-20 | |
| 4-Nitrophenol | 200.0 | 65.62 | 33 | 66.21 | 33 | 14-120 | 0-138 | 1 | 0-20 | |
| Pentachlorophenol | 200.0 | 139.5 | 70 | 140.5 | 70 | 31-127 | 15-143 | 1 | 0-20 | |
| Phenol | 200.0 | 65.48 | 33 | 65.97 | 33 | 17-120 | 0-137 | 1 | 0-20 | |
| Pyrene | 200.0 | 144.9 | 72 | 147.3 | 74 | 58-124 | 47-135 | 2 | 0-20 | |
| 1,2,4-Trichlorobenzene | 200.0 | 148.3 | 74 | 151.5 | 76 | 49-120 | 37-132 | 2 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-16243 | LCS | Aqueous | GC/MS JJ | 01/26/15 | 01/26/15 14:09 | 150126L019 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 47.84 | 96 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 48.61 | 97 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 46.78 | 94 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 48.32 | 97 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 47.18 | 94 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 49.34 | 99 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 42.64 | 85 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 48.45 | 97 | 80-123 | 73-130 | |
| Toluene | 50.00 | 48.10 | 96 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 50.19 | 100 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 36.64 | 73 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 96.66 | 97 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 47.41 | 95 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 39.65 | 79 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/23/15
Work Order: 15-01-1404
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-16235 | LCS | Aqueous | GC/MS LL | 01/23/15 | 01/23/15 15:13 | 150123L019 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 48.75 | 97 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 46.96 | 94 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 47.58 | 95 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 48.06 | 96 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 46.98 | 94 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 50.76 | 102 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 46.07 | 92 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 48.04 | 96 | 80-123 | 73-130 | |
| Toluene | 50.00 | 48.71 | 97 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 46.15 | 92 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 47.38 | 95 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 95.69 | 96 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 48.05 | 96 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 47.30 | 95 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Sample Analysis Summary Report

Work Order: 15-01-1404

Page 1 of 1

| <u>Method</u> | <u>Extraction</u> | <u>Chemist ID</u> | <u>Instrument</u> | <u>Analytical Location</u> |
|-----------------------|-------------------|-------------------|-------------------|----------------------------|
| EPA 1694 (M) Caffeine | EPA 1694 | 262 | LC/TQ 2 | 1 |
| EPA 300.0 | N/A | 650 | IC 15 | 1 |
| EPA 365.1 | N/A | 735 | ACA 1 | 1 |
| EPA 6010B | EPA 3005A Filt. | 771 | ICP 7300 | 1 |
| EPA 6020 | EPA 3005A Filt. | 776 | ICP/MS 03 | 1 |
| EPA 7470A | EPA 7470A Filt. | 915 | Mercury 04 | 1 |
| EPA 8081A | EPA 3510C | 669 | GC 51 | 1 |
| EPA 8082 | EPA 3510C | 944 | GC 31 | 1 |
| EPA 8141A | EPA 3510C | 949 | GC 26 | 1 |
| EPA 8151A | EPA 8151A | 944 | GC 40 | 1 |
| EPA 8260B | EPA 5030C | 959 | GC/MS JJ | 2 |
| EPA 8260B | EPA 5030C | 959 | GC/MS LL | 2 |
| EPA 8270C | EPA 3510C | 923 | GC/MS TT | 1 |
| SM 2320B | N/A | 885 | PH1/BUR03 | 1 |
| SM 2540 C | N/A | 689 | SC 5 | 1 |
| SM 2540 D | N/A | 689 | N/A | 1 |
| SM 4500 H+ B | N/A | 688 | PH 1 | 1 |
| SM 4500 N Org B | N/A | 685 | BUR05 | 1 |
| SM 5310 B | N/A | 735 | TOC 8 | 1 |

Return to Contents

Location 1: 7440 Lincoln Way, Garden Grove, CA 92841

Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841

Glossary of Terms and Qualifiers

Work Order: 15-01-1404

Page 1 of 1

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|--|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDS or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.



Calscience

7440 Lincoln Way, Garden Grove, CA 92641-1427 • (714) 895-5494
For courier service / sample drop off information, contact us26_sales@eurofinsus.com or call us.

CHAIN OF CUSTODY RECORD

WO# / LAB USE ONLY
15-01-1404

DATE: 1/22/15
PAGE: 1 OF 3

| | | | | | |
|---|---------------------------|---|--|--|--|
| LABORATORY CLIENT: SWCA Environmental Consultants | | CLIENT PROJECT NAME / NUMBER: EAA 27122 | | P.O. NO.: | |
| ADDRESS: 6200 UTSA Blvd. Suite 102 | | PROJECT CONTACT: Philip Pearce | | SAMPLER(S) (PRINT) Philip Pearce | |
| CITY: San Antonio | STATE: TX | ZIP: 78249-1618 | | | |
| TEL: 210.877.2847 | E-MAIL: P Pearce@swca.com | | | | |

| | | | | | | | | |
|---|--------------|---------------|-------|------------------------------------|--------------|-------------|-----------|----------------|
| TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD"): | | | | LOG CODE: | | | | |
| <input type="checkbox"/> SAME DAY <input type="checkbox"/> 24 HR <input type="checkbox"/> 48 HR <input type="checkbox"/> 72 HR <input checked="" type="checkbox"/> 10 Days (standard) | | | | <input type="checkbox"/> COELT EDF | | | | |
| GLOBAL ID: | | | | SPECIAL INSTRUCTIONS: | | | | |
| Samples for EPA 6010B, EPA 6020, and EPA 7470 Metals are field filtered EPA 6020 Metals: Zn, Ti, Se, Sb, Pb, Ni, Cu, Cr, Cd, Be, Ba, As, Al, Ag, Fe, Mn ★ Please also analyze each sample for pH. ★ | | | | | | | | |
| LAB. USE ONLY | SAMPLE ID | SAMPLING DATE | TIME | MATRIX | NO. OF CONT. | Unpreserved | Preserved | Field Filtered |
| 1 | HCS 210 Lead | 1/22/15 | 00:36 | NPW | 9 | 2 | 7 | 1 |
| 2 | HCS 240 Lead | 1/22/15 | 00:51 | NPW | 9 | 2 | 7 | 1 |
| 3 | HCS 250 Lead | 1/22/15 | 00:35 | NPW | 9 | 2 | 7 | 1 |
| 4 | HCS 260 Lead | 1/22/15 | 1:10 | NPW | 9 | 2 | 7 | 1 |
| 5 | HCS 270 Lead | 1/22/15 | 00:54 | NPW | 9 | 2 | 7 | 1 |

| | | | |
|------------------------------|--------------------------------------|---------------|-------------|
| Relinquished by: (Signature) | Received by: (Signature/Affiliation) | Date: 1/22/15 | Time: 09:13 |
| Relinquished by: (Signature) | Received by: (Signature/Affiliation) | Date: | Time: |
| Relinquished by: (Signature) | Received by: (Signature/Affiliation) | Date: 1/23/15 | Time: 1000 |



SWCA INC
6200 UTSA BLVD STE 102
SAN ANTONIO, TX 782491618
UNITED STATES US

ACTWGT: 54.2 LB
CAD: /POS1525
DIMS: 23x13x13 IN
BILL SENDER

SWCA INC
6200 UTSA BLVD STE 102
SAN ANTONIO, TX 782491618
UNITED STATES US

ACTWGT: 54.2 LB
CAD: /POS1525
DIMS: 23x13x13 IN
BILL SENDER

CAL SCIENCE
7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895-5494

REF:

DEPT:

TO

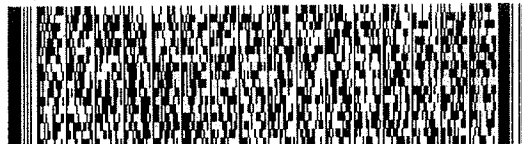
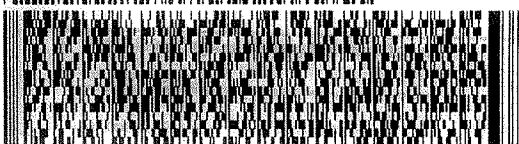
CAL SCIENCE
7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895-5494

REF:

DEPT:



1 of 5
TRK# 8066 6158 0190
0200
MASTER

FRI - 23 JAN AA
STANDARD OVERNIGHT

A7 APVA

92841
CA-US SNA

2 of 5
MPS# 7801 6964 1730
0681
Mstr# 8066 6158 0190

FRI - 23 JAN AA
STANDARD OVERNIGHT

A7 APVA

92841
CA-US SNA



ORIGIN ID:NIRA (210) 877-2847
SWCA INC

6200 UTSA BLVD STE 102

SAN ANTONIO, TX 782491618
UNITED STATES US

P DATE: 22JAN15
ACTWGT: 53.0 LB
ID: /POS1525
DIMS: 23x13x13 IN

JILL SENDER

ORIGIN ID:NIRA (210) 877-2847
SWCA INC

6200 UTSA BLVD STE 102

SAN ANTONIO, TX 782491618
UNITED STATES US

SHIP DATE: 22JAN15
ACTWGT: 51.1 LB
CAD: /POS1525
DIMS: 23x13x13 IN
BILL SENDER

TO

CAL SCIENCE
7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895-5494

REF:

DEPT:

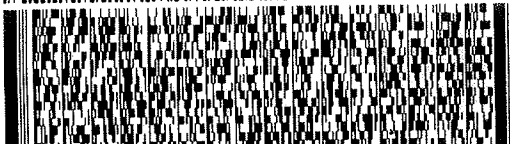
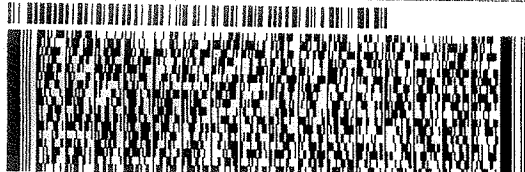
CAL SCIENCE
7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895-5494

REF:

DEPT:



3 of 5
MPS# 7801 6964 1741
0681
Mstr# 8066 6158 0190

FRI - 23 JAN AA
STANDARD OVERNIGHT

A7 APVA

92841
CA-US SNA

4 of 5
MPS# 7801 6964 1752
0681
Mstr# 8066 6158 0190

FRI - 23 JAN AA
STANDARD OVERNIGHT

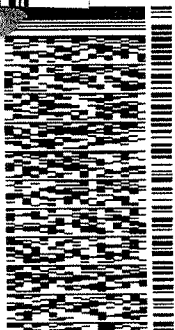
A7 APVA

92841
CA-US SNA



A7 APVA

5 of 5
MPS# 7801 6964 1763
0681
Mstr# 8066 6158 0190



GARDEN GROVE CA 92841

CAL SCIENCE
7440 LINCOLN WAY

TO

ORIGIN ID:NIRA (210) 877-2847
SWCA INC
6200 UTSA BLVD STE 102
SAN ANTONIO, TX 782491618
UNITED STATES US

Calscience

WORK ORDER #: 15-01-1404

SAMPLE RECEIPT FORM

Cooler 1 of 5

CLIENT: SWCA

DATE: 01/23/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 1.7 °C + 0.2 °C (CF) = 1.9 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 15

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 619

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|--------------------------|-------------------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
|--|--------------------------|-------------------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☒ ☐

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBznna ☐ 100PJ ☐ 100PJna₂ ☒ 2.5 gal. cube ☐ _____ ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 619

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 619

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure znna: ZnAc₂+NaOH f: Filtered Scanned by: 619

Calscience

WORK ORDER #: 15-01-1404

SAMPLE RECEIPT FORM

Cooler 2 of 5

CLIENT: SWCA

DATE: 01/23/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Temperature 2.4 °C + 0.2°C (CF) = 2.6 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 15

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 659

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☒ ☐

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBz₂na ☐ 100PJ ☐ 100PJna₂ ☒ 2.5 gal cube ☐ _____ ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 659

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 8m

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure znna: ZnAc₂+NaOH f: Filtered Scanned by: 8m

Calscience

WORK ORDER #: 15-01-1404

SAMPLE RECEIPT FORM

Cooler 3 of 5

CLIENT: SWCA

DATE: 01/23/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 2.7 °C + 0.2 °C (CF) = 2.9 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 15

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 659

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

| | | | |
|--|--------------------------|-------------------------------------|--------------------------|
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfides <input type="checkbox"/> Dissolved Oxygen..... | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
|--|--------------------------|-------------------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBznna ☐ 100PJ ☐ 100PJna₂ ☒ 2.5 gal cube ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 659

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 846

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure znna: ZnAc₂+NaOH f: Filtered Scanned by: 846

Calscience

WORK ORDER #: 15-01-1404

SAMPLE RECEIPT FORM

Cooler 4 of 5

CLIENT: SWCA

DATE: 01/23/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 2.0 °C + 0.2 °C (CF) = 2.2 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 15

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 689

SAMPLE CONDITION:

Yes No N/A

Chain-Of-Custody (COC) document(s) received with samples..... ☒ ☐ ☐

COC document(s) received complete..... ☒ ☐ ☐
☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

Sampler's name indicated on COC..... ☒ ☐ ☐

Sample container label(s) consistent with COC..... ☒ ☐ ☐

Sample container(s) intact and good condition..... ☒ ☐ ☐

Proper containers and sufficient volume for analyses requested..... ☒ ☐ ☐

Analyses received within holding time..... ☒ ☐ ☐

Aqueous samples received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☒ ☐

Proper preservation noted on COC or sample container..... ☒ ☐ ☐
☐ Unpreserved vials received for Volatiles analysis

Volatile analysis container(s) free of headspace..... ☒ ☐ ☐

Tedlar bag(s) free of condensation..... ☐ ☐ ☒

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOA³h ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ² ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBznna ☐ 100PJ ☐ 100PJna₂ ☒ ^{2.5 gnl} cube ☐ _____ ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 689

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 826

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure znna: ZnAc₂+NaOH f: Filtered Scanned by: 826

Calscience

WORK ORDER #: 15-01-1404

SAMPLE RECEIPT FORM

Cooler 5 of 5

CLIENT: SWCA

DATE: 01/23/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 2.3 °C + 0.2 °C (CF) = 2.5 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 15

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 659

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☒ ☐

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBz₂na ☐ 100PJ ☐ 100PJna₂ ☒ 2.5 gal cube ☐ _____ ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 659

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 846

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure z₂na: ZnAc₂+NaOH f: Filtered Scanned by: 846



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WORK ORDER #: 15-01-1404

SAMPLE ANOMALY FORM

SAMPLES - CONTAINERS & LABELS:

Comments:

- ☐ Sample(s) NOT RECEIVED but listed on COC
☐ Sample(s) received but NOT LISTED on COC
☒ Holding time expired – list sample ID(s) and test
☐ Insufficient quantities for analysis – list test
☐ Improper container(s) used – list test
☐ Improper preservative used – list test
☐ No preservative noted on COC or label – list test & notify lab
☐ Sample labels illegible – note test/container type
☒ Sample label(s) do not match COC – Note in comments
 - ☐ Sample ID
 - ☒ Date and/or Time Collected
 - ☐ Project Information
 - ☐ # of Container(s)
 - ☐ Analysis☐ Sample container(s) compromised – Note in comments
 - ☐ Water present in sample container
 - ☐ Broken☐ Sample container(s) not labeled
☐ Air sample container(s) compromised – Note in comments
 - ☐ Flat
 - ☐ Very low in volume
 - ☐ Leaking (Not transferred - duplicate bag submitted)
 - ☐ Leaking (transferred into Calscience Tedlar® Bag*)
 - ☐ Leaking (transferred into Client's Tedlar® Bag*)☐ Other: _____

(-1) through (-5) pH expired.

(-4) Collection time per label is 00:10

HEADSPACE – Containers with Bubble > 6mm or ¼ inch:

| Sample # | Container ID(s) | # of Vials Received | Sample # | Container ID(s) | # of Vials Received | Sample # | Container ID(s) | # of Cont. received | Analysis |
|----------|-----------------|---------------------|----------|-----------------|---------------------|----------|-----------------|---------------------|----------|
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |

Comments: _____

*Transferred at Client's request.

Initial / Date: 6/3 01/23/15



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WORK ORDER NUMBER: 15-01-1511

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: SWCA Environmental Consultants

Client Project Name: EAA 27122

Attention: Philip Pearce
6200 UTSA Blvd.
Suite 102
San Antonio, TX 78249-1618

Approved for release on 02/10/2015 by:
Don Burley
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

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Work Order Number: 15-01-1511

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Work Order Narrative

Work Order: 15-01-1511

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Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 01/24/15. They were assigned to Work Order 15-01-1511.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

EPA 6010B Silicon LCS recovery was below the control limits. Silicon was detected in each of the samples. The results may be biased low.



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Sample Summary

| | | |
|--|-----------------------|----------------|
| Client: SWCA Environmental Consultants | Work Order: | 15-01-1511 |
| 6200 UTSA Blvd., Suite 102 | Project Name: | EAA 27122 |
| San Antonio, TX 78249-1618 | PO Number: | |
| | Date/Time Received: | 01/24/15 10:00 |
| | Number of Containers: | 109 |
| Attn: Philip Pearce | | |

| Sample Identification | Lab Number | Collection Date and Time | Number of Containers | Matrix |
|-----------------------|---------------|--------------------------|----------------------|---------|
| HCS 210 Peak | 15-01-1511-1 | 01/22/15 16:28 | 9 | Aqueous |
| HCS 240 Peak | 15-01-1511-2 | 01/22/15 16:51 | 9 | Aqueous |
| HCS 250 Peak | 15-01-1511-3 | 01/22/15 16:40 | 9 | Aqueous |
| HCS 260 Peak | 15-01-1511-4 | 01/22/15 17:03 | 9 | Aqueous |
| HCS 270 Peak | 15-01-1511-5 | 01/22/15 16:55 | 9 | Aqueous |
| HCS 210 Trail | 15-01-1511-6 | 01/23/15 03:01 | 9 | Aqueous |
| HCS 240 Trail | 15-01-1511-7 | 01/23/15 03:21 | 9 | Aqueous |
| HCS 250 Trail | 15-01-1511-8 | 01/23/15 03:03 | 9 | Aqueous |
| HCS 260 Trail | 15-01-1511-9 | 01/23/15 03:36 | 9 | Aqueous |
| HCS 270 Trail | 15-01-1511-10 | 01/23/15 03:30 | 9 | Aqueous |
| FDHCS 260 Trail | 15-01-1511-11 | 01/23/15 03:36 | 9 | Aqueous |
| FDHCS 270 Trail | 15-01-1511-12 | 01/23/15 03:30 | 9 | Aqueous |
| TB01 | 15-01-1511-13 | 01/23/15 00:00 | 1 | Aqueous |

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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HCS 210 Peak | 15-01-1511-1-D | 01/22/15 16:28 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 17:55 | 150126L6FF |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 19.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 2.17 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.56 | 0.500 | 0.103 | 1.00 | |
| Sodium | 1.04 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.0978 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 1.07 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HCS 240 Peak | 15-01-1511-2-D | 01/22/15 16:51 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:00 | 150126L6FF |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 75.6 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 14.6 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.50 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.7 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.642 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.45 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HCS 250 Peak | 15-01-1511-3-D | 01/22/15 16:40 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:01 | 150126L6FF |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 70.9 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 13.8 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.68 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.0 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.597 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.24 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HCS 260 Peak | 15-01-1511-4-D | 01/22/15 17:03 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:09 | 150126L6FF |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 47.0 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 5.45 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 5.26 | 0.500 | 0.103 | 1.00 | |
| Sodium | 9.44 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.262 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 6.77 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HCS 270 Peak | 15-01-1511-5-D | 01/22/15 16:55 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:11 | 150126L6FF |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 40.4 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 5.15 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 4.92 | 0.500 | 0.103 | 1.00 | |
| Sodium | 9.94 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.262 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.09 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|----------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HCS 210 Trail | 15-01-1511-6-D | 01/23/15 03:01 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:12 | 150126L6FF |
|----------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 22.4 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 2.76 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.77 | 0.500 | 0.103 | 1.00 | |
| Sodium | 1.53 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.128 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 0.956 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-01-1511-7-D | 01/23/15 03:21 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:14 | 150126L6FF |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 81.3 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 15.7 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.43 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.3 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.682 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.70 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS 250 Trail | 15-01-1511-8-D | 01/23/15 03:03 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:16 | 150126L6FF |
|---------------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 79.5 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 14.6 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.55 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.675 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.40 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS 260 Trail | 15-01-1511-9-D | 01/23/15 03:36 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:17 | 150126L6FF |
|---------------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 57.6 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 8.87 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 3.91 | 0.500 | 0.103 | 1.00 | |
| Sodium | 10.2 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.409 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.52 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-01-1511-10-D | 01/23/15 03:30 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:19 | 150126L6FF |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 57.8 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 8.83 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 4.06 | 0.500 | 0.103 | 1.00 | |
| Sodium | 10.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.414 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.48 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|-----------------|-----------------|----------------|---------|----------|----------|----------------|------------|
| FDHCS 260 Trail | 15-01-1511-11-D | 01/23/15 03:36 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:21 | 150126L6FF |
|-----------------|-----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 56.0 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 8.74 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 3.82 | 0.500 | 0.103 | 1.00 | |
| Sodium | 10.0 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.401 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.68 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|-----------------|-----------------|----------------|---------|----------|----------|----------------|------------|
| FDHCS 270 Trail | 15-01-1511-12-D | 01/23/15 03:30 | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 18:23 | 150126L6FF |
|-----------------|-----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 55.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 8.93 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 3.86 | 0.500 | 0.103 | 1.00 | |
| Sodium | 9.77 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.399 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.64 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-683-1120 | N/A | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 17:52 | 150126L6FF |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | ND | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 0.0522 | 0.100 | 0.00336 | 1.00 | J |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Peak | 15-01-1511-1-D | 01/22/15 16:28 | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 08:45 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000137 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00110 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0115 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000452 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00140 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000447 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.000703 | 0.00100 | 0.000132 | 1.00 | J |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00893 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.135 | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.100 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00275 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-01-1511-2-D | 01/22/15 16:51 | Aqueous | ICP/MS 04 | 01/26/15 | 01/28/15 11:59 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|-----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0525 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00126 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.0000962 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00137 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0188 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0212 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0408 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00131 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Peak | 15-01-1511-3-D | 01/22/15 16:40 | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 08:59 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | 0.000565 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | 0.0494 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000832 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00156 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000238 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00538 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00453 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0385 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00189 | 0.00100 | 0.000139 | 1.00 | |

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Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-01-1511-4-D | 01/22/15 17:03 | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 09:03 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000177 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00237 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0481 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.00276 | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00280 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.00139 | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00270 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0514 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 2.25 | 0.0500 | 0.00331 | 1.00 | |
| Iron | 1.25 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0352 | 0.00100 | 0.000139 | 1.00 | |

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Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Peak | 15-01-1511-5-D | 01/22/15 16:55 | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 09:06 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000146 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00182 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0378 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00157 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00167 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000175 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00561 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0850 | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.101 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0126 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Trail | 15-01-1511-6-D | 01/23/15 03:01 | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 09:09 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | 0.000818 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | 0.0141 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00135 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.000886 | 0.00100 | 0.000132 | 1.00 | J |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00793 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0424 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00164 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-01-1511-7-D | 01/23/15 03:21 | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 09:13 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | 0.000729 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | 0.0575 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00100 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00230 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000395 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00392 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0315 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.000980 | 0.00100 | 0.000139 | 1.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Trail | 15-01-1511-8-D | 01/23/15 03:03 | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 09:16 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | 0.000486 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | 0.0576 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00102 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00184 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000379 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0237 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0378 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00155 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Trail | 15-01-1511-9-D | 01/23/15 03:36 | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 09:20 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000118 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00140 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0471 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00142 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00191 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000207 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00978 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0224 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0678 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00767 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-01-1511-10-D | 01/23/15 03:30 | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 09:23 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000110 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00109 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0498 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00142 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00180 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000171 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00605 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0535 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00408 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 260 Trail | 15-01-1511-11-D | 01/23/15 03:36 | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 09:27 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000104 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00134 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0466 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000525 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00141 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000110 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00171 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000180 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0135 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.188 | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.129 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00662 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 270 Trail | 15-01-1511-12-D | 01/23/15 03:30 | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 09:30 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000108 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00121 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0467 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000417 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00186 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00204 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000258 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0105 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0847 | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0808 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00446 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-719 | N/A | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 08:18 | 150126L03F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

Page 1 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Peak | 15-01-1511-1-D | 01/22/15 16:28 | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 17:47 | 150130L06F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 240 Peak | 15-01-1511-2-D | 01/22/15 16:51 | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 17:50 | 150130L06F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 250 Peak | 15-01-1511-3-D | 01/22/15 16:40 | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 17:52 | 150130L06F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 260 Peak | 15-01-1511-4-D | 01/22/15 17:03 | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 17:54 | 150130L06F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 270 Peak | 15-01-1511-5-D | 01/22/15 16:55 | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 17:56 | 150130L06F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 210 Trail | 15-01-1511-6-D | 01/23/15 03:01 | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 17:58 | 150130L06F |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-01-1511-7-D | 01/23/15 03:21 | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 18:05 | 150130L06F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 250 Trail | 15-01-1511-8-D | 01/23/15 03:03 | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 18:07 | 150130L06F |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 260 Trail | 15-01-1511-9-D | 01/23/15 03:36 | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 18:10 | 150130L06F |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 270 Trail | 15-01-1511-10-D | 01/23/15 03:30 | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 18:12 | 150130L06F |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|------------|
| FDHCS 260 Trail | 15-01-1511-11-D | 01/23/15 03:36 | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 18:14 | 150130L06F |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|------------|
| FDHCS 270 Trail | 15-01-1511-12-D | 01/23/15 03:30 | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 18:17 | 150130L06F |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-763-485 | N/A | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 16:47 | 150130L06F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

Page 1 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|-----------------------|-----------------------|----------------|-----------------------|-----------------|-----------------------|-------------------|
| HCS 210 Peak | 15-01-1511-1-i | 01/22/15 16:28 | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 18:03 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 95 | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 52 | | 31-200 | | | |
| HCS 240 Peak | 15-01-1511-2-i | 01/22/15 16:51 | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 18:11 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 69 | | 31-200 | | | |
| HCS 250 Peak | 15-01-1511-3-i | 01/22/15 16:40 | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 18:19 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 38 | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 48 | | 31-200 | | | |
| HCS 260 Peak | 15-01-1511-4-i | 01/22/15 17:03 | Aqueous | LC/TQ 2 | 01/29/15 | 02/09/15 14:32 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 130 | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 37 | | 31-200 | | | |
| HCS 270 Peak | 15-01-1511-5-i | 01/22/15 16:55 | Aqueous | LC/TQ 2 | 01/29/15 | 02/02/15 17:42 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 130 | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 35 | | 31-200 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|------------------------|-----------------------|----------------|-----------------------|-----------------|-----------------------|-------------------|
| HCS 210 Trail | 15-01-1511-6-i | 01/23/15 03:01 | Aqueous | LC/TQ 2 | 01/29/15 | 02/02/15 17:51 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 140 | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 44 | | 31-200 | | | |
| HCS 240 Trail | 15-01-1511-7-i | 01/23/15 03:21 | Aqueous | LC/TQ 2 | 01/29/15 | 02/02/15 17:59 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 18 | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 49 | | 31-200 | | | |
| HCS 250 Trail | 15-01-1511-8-i | 01/23/15 03:03 | Aqueous | LC/TQ 2 | 01/29/15 | 02/02/15 18:07 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 11 | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 45 | | 31-200 | | | |
| HCS 260 Trail | 15-01-1511-9-i | 01/23/15 03:36 | Aqueous | LC/TQ 2 | 01/29/15 | 02/02/15 18:16 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 110 | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 38 | | 31-200 | | | |
| HCS 270 Trail | 15-01-1511-10-i | 01/23/15 03:30 | Aqueous | LC/TQ 2 | 01/29/15 | 02/02/15 18:24 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 100 | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 41 | | 31-200 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|------------------------|-----------------------|----------------|-----------------------|-----------------|-----------------------|-------------------|
| FDHCS 260 Trail | 15-01-1511-11-i | 01/23/15 03:36 | Aqueous | LC/TQ 2 | 01/29/15 | 02/02/15 18:32 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 67 | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 42 | | 31-200 | | | |
| FDHCS 270 Trail | 15-01-1511-12-i | 01/23/15 03:30 | Aqueous | LC/TQ 2 | 01/29/15 | 02/02/15 18:41 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 65 | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 39 | | 31-200 | | | |
| Method Blank | 099-16-376-10 | N/A | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 17:06 | 150129L19 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 49 | | 31-200 | | | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Peak | 15-01-1511-1-I | 01/22/15 16:28 | Aqueous | GC 51 | 01/24/15 | 02/04/15 05:05 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.098 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.098 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.098 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.098 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.098 | 0.028 | 1.00 | |
| Aldrin | ND | 0.098 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.098 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.098 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.098 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.098 | 0.026 | 1.00 | |
| Endrin | ND | 0.098 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.098 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.098 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.098 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.098 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.098 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.098 | 0.025 | 1.00 | |
| Chlordane | ND | 0.98 | 0.32 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.58 | 1.00 | |
| Endrin Ketone | ND | 0.098 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 94 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 98 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

Page 2 of 13

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-01-1511-2-I | 01/22/15 16:51 | Aqueous | GC 51 | 01/24/15 | 02/04/15 05:19 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.097 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.097 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.097 | 0.028 | 1.00 | |
| Aldrin | ND | 0.097 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.097 | 0.024 | 1.00 | |
| Endosulfan I | ND | 0.097 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.097 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.097 | 0.026 | 1.00 | |
| Endrin | ND | 0.097 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan II | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDT | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.097 | 0.028 | 1.00 | |
| Methoxychlor | ND | 0.097 | 0.024 | 1.00 | |
| Chlordane | ND | 0.97 | 0.32 | 1.00 | |
| Toxaphene | ND | 1.9 | 0.57 | 1.00 | |
| Endrin Ketone | ND | 0.097 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 104 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 103 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Peak | 15-01-1511-3-I | 01/22/15 16:40 | Aqueous | GC 51 | 01/24/15 | 02/04/15 05:34 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.097 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.097 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.097 | 0.028 | 1.00 | |
| Aldrin | ND | 0.097 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.097 | 0.024 | 1.00 | |
| Endosulfan I | ND | 0.097 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.097 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.097 | 0.026 | 1.00 | |
| Endrin | ND | 0.097 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan II | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDT | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.097 | 0.028 | 1.00 | |
| Methoxychlor | ND | 0.097 | 0.024 | 1.00 | |
| Chlordane | ND | 0.97 | 0.32 | 1.00 | |
| Toxaphene | ND | 1.9 | 0.57 | 1.00 | |
| Endrin Ketone | ND | 0.097 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 105 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 105 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-01-1511-4-I | 01/22/15 17:03 | Aqueous | GC 51 | 01/24/15 | 02/04/15 05:48 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 95 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 97 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Peak | 15-01-1511-5-I | 01/22/15 16:55 | Aqueous | GC 51 | 01/24/15 | 02/04/15 06:02 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.098 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.098 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.098 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.098 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.098 | 0.028 | 1.00 | |
| Aldrin | ND | 0.098 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.098 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.098 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.098 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.098 | 0.026 | 1.00 | |
| Endrin | ND | 0.098 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.098 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.098 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.098 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.098 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.098 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.098 | 0.025 | 1.00 | |
| Chlordane | ND | 0.98 | 0.32 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.58 | 1.00 | |
| Endrin Ketone | ND | 0.098 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 84 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Trail | 15-01-1511-6-I | 01/23/15 03:01 | Aqueous | GC 51 | 01/24/15 | 02/04/15 06:17 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 98 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 99 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-01-1511-7-I | 01/23/15 03:21 | Aqueous | GC 51 | 01/24/15 | 02/04/15 06:31 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.097 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.097 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.097 | 0.028 | 1.00 | |
| Aldrin | ND | 0.097 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.097 | 0.024 | 1.00 | |
| Endosulfan I | ND | 0.097 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.097 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.097 | 0.026 | 1.00 | |
| Endrin | ND | 0.097 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan II | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDT | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.097 | 0.028 | 1.00 | |
| Methoxychlor | ND | 0.097 | 0.024 | 1.00 | |
| Chlordane | ND | 0.97 | 0.32 | 1.00 | |
| Toxaphene | ND | 1.9 | 0.57 | 1.00 | |
| Endrin Ketone | ND | 0.097 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 97 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 98 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Trail | 15-01-1511-8-I | 01/23/15 03:03 | Aqueous | GC 51 | 01/24/15 | 02/04/15 06:45 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.096 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.096 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.096 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.096 | 0.025 | 1.00 | |
| Delta-BHC | ND | 0.096 | 0.027 | 1.00 | |
| Aldrin | ND | 0.096 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.096 | 0.024 | 1.00 | |
| Endosulfan I | ND | 0.096 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.096 | 0.027 | 1.00 | |
| 4,4'-DDE | ND | 0.096 | 0.026 | 1.00 | |
| Endrin | ND | 0.096 | 0.029 | 1.00 | |
| Endrin Aldehyde | ND | 0.096 | 0.025 | 1.00 | |
| 4,4'-DDD | ND | 0.096 | 0.026 | 1.00 | |
| Endosulfan II | ND | 0.096 | 0.026 | 1.00 | |
| 4,4'-DDT | ND | 0.096 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.096 | 0.028 | 1.00 | |
| Methoxychlor | ND | 0.096 | 0.024 | 1.00 | |
| Chlordane | ND | 0.96 | 0.32 | 1.00 | |
| Toxaphene | ND | 1.9 | 0.57 | 1.00 | |
| Endrin Ketone | ND | 0.096 | 0.023 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 89 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Trail | 15-01-1511-9-I | 01/23/15 03:36 | Aqueous | GC 51 | 01/24/15 | 02/04/15 07:00 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.097 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.097 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.097 | 0.028 | 1.00 | |
| Aldrin | ND | 0.097 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.097 | 0.024 | 1.00 | |
| Endosulfan I | ND | 0.097 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.097 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.097 | 0.026 | 1.00 | |
| Endrin | ND | 0.097 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan II | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDT | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.097 | 0.028 | 1.00 | |
| Methoxychlor | ND | 0.097 | 0.024 | 1.00 | |
| Chlordane | ND | 0.97 | 0.32 | 1.00 | |
| Toxaphene | ND | 1.9 | 0.57 | 1.00 | |
| Endrin Ketone | ND | 0.097 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 93 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 94 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-01-1511-10-I | 01/23/15 03:30 | Aqueous | GC 51 | 01/24/15 | 02/04/15 07:14 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.097 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.097 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.097 | 0.028 | 1.00 | |
| Aldrin | ND | 0.097 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.097 | 0.024 | 1.00 | |
| Endosulfan I | ND | 0.097 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.097 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.097 | 0.026 | 1.00 | |
| Endrin | ND | 0.097 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan II | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDT | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.097 | 0.028 | 1.00 | |
| Methoxychlor | ND | 0.097 | 0.024 | 1.00 | |
| Chlordane | ND | 0.97 | 0.32 | 1.00 | |
| Toxaphene | ND | 1.9 | 0.57 | 1.00 | |
| Endrin Ketone | ND | 0.097 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 92 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 94 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 260 Trail | 15-01-1511-11-I | 01/23/15 03:36 | Aqueous | GC 51 | 01/24/15 | 02/04/15 07:28 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 93 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 93 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 270 Trail | 15-01-1511-12-I | 01/23/15 03:30 | Aqueous | GC 51 | 01/24/15 | 02/04/15 07:43 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|-------|-------|------|------------|
| Alpha-BHC | ND | 0.097 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.097 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.097 | 0.028 | 1.00 | |
| Aldrin | ND | 0.097 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.097 | 0.024 | 1.00 | |
| Endosulfan I | ND | 0.097 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.097 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.097 | 0.026 | 1.00 | |
| Endrin | ND | 0.097 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan II | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDT | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.097 | 0.028 | 1.00 | |
| Methoxychlor | ND | 0.097 | 0.024 | 1.00 | |
| Chlordane | ND | 0.97 | 0.32 | 1.00 | |
| Toxaphene | ND | 1.9 | 0.57 | 1.00 | |
| Endrin Ketone | ND | 0.097 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 83 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 80 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-778 | N/A | Aqueous | GC 51 | 01/24/15 | 02/04/15 04:50 | 150124L08A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 97 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 90 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

Page 1 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Peak | 15-01-1511-1-I | 01/22/15 16:28 | Aqueous | GC 31 | 01/24/15 | 02/04/15 16:07 | 150124L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|------|------|------|------------|
| Aroclor-1016 | ND | 0.98 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.98 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 0.98 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.98 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 0.98 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.98 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.98 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.98 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 68 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 80 | 50-135 | |

| HCS 240 Peak | 15-01-1511-2-I | 01/22/15 16:51 | Aqueous | GC 31 | 01/24/15 | 02/04/15 16:26 | 150124L09 |
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|------|------|------|------------|
| Aroclor-1016 | ND | 0.97 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.97 | 0.27 | 1.00 | |
| Aroclor-1232 | ND | 0.97 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.97 | 0.17 | 1.00 | |
| Aroclor-1248 | ND | 0.97 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.97 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.97 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.97 | 0.25 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 78 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 85 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

Page 2 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HCS 250 Peak | 15-01-1511-3-I | 01/22/15 16:40 | Aqueous | GC 31 | 01/24/15 | 02/05/15 17:30 | 150124L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 0.97 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.97 | 0.27 | 1.00 | |
| Aroclor-1232 | ND | 0.97 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.97 | 0.17 | 1.00 | |
| Aroclor-1248 | ND | 0.97 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.97 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.97 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.97 | 0.25 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 81 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 95 | 50-135 | |

| | | | | | | | |
|---------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HCS 260 Peak | 15-01-1511-4-I | 01/22/15 17:03 | Aqueous | GC 31 | 01/24/15 | 02/05/15 17:49 | 150124L09 |
|---------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 93 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 99 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HCS 270 Peak | 15-01-1511-5-I | 01/22/15 16:55 | Aqueous | GC 31 | 01/24/15 | 02/05/15 18:08 | 150124L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 0.98 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.98 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 0.98 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.98 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 0.98 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.98 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.98 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.98 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 75 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 88 | 50-135 | |

| | | | | | | | |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HCS 210 Trail | 15-01-1511-6-I | 01/23/15 03:01 | Aqueous | GC 31 | 01/24/15 | 02/05/15 18:27 | 150124L09 |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 90 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 103 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-01-1511-7-I | 01/23/15 03:21 | Aqueous | GC 31 | 01/24/15 | 02/05/15 18:46 | 150124L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|------|------|------|------------|
| Aroclor-1016 | ND | 0.97 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.97 | 0.27 | 1.00 | |
| Aroclor-1232 | ND | 0.97 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.97 | 0.17 | 1.00 | |
| Aroclor-1248 | ND | 0.97 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.97 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.97 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.97 | 0.25 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 111 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 122 | 50-135 | |

| HCS 250 Trail | 15-01-1511-8-I | 01/23/15 03:03 | Aqueous | GC 31 | 01/24/15 | 02/05/15 19:05 | 150124L09 |
|---------------|----------------|----------------|---------|-------|----------|----------------|-----------|
|---------------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|------|------|------|------------|
| Aroclor-1016 | ND | 0.96 | 0.28 | 1.00 | |
| Aroclor-1221 | ND | 0.96 | 0.27 | 1.00 | |
| Aroclor-1232 | ND | 0.96 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.96 | 0.17 | 1.00 | |
| Aroclor-1248 | ND | 0.96 | 0.19 | 1.00 | |
| Aroclor-1254 | ND | 0.96 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.96 | 0.25 | 1.00 | |
| Aroclor-1262 | ND | 0.96 | 0.25 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 81 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 90 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Trail | 15-01-1511-9-I | 01/23/15 03:36 | Aqueous | GC 31 | 01/24/15 | 02/05/15 19:24 | 150124L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|------|------|------|------------|
| Aroclor-1016 | ND | 0.97 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.97 | 0.27 | 1.00 | |
| Aroclor-1232 | ND | 0.97 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.97 | 0.17 | 1.00 | |
| Aroclor-1248 | ND | 0.97 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.97 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.97 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.97 | 0.25 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 137 | 50-135 | 2,7 |
| 2,4,5,6-Tetrachloro-m-Xylene | 151 | 50-135 | 2,7 |

| HCS 270 Trail | 15-01-1511-10-I | 01/23/15 03:30 | Aqueous | GC 31 | 01/24/15 | 02/05/15 19:43 | 150124L09 |
|---------------|-----------------|-------------------|---------|-------|----------|-------------------|-----------|
|---------------|-----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|------|------|------|------------|
| Aroclor-1016 | ND | 0.97 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.97 | 0.27 | 1.00 | |
| Aroclor-1232 | ND | 0.97 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.97 | 0.17 | 1.00 | |
| Aroclor-1248 | ND | 0.97 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.97 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.97 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.97 | 0.25 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 82 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 87 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 260 Trail | 15-01-1511-11-I | 01/23/15 03:36 | Aqueous | GC 31 | 01/24/15 | 02/05/15 20:02 | 150124L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 93 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 101 | 50-135 | |

| FDHCS 270 Trail | 15-01-1511-12-I | 01/23/15 03:30 | Aqueous | GC 31 | 01/24/15 | 02/05/15 20:21 | 150124L09 |
|-----------------|-----------------|----------------|---------|-------|----------|----------------|-----------|
|-----------------|-----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|------|------|------|------------|
| Aroclor-1016 | ND | 0.97 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.97 | 0.27 | 1.00 | |
| Aroclor-1232 | ND | 0.97 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.97 | 0.17 | 1.00 | |
| Aroclor-1248 | ND | 0.97 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.97 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.97 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.97 | 0.25 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 87 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 96 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-533-1000 | N/A | Aqueous | GC 31 | 01/24/15 | 02/04/15 11:03 | 150124L09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 71 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 79 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Peak | 15-01-1511-1-I | 01/22/15 16:28 | Aqueous | GC 26 | 01/26/15 | 02/08/15 15:05 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0049 | 0.0028 | 1.00 | |
| Bolstar | ND | 0.0049 | 0.0028 | 1.00 | |
| Chlorpyrifos | ND | 0.0049 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0049 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0049 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0049 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.0098 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0049 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0049 | 0.0028 | 1.00 | |
| Fenthion | ND | 0.0049 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0049 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0049 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0049 | 0.0027 | 1.00 | |
| Naled | ND | 0.039 | 0.019 | 1.00 | |
| Phorate | ND | 0.0049 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0049 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0086 | 1.00 | |
| Tokuthion | ND | 0.0049 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0049 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0049 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 113 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-01-1511-2-I | 01/22/15 16:51 | Aqueous | GC 26 | 01/26/15 | 02/08/15 15:49 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0048 | 0.0027 | 1.00 | |
| Bolstar | ND | 0.0048 | 0.0027 | 1.00 | |
| Chlorpyrifos | ND | 0.0048 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0048 | 0.0023 | 1.00 | |
| Diazinon | ND | 0.0048 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0048 | 0.0035 | 1.00 | |
| Disulfoton | ND | 0.0096 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0048 | 0.0024 | 1.00 | |
| Fensulfothion | ND | 0.0048 | 0.0027 | 1.00 | |
| Fenthion | ND | 0.0048 | 0.0025 | 1.00 | |
| Merphos | ND | 0.0048 | 0.0025 | 1.00 | |
| Methyl Parathion | ND | 0.0048 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0048 | 0.0026 | 1.00 | |
| Naled | ND | 0.038 | 0.019 | 1.00 | |
| Phorate | ND | 0.0048 | 0.0024 | 1.00 | |
| Ronnel | ND | 0.0048 | 0.0031 | 1.00 | |
| Stirophos | ND | 0.019 | 0.0085 | 1.00 | |
| Tokuthion | ND | 0.0048 | 0.0027 | 1.00 | |
| Trichloronate | ND | 0.0048 | 0.0020 | 1.00 | |
| Demeton-o/s | ND | 0.0048 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 126 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 3 of 13

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Peak | 15-01-1511-3-I | 01/22/15 16:40 | Aqueous | GC 26 | 01/26/15 | 02/08/15 16:33 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0049 | 0.0028 | 1.00 | |
| Bolstar | ND | 0.0049 | 0.0028 | 1.00 | |
| Chlorpyrifos | ND | 0.0049 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0049 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0049 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0049 | 0.0035 | 1.00 | |
| Disulfoton | ND | 0.0097 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0049 | 0.0024 | 1.00 | |
| Fensulfothion | ND | 0.0049 | 0.0028 | 1.00 | |
| Fenthion | ND | 0.0049 | 0.0025 | 1.00 | |
| Merphos | ND | 0.0049 | 0.0025 | 1.00 | |
| Methyl Parathion | ND | 0.0049 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0049 | 0.0027 | 1.00 | |
| Naled | ND | 0.039 | 0.019 | 1.00 | |
| Phorate | ND | 0.0049 | 0.0024 | 1.00 | |
| Ronnel | ND | 0.0049 | 0.0031 | 1.00 | |
| Stirophos | ND | 0.019 | 0.0086 | 1.00 | |
| Tokuthion | ND | 0.0049 | 0.0027 | 1.00 | |
| Trichloronate | ND | 0.0049 | 0.0020 | 1.00 | |
| Demeton-o/s | ND | 0.0049 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 127 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-01-1511-4-I | 01/22/15 17:03 | Aqueous | GC 26 | 01/26/15 | 02/08/15 17:17 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0048 | 0.0027 | 1.00 | |
| Bolstar | ND | 0.0048 | 0.0027 | 1.00 | |
| Chlorpyrifos | ND | 0.0048 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0048 | 0.0023 | 1.00 | |
| Diazinon | ND | 0.0048 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0048 | 0.0035 | 1.00 | |
| Disulfoton | ND | 0.0096 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0048 | 0.0024 | 1.00 | |
| Fensulfothion | ND | 0.0048 | 0.0027 | 1.00 | |
| Fenthion | ND | 0.0048 | 0.0025 | 1.00 | |
| Merphos | ND | 0.0048 | 0.0025 | 1.00 | |
| Methyl Parathion | ND | 0.0048 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0048 | 0.0026 | 1.00 | |
| Naled | ND | 0.038 | 0.019 | 1.00 | |
| Phorate | ND | 0.0048 | 0.0024 | 1.00 | |
| Ronnel | ND | 0.0048 | 0.0031 | 1.00 | |
| Stirophos | ND | 0.019 | 0.0085 | 1.00 | |
| Tokuthion | ND | 0.0048 | 0.0027 | 1.00 | |
| Trichloronate | ND | 0.0048 | 0.0020 | 1.00 | |
| Demeton-o/s | ND | 0.0048 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 118 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Peak | 15-01-1511-5-I | 01/22/15 16:55 | Aqueous | GC 26 | 01/26/15 | 02/08/15 18:01 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0048 | 0.0027 | 1.00 | |
| Bolstar | ND | 0.0048 | 0.0027 | 1.00 | |
| Chlorpyrifos | ND | 0.0048 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0048 | 0.0023 | 1.00 | |
| Diazinon | ND | 0.0048 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0048 | 0.0035 | 1.00 | |
| Disulfoton | ND | 0.0096 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0048 | 0.0024 | 1.00 | |
| Fensulfothion | ND | 0.0048 | 0.0027 | 1.00 | |
| Fenthion | ND | 0.0048 | 0.0025 | 1.00 | |
| Merphos | ND | 0.0048 | 0.0025 | 1.00 | |
| Methyl Parathion | ND | 0.0048 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0048 | 0.0026 | 1.00 | |
| Naled | ND | 0.038 | 0.019 | 1.00 | |
| Phorate | ND | 0.0048 | 0.0024 | 1.00 | |
| Ronnel | ND | 0.0048 | 0.0031 | 1.00 | |
| Stirophos | ND | 0.019 | 0.0085 | 1.00 | |
| Tokuthion | ND | 0.0048 | 0.0027 | 1.00 | |
| Trichloronate | ND | 0.0048 | 0.0020 | 1.00 | |
| Demeton-o/s | ND | 0.0048 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 127 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Trail | 15-01-1511-6-I | 01/23/15 03:01 | Aqueous | GC 26 | 01/26/15 | 02/08/15 18:46 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0028 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0028 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.0099 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0028 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0087 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 127 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-01-1511-7-I | 01/23/15 03:21 | Aqueous | GC 26 | 01/26/15 | 02/08/15 19:30 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0048 | 0.0027 | 1.00 | |
| Bolstar | ND | 0.0048 | 0.0027 | 1.00 | |
| Chlorpyrifos | ND | 0.0048 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0048 | 0.0023 | 1.00 | |
| Diazinon | ND | 0.0048 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0048 | 0.0035 | 1.00 | |
| Disulfoton | ND | 0.0096 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0048 | 0.0024 | 1.00 | |
| Fensulfothion | ND | 0.0048 | 0.0027 | 1.00 | |
| Fenthion | ND | 0.0048 | 0.0025 | 1.00 | |
| Merphos | ND | 0.0048 | 0.0025 | 1.00 | |
| Methyl Parathion | ND | 0.0048 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0048 | 0.0026 | 1.00 | |
| Naled | ND | 0.038 | 0.019 | 1.00 | |
| Phorate | ND | 0.0048 | 0.0024 | 1.00 | |
| Ronnel | ND | 0.0048 | 0.0031 | 1.00 | |
| Stirophos | ND | 0.019 | 0.0085 | 1.00 | |
| Tokuthion | ND | 0.0048 | 0.0027 | 1.00 | |
| Trichloronate | ND | 0.0048 | 0.0020 | 1.00 | |
| Demeton-o/s | ND | 0.0048 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 129 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Trail | 15-01-1511-8-I | 01/23/15 03:03 | Aqueous | GC 26 | 01/26/15 | 02/08/15 20:14 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 130 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Trail | 15-01-1511-9-I | 01/23/15 03:36 | Aqueous | GC 26 | 01/26/15 | 02/08/15 20:58 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0049 | 0.0028 | 1.00 | |
| Bolstar | ND | 0.0049 | 0.0028 | 1.00 | |
| Chlorpyrifos | ND | 0.0049 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0049 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0049 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0049 | 0.0035 | 1.00 | |
| Disulfoton | ND | 0.0097 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0049 | 0.0024 | 1.00 | |
| Fensulfothion | ND | 0.0049 | 0.0028 | 1.00 | |
| Fenthion | ND | 0.0049 | 0.0025 | 1.00 | |
| Merphos | ND | 0.0049 | 0.0025 | 1.00 | |
| Methyl Parathion | ND | 0.0049 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0049 | 0.0027 | 1.00 | |
| Naled | ND | 0.039 | 0.019 | 1.00 | |
| Phorate | ND | 0.0049 | 0.0024 | 1.00 | |
| Ronnel | ND | 0.0049 | 0.0031 | 1.00 | |
| Stirophos | ND | 0.019 | 0.0086 | 1.00 | |
| Tokuthion | ND | 0.0049 | 0.0027 | 1.00 | |
| Trichloronate | ND | 0.0049 | 0.0020 | 1.00 | |
| Demeton-o/s | ND | 0.0049 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 127 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-01-1511-10-I | 01/23/15 03:30 | Aqueous | GC 26 | 01/26/15 | 02/08/15 21:42 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0049 | 0.0028 | 1.00 | |
| Bolstar | ND | 0.0049 | 0.0028 | 1.00 | |
| Chlorpyrifos | ND | 0.0049 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0049 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0049 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0049 | 0.0035 | 1.00 | |
| Disulfoton | ND | 0.0097 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0049 | 0.0024 | 1.00 | |
| Fensulfothion | ND | 0.0049 | 0.0028 | 1.00 | |
| Fenthion | ND | 0.0049 | 0.0025 | 1.00 | |
| Merphos | ND | 0.0049 | 0.0025 | 1.00 | |
| Methyl Parathion | ND | 0.0049 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0049 | 0.0027 | 1.00 | |
| Naled | ND | 0.039 | 0.019 | 1.00 | |
| Phorate | ND | 0.0049 | 0.0024 | 1.00 | |
| Ronnel | ND | 0.0049 | 0.0031 | 1.00 | |
| Stirophos | ND | 0.019 | 0.0086 | 1.00 | |
| Tokuthion | ND | 0.0049 | 0.0027 | 1.00 | |
| Trichloronate | ND | 0.0049 | 0.0020 | 1.00 | |
| Demeton-o/s | ND | 0.0049 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 128 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 260 Trail | 15-01-1511-11-I | 01/23/15 03:36 | Aqueous | GC 26 | 01/26/15 | 02/08/15 22:26 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0049 | 0.0028 | 1.00 | |
| Bolstar | ND | 0.0049 | 0.0028 | 1.00 | |
| Chlorpyrifos | ND | 0.0049 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0049 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0049 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0049 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.0098 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0049 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0049 | 0.0028 | 1.00 | |
| Fenthion | ND | 0.0049 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0049 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0049 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0049 | 0.0027 | 1.00 | |
| Naled | ND | 0.039 | 0.019 | 1.00 | |
| Phorate | ND | 0.0049 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0049 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0086 | 1.00 | |
| Tokuthion | ND | 0.0049 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0049 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0049 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 123 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 270 Trail | 15-01-1511-12-I | 01/23/15 03:30 | Aqueous | GC 26 | 01/26/15 | 02/08/15 23:10 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0049 | 0.0028 | 1.00 | |
| Bolstar | ND | 0.0049 | 0.0028 | 1.00 | |
| Chlorpyrifos | ND | 0.0049 | 0.0023 | 1.00 | |
| Coumaphos | ND | 0.0049 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0049 | 0.0028 | 1.00 | |
| Dichlorvos | ND | 0.0049 | 0.0035 | 1.00 | |
| Disulfoton | ND | 0.0097 | 0.0025 | 1.00 | |
| Ethoprop | ND | 0.0049 | 0.0024 | 1.00 | |
| Fensulfothion | ND | 0.0049 | 0.0028 | 1.00 | |
| Fenthion | ND | 0.0049 | 0.0025 | 1.00 | |
| Merphos | ND | 0.0049 | 0.0025 | 1.00 | |
| Methyl Parathion | ND | 0.0049 | 0.0029 | 1.00 | |
| Mevinphos | ND | 0.0049 | 0.0027 | 1.00 | |
| Naled | ND | 0.039 | 0.019 | 1.00 | |
| Phorate | ND | 0.0049 | 0.0024 | 1.00 | |
| Ronnel | ND | 0.0049 | 0.0031 | 1.00 | |
| Stirophos | ND | 0.019 | 0.0086 | 1.00 | |
| Tokuthion | ND | 0.0049 | 0.0027 | 1.00 | |
| Trichloronate | ND | 0.0049 | 0.0020 | 1.00 | |
| Demeton-o/s | ND | 0.0049 | 0.0027 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 129 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-83 | N/A | Aqueous | GC 26 | 01/26/15 | 02/08/15 14:21 | 150126L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 100 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

Page 1 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HCS 210 Peak | 15-01-1511-1-I | 01/22/15 16:28 | Aqueous | GC 40 | 01/26/15 | 01/29/15 22:12 | 150126L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 66 | 0-123 | |

| | | | | | | | |
|---------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HCS 240 Peak | 15-01-1511-2-I | 01/22/15 16:51 | Aqueous | GC 40 | 01/26/15 | 01/29/15 22:35 | 150126L04 |
|---------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 63 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HCS 250 Peak | 15-01-1511-3-I | 01/22/15 16:40 | Aqueous | GC 40 | 01/26/15 | 01/29/15 22:58 | 150126L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 68 | 0-123 | |

| | | | | | | | |
|---------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HCS 260 Peak | 15-01-1511-4-I | 01/22/15 17:03 | Aqueous | GC 40 | 01/26/15 | 01/29/15 23:21 | 150126L04 |
|---------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 65 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Peak | 15-01-1511-5-I | 01/22/15 16:55 | Aqueous | GC 40 | 01/26/15 | 01/29/15 23:44 | 150126L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 66 | 0-123 | |

| HCS 210 Trail | 15-01-1511-6-I | 01/23/15 03:01 | Aqueous | GC 40 | 01/26/15 | 01/30/15 00:07 | 150126L04 |
|---------------|----------------|----------------|---------|-------|----------|----------------|-----------|
|---------------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 62 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-01-1511-7-I | 01/23/15 03:21 | Aqueous | GC 40 | 01/26/15 | 01/30/15 00:30 | 150126L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 62 | 0-123 | |

| HCS 250 Trail | 15-01-1511-8-I | 01/23/15 03:03 | Aqueous | GC 40 | 01/26/15 | 01/30/15 00:53 | 150126L04 |
|---------------|----------------|----------------|---------|-------|----------|----------------|-----------|
|---------------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 60 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Trail | 15-01-1511-9-I | 01/23/15 03:36 | Aqueous | GC 40 | 01/26/15 | 01/30/15 01:16 | 150126L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 61 | 0-123 | |

| HCS 270 Trail | 15-01-1511-10-I | 01/23/15 03:30 | Aqueous | GC 40 | 01/26/15 | 01/30/15 01:39 | 150126L04 |
|---------------|-----------------|-------------------|---------|-------|----------|-------------------|-----------|
|---------------|-----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 63 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| FDHCS 260 Trail | 15-01-1511-11-I | 01/23/15 03:36 | Aqueous | GC 40 | 01/26/15 | 01/30/15 02:02 | 150126L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 67 | 0-123 | |

| | | | | | | | |
|------------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| FDHCS 270 Trail | 15-01-1511-12-I | 01/23/15 03:30 | Aqueous | GC 40 | 01/26/15 | 01/30/15 02:48 | 150126L04 |
|------------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 64 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-034-634 | N/A | Aqueous | GC 40 | 01/26/15 | 01/29/15 19:53 | 150126L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 60 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Peak | 15-01-1511-1-G | 01/22/15 16:28 | Aqueous | GC/MS CCC | 01/26/15 | 02/01/15 14:17 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 67 | 33-120 | |
| 2-Fluorophenol | 47 | 24-120 | |
| Nitrobenzene-d5 | 64 | 38-120 | |
| p-Terphenyl-d14 | 73 | 41-137 | |
| Phenol-d6 | 31 | 16-120 | |
| 2,4,6-Tribromophenol | 86 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-01-1511-2-G | 01/22/15 16:51 | Aqueous | GC/MS CCC | 01/26/15 | 02/01/15 14:35 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 66 | 33-120 | |
| 2-Fluorophenol | 49 | 24-120 | |
| Nitrobenzene-d5 | 64 | 38-120 | |
| p-Terphenyl-d14 | 70 | 41-137 | |
| Phenol-d6 | 31 | 16-120 | |
| 2,4,6-Tribromophenol | 81 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Peak | 15-01-1511-3-G | 01/22/15 16:40 | Aqueous | GC/MS CCC | 01/26/15 | 02/01/15 14:53 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 67 | 33-120 | |
| 2-Fluorophenol | 45 | 24-120 | |
| Nitrobenzene-d5 | 66 | 38-120 | |
| p-Terphenyl-d14 | 78 | 41-137 | |
| Phenol-d6 | 30 | 16-120 | |
| 2,4,6-Tribromophenol | 92 | 27-159 | |



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-01-1511-4-G | 01/22/15 17:03 | Aqueous | GC/MS CCC | 01/26/15 | 02/01/15 15:11 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 74 | 33-120 | |
| 2-Fluorophenol | 48 | 24-120 | |
| Nitrobenzene-d5 | 71 | 38-120 | |
| p-Terphenyl-d14 | 78 | 41-137 | |
| Phenol-d6 | 33 | 16-120 | |
| 2,4,6-Tribromophenol | 85 | 27-159 | |



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Peak | 15-01-1511-5-G | 01/22/15 16:55 | Aqueous | GC/MS CCC | 01/26/15 | 02/01/15 15:30 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | 3.4 | 9.5 | 2.4 | 1.00 | J |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 66 | 33-120 | |
| 2-Fluorophenol | 43 | 24-120 | |
| Nitrobenzene-d5 | 65 | 38-120 | |
| p-Terphenyl-d14 | 71 | 41-137 | |
| Phenol-d6 | 29 | 16-120 | |
| 2,4,6-Tribromophenol | 80 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Trail | 15-01-1511-6-G | 01/23/15 03:01 | Aqueous | GC/MS CCC | 01/26/15 | 02/01/15 15:48 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 63 | 33-120 | |
| 2-Fluorophenol | 42 | 24-120 | |
| Nitrobenzene-d5 | 62 | 38-120 | |
| p-Terphenyl-d14 | 73 | 41-137 | |
| Phenol-d6 | 28 | 16-120 | |
| 2,4,6-Tribromophenol | 84 | 27-159 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-01-1511-7-G | 01/23/15 03:21 | Aqueous | GC/MS CCC | 01/26/15 | 02/01/15 16:06 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 74 | 33-120 | |
| 2-Fluorophenol | 53 | 24-120 | |
| Nitrobenzene-d5 | 72 | 38-120 | |
| p-Terphenyl-d14 | 77 | 41-137 | |
| Phenol-d6 | 35 | 16-120 | |
| 2,4,6-Tribromophenol | 91 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Trail | 15-01-1511-8-G | 01/23/15 03:03 | Aqueous | GC/MS CCC | 01/26/15 | 01/29/15 16:59 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 78 | 33-120 | |
| 2-Fluorophenol | 58 | 24-120 | |
| Nitrobenzene-d5 | 76 | 38-120 | |
| p-Terphenyl-d14 | 85 | 41-137 | |
| Phenol-d6 | 38 | 16-120 | |
| 2,4,6-Tribromophenol | 92 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Trail | 15-01-1511-9-G | 01/23/15 03:36 | Aqueous | GC/MS CCC | 01/26/15 | 01/29/15 17:17 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 75 | 33-120 | |
| 2-Fluorophenol | 57 | 24-120 | |
| Nitrobenzene-d5 | 71 | 38-120 | |
| p-Terphenyl-d14 | 82 | 41-137 | |
| Phenol-d6 | 40 | 16-120 | |
| 2,4,6-Tribromophenol | 94 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-01-1511-10-G | 01/23/15 03:30 | Aqueous | GC/MS CCC | 01/26/15 | 01/29/15 17:35 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 14 | 4.0 | 1.00 | |
| Acenaphthylene | ND | 14 | 4.1 | 1.00 | |
| Aniline | ND | 14 | 2.1 | 1.00 | |
| Anthracene | ND | 14 | 4.3 | 1.00 | |
| Azobenzene | ND | 14 | 3.8 | 1.00 | |
| Benzidine | ND | 71 | 9.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 14 | 6.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 14 | 3.5 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 14 | 3.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 14 | 3.6 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 14 | 4.6 | 1.00 | |
| Benzoic Acid | ND | 71 | 17 | 1.00 | |
| Benzyl Alcohol | ND | 14 | 3.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 14 | 3.6 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 36 | 3.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 14 | 4.6 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 14 | 4.5 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 14 | 3.9 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 14 | 3.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 14 | 3.4 | 1.00 | |
| 4-Chloroaniline | ND | 14 | 2.8 | 1.00 | |
| 2-Chloronaphthalene | ND | 14 | 3.9 | 1.00 | |
| 2-Chlorophenol | ND | 14 | 3.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 14 | 3.8 | 1.00 | |
| Chrysene | ND | 14 | 4.1 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 14 | 4.2 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 14 | 3.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 14 | 3.6 | 1.00 | |
| Dibenzofuran | ND | 14 | 4.0 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 14 | 4.3 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 14 | 4.4 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 14 | 4.1 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 36 | 3.7 | 1.00 | |
| 2,4-Dichlorophenol | ND | 14 | 3.6 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 14 | 4.0 | 1.00 | |
| Dimethyl Phthalate | ND | 14 | 3.7 | 1.00 | |
| 2,4-Dimethylphenol | ND | 14 | 3.5 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 71 | 20 | 1.00 | |
| 2,4-Dinitrophenol | ND | 71 | 19 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 14 | 3.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 14 | 3.4 | 1.00 | |
| Fluoranthene | ND | 14 | 4.4 | 1.00 | |
| Fluorene | ND | 14 | 3.9 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 14 | 4.1 | 1.00 | |
| Hexachlorobenzene | ND | 14 | 4.4 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 36 | 9.9 | 1.00 | |
| Hexachloroethane | ND | 14 | 4.3 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 14 | 3.1 | 1.00 | |
| Isophorone | ND | 14 | 3.6 | 1.00 | |
| 2-Methylnaphthalene | ND | 14 | 4.0 | 1.00 | |
| 1-Methylnaphthalene | ND | 14 | 4.0 | 1.00 | |
| 2-Methylphenol | ND | 14 | 3.0 | 1.00 | |
| 3/4-Methylphenol | ND | 14 | 3.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 14 | 3.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 14 | 4.6 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 14 | 3.9 | 1.00 | |
| Naphthalene | ND | 14 | 4.1 | 1.00 | |
| 4-Nitroaniline | ND | 14 | 3.0 | 1.00 | |
| 3-Nitroaniline | ND | 14 | 3.2 | 1.00 | |
| 2-Nitroaniline | ND | 14 | 3.2 | 1.00 | |
| Nitrobenzene | ND | 36 | 4.3 | 1.00 | |
| 4-Nitrophenol | ND | 14 | 2.3 | 1.00 | |
| 2-Nitrophenol | ND | 14 | 3.7 | 1.00 | |
| Pentachlorophenol | ND | 14 | 6.6 | 1.00 | |
| Phenanthrene | ND | 14 | 4.2 | 1.00 | |
| Phenol | ND | 14 | 2.9 | 1.00 | |
| Pyrene | ND | 14 | 4.3 | 1.00 | |
| Pyridine | ND | 14 | 4.3 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 14 | 4.1 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 14 | 3.6 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 14 | 3.6 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 63 | 33-120 | |
| 2-Fluorophenol | 60 | 24-120 | |
| Nitrobenzene-d5 | 69 | 38-120 | |
| p-Terphenyl-d14 | 77 | 41-137 | |
| Phenol-d6 | 44 | 16-120 | |
| 2,4,6-Tribromophenol | 90 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 260 Trail | 15-01-1511-11-G | 01/23/15 03:36 | Aqueous | GC/MS CCC | 01/26/15 | 02/01/15 16:24 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 63 | 33-120 | |
| 2-Fluorophenol | 47 | 24-120 | |
| Nitrobenzene-d5 | 68 | 38-120 | |
| p-Terphenyl-d14 | 75 | 41-137 | |
| Phenol-d6 | 31 | 16-120 | |
| 2,4,6-Tribromophenol | 86 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 270 Trail | 15-01-1511-12-G | 01/23/15 03:30 | Aqueous | GC/MS CCC | 01/26/15 | 01/29/15 17:53 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 65 | 33-120 | |
| 2-Fluorophenol | 55 | 24-120 | |
| Nitrobenzene-d5 | 73 | 38-120 | |
| p-Terphenyl-d14 | 78 | 41-137 | |
| Phenol-d6 | 38 | 16-120 | |
| 2,4,6-Tribromophenol | 92 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-3987 | N/A | Aqueous | GC/MS CCC | 01/26/15 | 01/29/15 16:41 | 150126L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 80 | 33-120 | |
| 2-Fluorophenol | 62 | 24-120 | |
| Nitrobenzene-d5 | 77 | 38-120 | |
| p-Terphenyl-d14 | 85 | 41-137 | |
| Phenol-d6 | 41 | 16-120 | |
| 2,4,6-Tribromophenol | 99 | 27-159 | |



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 1 of 30

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Peak | 15-01-1511-1-A | 01/22/15 16:28 | Aqueous | GC/MS O | 01/24/15 | 01/24/15 23:56 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 2 of 30

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|-----------------|-----------------------|-------------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | |
| 1,4-Bromofluorobenzene | 95 | 80-120 | | | |
| Dibromofluoromethane | 111 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 105 | 75-135 | | | |
| Toluene-d8 | 99 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-01-1511-2-A | 01/22/15 16:51 | Aqueous | GC/MS O | 01/24/15 | 01/25/15 00:23 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 110 | 78-126 | |
| 1,2-Dichloroethane-d4 | 105 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Peak | 15-01-1511-3-A | 01/22/15 16:40 | Aqueous | GC/MS O | 01/24/15 | 01/25/15 00:51 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 97 | 80-120 | |
| Dibromofluoromethane | 112 | 78-126 | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 7 of 30

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-01-1511-4-A | 01/22/15 17:03 | Aqueous | GC/MS O | 01/24/15 | 01/25/15 01:19 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 8 of 30

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 95 | 80-120 | |
| Dibromofluoromethane | 112 | 78-126 | |
| 1,2-Dichloroethane-d4 | 108 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 9 of 30

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Peak | 15-01-1511-5-A | 01/22/15 16:55 | Aqueous | GC/MS O | 01/24/15 | 01/25/15 01:46 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 108 | 78-126 | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Trail | 15-01-1511-6-A | 01/23/15 03:01 | Aqueous | GC/MS O | 01/24/15 | 01/25/15 02:14 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 95 | 80-120 | |
| Dibromofluoromethane | 111 | 78-126 | |
| 1,2-Dichloroethane-d4 | 107 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 13 of 30

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-01-1511-7-A | 01/23/15 03:21 | Aqueous | GC/MS O | 01/24/15 | 01/25/15 02:42 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 14 of 30

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 95 | 80-120 | |
| Dibromofluoromethane | 111 | 78-126 | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 15 of 30

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Trail | 15-01-1511-8-A | 01/23/15 03:03 | Aqueous | GC/MS O | 01/24/15 | 01/25/15 03:09 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 96 | 80-120 | |
| Dibromofluoromethane | 112 | 78-126 | |
| 1,2-Dichloroethane-d4 | 107 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 17 of 30

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Trail | 15-01-1511-9-A | 01/23/15 03:36 | Aqueous | GC/MS O | 01/24/15 | 01/25/15 03:37 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 95 | 80-120 | |
| Dibromofluoromethane | 112 | 78-126 | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-01-1511-10-A | 01/23/15 03:30 | Aqueous | GC/MS O | 01/24/15 | 01/25/15 04:04 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 20 of 30

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 96 | 80-120 | |
| Dibromofluoromethane | 112 | 78-126 | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | |
| Toluene-d8 | 103 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 21 of 30

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 260 Trail | 15-01-1511-11-A | 01/23/15 03:36 | Aqueous | GC/MS O | 01/24/15 | 01/25/15 04:32 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|-----------------|-----------------------|-------------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | |
| 1,4-Bromofluorobenzene | 95 | 80-120 | | | |
| Dibromofluoromethane | 114 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 105 | 75-135 | | | |
| Toluene-d8 | 101 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 270 Trail | 15-01-1511-12-A | 01/23/15 03:30 | Aqueous | GC/MS O | 01/24/15 | 01/25/15 05:00 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 97 | 80-120 | |
| Dibromofluoromethane | 111 | 78-126 | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | |
| Toluene-d8 | 103 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| TB01 | 15-01-1511-13-A | 01/23/15 00:00 | Aqueous | GC/MS O | 01/30/15 | 01/30/15 16:08 | 150130L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 88 | 80-120 | |
| Dibromofluoromethane | 110 | 78-126 | |
| 1,2-Dichloroethane-d4 | 113 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 27 of 30

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-16246 | N/A | Aqueous | GC/MS O | 01/24/15 | 01/24/15 21:10 | 150124L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 95 | 80-120 | |
| Dibromofluoromethane | 109 | 78-126 | |
| 1,2-Dichloroethane-d4 | 100 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-16277 | N/A | Aqueous | GC/MS O | 01/30/15 | 01/30/15 13:50 | 150130L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 90 | 80-120 | |
| Dibromofluoromethane | 107 | 78-126 | |
| 1,2-Dichloroethane-d4 | 113 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

01/24/15

Work Order:

15-01-1511

Page 1 of 7

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 210 Peak | 15-01-1511-1 | 01/22/15 16:28 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.068 | 0.10 | 0.025 | 1.00 | J | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | 1.4 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | 0.26 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | 3.3 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.084 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 58.0 | 1.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 58.0 | 1.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | 165 | 1.00 | 0.820 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | 18 | 1.0 | 0.95 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| pH (24) | 7.17 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.0 | 0.50 | 0.24 | 1.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 5.4 | 0.50 | 0.24 | 1.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

| | | | |
|---------------------|---------------------|-----------------------|----------------|
| HCS 240 Peak | 15-01-1511-2 | 01/22/15 16:51 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.24 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | 17 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | 0.12 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.7 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | 29 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 220 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 220 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | 340 | 1.00 | 0.820 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | 2.8 | 1.0 | 0.95 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| pH (24) | 7.56 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 8.6 | 0.50 | 0.24 | 1.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 5.6 | 0.50 | 0.24 | 1.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

01/24/15

Work Order:

15-01-1511

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 250 Peak | 15-01-1511-3 | 01/22/15 16:40 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.21 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | 16 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | 0.093 | 0.10 | 0.037 | 1.00 | J | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.5 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | 27 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 204 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 204 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | 440 | 1.00 | 0.820 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | 3.7 | 1.0 | 0.95 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| pH (24) | 7.63 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 4.7 | 0.50 | 0.24 | 1.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 5.6 | 0.50 | 0.24 | 1.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

| | | | |
|---------------------|---------------------|-----------------------|----------------|
| HCS 260 Peak | 15-01-1511-4 | 01/22/15 17:03 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.20 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | 9.4 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | 0.062 | 0.10 | 0.037 | 1.00 | J | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | 0.94 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | 18 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.12 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 110 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 110 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | 205 | 1.00 | 0.820 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | 368 | 1.00 | 0.950 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| pH (24) | 7.65 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.0 | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 11 | 0.50 | 0.24 | 1.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 8.8 | 0.50 | 0.24 | 1.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

01/24/15

Work Order:

15-01-1511

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 270 Peak | 15-01-1511-5 | 01/22/15 16:55 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.19 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | 11 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | 0.078 | 0.10 | 0.037 | 1.00 | J | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | 0.98 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | 22 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.12 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 122 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 122 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | 265 | 1.00 | 0.820 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | 409 | 1.00 | 0.950 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| pH (24) | 7.70 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.5 | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 14 | 2.5 | 1.2 | 5.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 10 | 2.5 | 1.2 | 5.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

| HCS 210 Trail | 15-01-1511-6 | 01/23/15 03:01 | Aqueous |
|----------------------|---------------------|-----------------------|----------------|
|----------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.067 | 0.10 | 0.025 | 1.00 | J | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | 1.8 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | 0.21 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | 3.9 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.13 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 71.0 | 1.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 71.0 | 1.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | 325 | 1.00 | 0.820 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | 39 | 1.0 | 0.95 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| pH (24) | 7.76 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.8 | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 17 | 0.50 | 0.24 | 1.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 5.9 | 0.50 | 0.24 | 1.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

01/24/15

Work Order:

15-01-1511

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 240 Trail | 15-01-1511-7 | 01/23/15 03:21 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.19 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | 18 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | 0.11 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.8 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | 29 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 233 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 233 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | 380 | 1.00 | 0.820 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.5 | 1.0 | 0.95 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| pH (24) | 7.64 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.6 | 0.50 | 0.24 | 1.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 3.7 | 0.50 | 0.24 | 1.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

| | | | |
|----------------------|---------------------|-----------------------|----------------|
| HCS 250 Trail | 15-01-1511-8 | 01/23/15 03:03 | Aqueous |
|----------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.23 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | 18 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | 0.12 | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.7 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | 30 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 228 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 228 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | 500 | 1.00 | 0.820 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | 3.6 | 1.0 | 0.95 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| pH (24) | 7.86 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 4.7 | 0.50 | 0.24 | 1.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 5.1 | 0.50 | 0.24 | 1.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

01/24/15

Work Order:

15-01-1511

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 260 Trail | 15-01-1511-9 | 01/23/15 03:36 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.20 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | 12 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | 0.077 | 0.10 | 0.037 | 1.00 | J | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.2 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | 25 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.11 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 155 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 155 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | 280 | 1.00 | 0.820 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | 182 | 1.00 | 0.950 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| pH (24) | 7.75 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.0 | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 11 | 0.50 | 0.24 | 1.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 4.7 | 0.50 | 0.24 | 1.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

| | | | |
|----------------------|----------------------|-----------------------|----------------|
| HCS 270 Trail | 15-01-1511-10 | 01/23/15 03:30 | Aqueous |
|----------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.20 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | 12 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | 0.082 | 0.10 | 0.037 | 1.00 | J | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.2 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | 24 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.096 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 152 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 152 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | 265 | 1.00 | 0.820 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | 114 | 1.00 | 0.950 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| pH (24) | 7.78 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 7.4 | 0.50 | 0.24 | 1.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 7.0 | 0.50 | 0.24 | 1.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

01/24/15

Work Order:

15-01-1511

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|------------------------|----------------------|-----------------------|----------------|
| FDHCS 260 Trail | 15-01-1511-11 | 01/23/15 03:36 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.19 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | 12 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | 0.073 | 0.10 | 0.037 | 1.00 | J | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.2 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | 25 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.10 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 157 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 157 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | 400 | 1.00 | 0.820 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | 89 | 1.0 | 0.95 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| pH (24) | 7.71 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.77 | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 9.5 | 0.50 | 0.24 | 1.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 5.6 | 0.50 | 0.24 | 1.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

| FDHCS 270 Trail | 15-01-1511-12 | 01/23/15 03:30 | Aqueous |
|-----------------|---------------|----------------|---------|
|-----------------|---------------|----------------|---------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Fluoride (24) | 0.18 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | 12 | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | 0.089 | 0.10 | 0.037 | 1.00 | J | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | 1.2 | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | 23 | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | 0.12 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 154 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 154 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | 275 | 1.00 | 0.820 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | 114 | 1.00 | 0.950 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| pH (24) | 7.77 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 01/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 8.2 | 0.50 | 0.24 | 1.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 7.8 | 0.50 | 0.24 | 1.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

01/24/15

Work Order:

15-01-1511

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|-------------------|---------------------|----------------|
| Method Blank | | N/A | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|-------|---------------|---------------|-----------------|
| Fluoride (24) | ND | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Chloride (24) | ND | 1.0 | 0.12 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Bromide (24) | ND | 0.10 | 0.037 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Nitrate (as N) (24) | ND | 0.10 | 0.025 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Sulfate (24) | ND | 1.0 | 0.19 | 1.00 | | mg/L | N/A | 01/24/15 | EPA 300.0 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 02/02/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 01/30/15 | SM 2320B |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.82 | 1.00 | | mg/L | 01/27/15 | 01/27/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.95 | 1.00 | | mg/L | 01/28/15 | 01/28/15 | SM 2540 D |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 02/06/15 | 02/06/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 02/07/15 | 02/07/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 01/24/15 | 01/25/15 | SM 5310 B |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| FDHCS 270 Trail | Sample | Aqueous | IC 15 | N/A | 01/24/15 16:47 | 150124S01A |
| FDHCS 270 Trail | Matrix Spike | Aqueous | IC 15 | N/A | 01/24/15 17:37 | 150124S01A |
| FDHCS 270 Trail | Matrix Spike Duplicate | Aqueous | IC 15 | N/A | 01/24/15 17:54 | 150124S01A |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.1777 | 250.0 | 241.3 | 96 | 242.4 | 97 | 80-120 | 0 | 0-20 | |
| Chloride | 11.95 | 5000 | 5111 | 102 | 5114 | 102 | 80-120 | 0 | 0-20 | |
| Bromide | ND | 500.0 | 513.1 | 103 | 513.6 | 103 | 80-120 | 0 | 0-20 | |
| Nitrate (as N) | 1.154 | 500.0 | 520.8 | 104 | 521.7 | 104 | 80-120 | 0 | 0-20 | |
| Sulfate | 23.48 | 5000 | 5083 | 101 | 5098 | 101 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|------------------|---------------------|------------|---------------|-------------------|
| FDHCS 270 Trail | Sample | | Aqueous | ACA 1 | N/A | 02/02/15 14:36 | 150202SO1A | | | |
| FDHCS 270 Trail | Matrix Spike | | Aqueous | ACA 1 | N/A | 02/02/15 14:36 | 150202SO1A | | | |
| FDHCS 270 Trail | Matrix Spike Duplicate | | Aqueous | ACA 1 | N/A | 02/02/15 14:36 | 150202SO1A | | | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Phosphorus, Total | 0.1192 | 0.2000 | 0.3152 | 98 | 0.3161 | 98 | 90-110 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|---------------------|-----------------|------------|---------------|-------------------|
| HCS 270 Peak | Sample | Aqueous | TOC 8 | 02/07/15 | 02/07/15 20:21 | F0207TOCS1 | | | | |
| HCS 270 Peak | Matrix Spike | Aqueous | TOC 8 | 02/07/15 | 02/07/15 20:21 | F0207TOCS1 | | | | |
| HCS 270 Peak | Matrix Spike Duplicate | Aqueous | TOC 8 | 02/07/15 | 02/07/15 20:21 | F0207TOCS1 | | | | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Carbon, Total Organic | 14.45 | 50.00 | 59.00 | 89 | 64.50 | 100 | 31-145 | 9 | 0-23 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HCS 270 Peak | Sample | Aqueous | TOC 8 | 01/24/15 | 01/25/15 08:45 | F0124DOCS1 |
| HCS 270 Peak | Matrix Spike | Aqueous | TOC 8 | 01/24/15 | 01/25/15 08:45 | F0124DOCS1 |
| HCS 270 Peak | Matrix Spike Duplicate | Aqueous | TOC 8 | 01/24/15 | 01/25/15 08:45 | F0124DOCS1 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|---------------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Carbon, Dissolved Organic | 10.45 | 50.00 | 60.50 | 100 | 58.00 | 95 | 31-145 | 4 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

Page 5 of 10

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HCS 210 Peak | Sample | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 17:55 | 150126SA6 |
| HCS 210 Peak | Matrix Spike | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 17:57 | 150126SA6 |
| HCS 210 Peak | Matrix Spike Duplicate | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 17:58 | 150126SA6 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Calcium | 19.19 | 0.5000 | 20.14 | 4X | 20.20 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 2.172 | 0.5000 | 2.657 | 4X | 2.674 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 1.560 | 5.000 | 6.465 | 98 | 6.337 | 96 | 83-131 | 2 | 0-7 | |
| Sodium | 1.042 | 5.000 | 5.732 | 94 | 5.907 | 97 | 73-127 | 3 | 0-9 | |
| Strontium | 0.09780 | 0.5000 | 0.6090 | 102 | 0.6223 | 105 | 81-123 | 2 | 0-6 | |
| Silicon | 1.065 | 0.5000 | 1.494 | 86 | 1.491 | 85 | 24-180 | 0 | 0-15 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|---------------------------|------------------------|-------------|----------|------------|---------------|----------------|---------------------|-----|--------|------------|
| HCS 240 Peak | Sample | | Aqueous | ICP/MS 04 | 01/26/15 | 01/28/15 11:59 | 150126S03 | | | |
| HCS 240 Peak | Matrix Spike | | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 08:25 | 150126S03 | | | |
| HCS 240 Peak | Matrix Spike Duplicate | | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 08:28 | 150126S03 | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.09735 | 97 | 0.09348 | 93 | 85-133 | 4 | 0-11 | |
| Arsenic | ND | 0.1000 | 0.09587 | 96 | 0.09385 | 94 | 73-127 | 2 | 0-11 | |
| Barium | 0.05252 | 0.1000 | 0.1547 | 102 | 0.1527 | 100 | 74-128 | 1 | 0-10 | |
| Beryllium | ND | 0.1000 | 0.09530 | 95 | 0.09080 | 91 | 56-122 | 5 | 0-11 | |
| Cadmium | ND | 0.1000 | 0.09322 | 93 | 0.08931 | 89 | 84-114 | 4 | 0-8 | |
| Chromium | ND | 0.1000 | 0.1030 | 103 | 0.1001 | 100 | 73-133 | 3 | 0-11 | |
| Copper | 0.001256 | 0.1000 | 0.09476 | 94 | 0.09231 | 91 | 72-108 | 3 | 0-10 | |
| Lead | ND | 0.1000 | 0.1001 | 100 | 0.09724 | 97 | 79-121 | 3 | 0-10 | |
| Nickel | 0.001371 | 0.1000 | 0.09097 | 90 | 0.08788 | 87 | 68-122 | 3 | 0-10 | |
| Selenium | ND | 0.1000 | 0.09183 | 92 | 0.08674 | 87 | 59-125 | 6 | 0-12 | |
| Silver | ND | 0.05000 | 0.04271 | 85 | 0.03964 | 79 | 68-128 | 7 | 0-14 | |
| Thallium | ND | 0.1000 | 0.09514 | 95 | 0.09216 | 92 | 73-121 | 3 | 0-11 | |
| Zinc | 0.01878 | 0.1000 | 0.1226 | 104 | 0.09249 | 74 | 43-145 | 28 | 0-39 | |
| Aluminum | ND | 0.1000 | 0.1196 | 120 | 0.1191 | 119 | 47-161 | 0 | 0-24 | |
| Iron | ND | 5.100 | 5.315 | 104 | 5.347 | 105 | 27-201 | 1 | 0-24 | |
| Manganese | 0.001311 | 0.1000 | 0.09499 | 94 | 0.09083 | 90 | 72-126 | 4 | 0-42 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 7470A Total
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-01-1471-1 | Sample | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 17:25 | 150130S06 |
| 15-01-1471-1 | Matrix Spike | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 17:03 | 150130S06 |
| 15-01-1471-1 | Matrix Spike Duplicate | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 17:05 | 150130S06 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | ND | 0.01000 | 0.01002 | 100 | 0.01014 | 101 | 57-141 | 1 | 0-10 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-01-1404-4 | Sample | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 13:33 | 150129S19 |
| 15-01-1404-4 | Matrix Spike | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 13:42 | 150129S19 |
| 15-01-1404-4 | Matrix Spike Duplicate | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 17:15 | 150129S19 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Caffeine | 57.25 | 100.0 | 176.7 | 119 | 171.2 | 114 | 70-130 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-01-1454-1 | Sample | Aqueous | GC/MS O | 01/24/15 | 01/24/15 21:38 | 150124S010 | | | | |
| 15-01-1454-1 | Matrix Spike | Aqueous | GC/MS O | 01/24/15 | 01/24/15 22:05 | 150124S010 | | | | |
| 15-01-1454-1 | Matrix Spike Duplicate | Aqueous | GC/MS O | 01/24/15 | 01/24/15 22:33 | 150124S010 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 50.55 | 101 | 47.70 | 95 | 74-122 | 6 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 52.34 | 105 | 49.68 | 99 | 60-144 | 5 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 49.27 | 99 | 45.96 | 92 | 73-120 | 7 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 47.18 | 94 | 43.74 | 87 | 80-122 | 8 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 43.07 | 86 | 40.50 | 81 | 70-120 | 6 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 50.32 | 101 | 46.38 | 93 | 64-142 | 8 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 47.55 | 95 | 46.35 | 93 | 52-136 | 3 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 52.85 | 106 | 49.26 | 99 | 77-125 | 7 | 0-24 | |
| Toluene | ND | 50.00 | 53.06 | 106 | 50.12 | 100 | 72-126 | 6 | 0-23 | |
| Trichloroethene | ND | 50.00 | 51.45 | 103 | 48.98 | 98 | 74-128 | 5 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 38.87 | 78 | 39.53 | 79 | 67-133 | 2 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 106.7 | 107 | 98.76 | 99 | 63-129 | 8 | 0-25 | |
| o-Xylene | ND | 50.00 | 52.89 | 106 | 48.86 | 98 | 62-128 | 8 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 40.58 | 81 | 40.37 | 81 | 68-134 | 1 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-01-1645-1 | Sample | Aqueous | GC/MS O | 01/30/15 | 01/30/15 14:17 | 150130S016 | | | | |
| 15-01-1645-1 | Matrix Spike | Aqueous | GC/MS O | 01/30/15 | 01/30/15 14:45 | 150130S016 | | | | |
| 15-01-1645-1 | Matrix Spike Duplicate | Aqueous | GC/MS O | 01/30/15 | 01/30/15 15:13 | 150130S016 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 49.06 | 98 | 51.29 | 103 | 74-122 | 4 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 49.15 | 98 | 50.31 | 101 | 60-144 | 2 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 48.72 | 97 | 50.09 | 100 | 73-120 | 3 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 47.60 | 95 | 49.24 | 98 | 80-122 | 3 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 46.84 | 94 | 49.27 | 99 | 70-120 | 5 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 48.00 | 96 | 50.12 | 100 | 64-142 | 4 | 0-20 | |
| 1,1-Dichloroethene | 2.693 | 50.00 | 45.01 | 85 | 47.39 | 89 | 52-136 | 5 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 50.89 | 102 | 52.75 | 105 | 77-125 | 4 | 0-24 | |
| Toluene | ND | 50.00 | 49.22 | 98 | 50.41 | 101 | 72-126 | 2 | 0-23 | |
| Trichloroethene | 1.449 | 50.00 | 49.02 | 95 | 50.25 | 98 | 74-128 | 2 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 40.50 | 81 | 44.71 | 89 | 67-133 | 10 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 106.3 | 106 | 108.2 | 108 | 63-129 | 2 | 0-25 | |
| o-Xylene | ND | 50.00 | 53.52 | 107 | 55.00 | 110 | 62-128 | 3 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 46.32 | 93 | 50.61 | 101 | 68-134 | 9 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|---------------------|--------------------|------------------|------------------|-----------------|-----------------------|
| HCS 240 Peak | Sample | Aqueous | ICP/MS 04 | 01/26/15 00:00 | 01/28/15 11:59 | 150126S03 |
| HCS 240 Peak | PDS | Aqueous | ICP/MS 04 | 01/26/15 00:00 | 01/27/15 08:31 | 150126S03 |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>PDS Conc.</u> | <u>PDS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Antimony | ND | 0.1000 | 0.09422 | 94 | 75-125 | |
| Arsenic | ND | 0.1000 | 0.09373 | 94 | 75-125 | |
| Barium | 0.05252 | 0.1000 | 0.1449 | 92 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.09330 | 93 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.08992 | 90 | 75-125 | |
| Chromium | ND | 0.1000 | 0.1011 | 101 | 75-125 | |
| Copper | 0.001256 | 0.1000 | 0.09289 | 92 | 75-125 | |
| Lead | ND | 0.1000 | 0.09506 | 95 | 75-125 | |
| Nickel | 0.001371 | 0.1000 | 0.08826 | 87 | 75-125 | |
| Selenium | ND | 0.1000 | 0.08392 | 84 | 75-125 | |
| Silver | ND | 0.05000 | 0.04708 | 94 | 75-125 | |
| Thallium | ND | 0.1000 | 0.09009 | 90 | 75-125 | |
| Zinc | 0.01878 | 0.1000 | 0.1007 | 82 | 75-125 | |
| Aluminum | ND | 0.1000 | 0.1196 | 120 | 75-125 | |
| Iron | ND | 5.100 | 4.713 | 92 | 75-125 | |
| Manganese | 0.001311 | 0.1000 | 0.08988 | 89 | 75-125 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---|------------------|---------------------|------------------|---------------|----------------|------------------------|
| HCS 240 Peak | Sample | Aqueous | PH1/BUR03 | N/A | 01/30/15 16:15 | F0130ALKD3 |
| HCS 240 Peak | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 01/30/15 16:15 | F0130ALKD3 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Alkalinity, Total (as CaCO ₃) | | 220.0 | 220.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|-------------------------------------|------------------|---------------------|------------------|---------------|----------------|------------------------|
| HCS 240 Peak | Sample | Aqueous | PH1/BUR03 | N/A | 01/30/15 16:15 | F0130HCOD3 |
| HCS 240 Peak | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 01/30/15 16:15 | F0130HCOD3 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Bicarbonate (as CaCO ₃) | | 220.0 | 220.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| HCS 240 Peak | Sample | Aqueous | PH1/BUR03 | N/A | 01/30/15 16:15 | F0130CO3D3 |
| HCS 240 Peak | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 01/30/15 16:15 | F0130CO3D3 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| HCS 210 Peak | Sample | Aqueous | SC 5 | 01/27/15 00:00 | 01/27/15 16:00 | F0127TDSD1 |
| HCS 210 Peak | Sample Duplicate | Aqueous | SC 5 | 01/27/15 00:00 | 01/27/15 16:00 | F0127TDSD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Dissolved | 165.0 | 160.0 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-01-1473-3 | Sample | Aqueous | N/A | 01/28/15 00:00 | 01/28/15 16:00 | F0128TSSD2 |
| 15-01-1473-3 | Sample Duplicate | Aqueous | N/A | 01/28/15 00:00 | 01/28/15 16:00 | F0128TSSD2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Suspended | 53.60 | 56.80 | 6 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------------------|------------------|---------------|----------------|------------------------|
| HCS 210 Peak | Sample | Aqueous | PH 1 | N/A | 01/24/15 11:57 | F0124PHD1 |
| HCS 210 Peak | Sample Duplicate | Aqueous | PH 1 | N/A | 01/24/15 11:57 | F0124PHD1 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| pH | | 7.170 | 7.220 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 4500 N Org B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|----------------|--------------|-----------------------|-----------------------|------------------------|
| FDHCS 270 Trail | Sample | Aqueous | BUR05 | 02/06/15 00:00 | 02/06/15 15:42 | F0206TKND1 |
| FDHCS 270 Trail | Sample Duplicate | Aqueous | BUR05 | 02/06/15 00:00 | 02/06/15 15:42 | F0206TKND1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Total Kjeldahl Nitrogen | 0.5600 | 0.6300 | 12 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|--------------------|------------------------|------------------|-----------------|-------------------|
| 099-12-906-5374 | LCS | Aqueous | IC 15 | N/A | 01/24/15 11:33 | 150124L01 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Fluoride | | 2.500 | 2.482 | 99 | 90-110 | |
| Chloride | | 50.00 | 50.36 | 101 | 90-110 | |
| Bromide | | 5.000 | 5.070 | 101 | 90-110 | |
| Nitrate (as N) | | 5.000 | 5.066 | 101 | 90-110 | |
| Sulfate | | 50.00 | 50.31 | 101 | 90-110 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 099-12-739-97 | LCS | Aqueous | ACA 1 | N/A | 02/02/15 14:36 | 150202LO1 |
| 099-12-739-97 | LCSD | Aqueous | ACA 1 | N/A | 02/02/15 14:36 | 150202LO1 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| Phosphorus, Total | 0.2000 | 0.1852 | 93 | 0.1862 | 93 | 90-110 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-859-576 | LCS | Aqueous | PH1/BUR03 | N/A | 01/30/15 16:15 | F0130ALKB3 | | | |
| 099-15-859-576 | LCSD | Aqueous | PH1/BUR03 | N/A | 01/30/15 16:15 | F0130ALKB3 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 100.0 | 98.00 | 98 | 98.00 | 98 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4401 | LCS | Aqueous | SC 5 | 01/27/15 | 01/27/15 16:00 | F0127TDSL1 | | | |
| 099-12-180-4401 | LCSD | Aqueous | SC 5 | 01/27/15 | 01/27/15 16:00 | F0127TDSL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 95.00 | 95 | 90.00 | 90 | 80-120 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-09-010-7025 | LCS | Aqueous | N/A | 01/28/15 | 01/28/15 16:00 | F0128TSSL2 | | | |
| 099-09-010-7025 | LCSD | Aqueous | N/A | 01/28/15 | 01/28/15 16:00 | F0128TSSL2 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Suspended | 100.0 | 101.0 | 101 | 98.00 | 98 | 80-120 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-097-5508 | LCS | Aqueous | TOC 8 | 02/07/15 | 02/07/15 20:21 | F0207TOCL1 | | | |
| 099-05-097-5508 | LCSD | Aqueous | TOC 8 | 02/07/15 | 02/07/15 20:21 | F0207TOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 10.00 | 9.910 | 99 | 10.50 | 105 | 80-120 | 6 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-115-1398 | LCS | Aqueous | TOC 8 | 01/24/15 | 01/25/15 08:45 | F0124DOCL1 | | | |
| 099-05-115-1398 | LCSD | Aqueous | TOC 8 | 01/24/15 | 01/25/15 08:45 | F0124DOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Dissolved Organic | 10.00 | 10.50 | 105 | 10.50 | 105 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|--------------------|------------------------|------------------|-----------------|-------------------|
| 099-15-683-1120 | LCS | Aqueous | ICP 7300 | 01/26/15 | 01/27/15 17:53 | 150126L6FF |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Calcium | | 0.5000 | 0.4279 | 86 | 80-120 | |
| Magnesium | | 0.5000 | 0.4097 | 82 | 80-120 | |
| Potassium | | 5.000 | 4.428 | 89 | 80-120 | |
| Sodium | | 5.000 | 4.045 | 81 | 80-120 | |
| Strontium | | 0.5000 | 0.4624 | 92 | 80-120 | |
| Silicon | | 0.5000 | 0.2593 | 52 | 80-120 | X |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-719 | LCS | Aqueous | ICP/MS 04 | 01/26/15 | 01/27/15 15:23 | 150126L03F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.1029 | 103 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.1052 | 105 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.09971 | 100 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.1018 | 102 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.1012 | 101 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.1053 | 105 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.09728 | 97 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.1023 | 102 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.1020 | 102 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.1022 | 102 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.05273 | 105 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.09673 | 97 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.1070 | 107 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.1175 | 118 | 80-120 | 73-127 | |
| Iron | 5.100 | 5.029 | 99 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.1022 | 102 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|------------------|
| 099-15-763-485 | LCS | Aqueous | Mercury 04 | 01/30/15 | 01/30/15 16:53 | 150130L06F |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.01000 | 0.01074 | 107 | 85-121 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|----------------|----------------|-----------------|-----------------------|------------------|
| 099-16-376-10 | LCS | Aqueous | LC/TQ 2 | 01/29/15 | 01/31/15 12:55 | 150129L19 |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Caffeine | 100.0 | 103.8 | 104 | 80-120 | |

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Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-529-778 | LCS | Aqueous | | GC 51 | 01/24/15 | 02/04/15 04:22 | 150124L08A | | | |
| 099-12-529-778 | LCSD | Aqueous | | GC 51 | 01/24/15 | 02/04/15 04:36 | 150124L08A | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Alpha-BHC | 0.5000 | 0.5324 | 106 | 0.5460 | 109 | 50-135 | 36-149 | 3 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.5466 | 109 | 0.5590 | 112 | 50-135 | 36-149 | 2 | 0-25 | |
| Beta-BHC | 0.5000 | 0.4827 | 97 | 0.4746 | 95 | 50-135 | 36-149 | 2 | 0-25 | |
| Heptachlor | 0.5000 | 0.5339 | 107 | 0.5402 | 108 | 50-135 | 36-149 | 1 | 0-25 | |
| Delta-BHC | 0.5000 | 0.6051 | 121 | 0.6324 | 126 | 50-135 | 36-149 | 4 | 0-25 | |
| Aldrin | 0.5000 | 0.4714 | 94 | 0.4712 | 94 | 50-135 | 36-149 | 0 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.5246 | 105 | 0.5276 | 106 | 50-135 | 36-149 | 1 | 0-25 | |
| Endosulfan I | 0.5000 | 0.5515 | 110 | 0.5555 | 111 | 50-135 | 36-149 | 1 | 0-25 | |
| Dieldrin | 0.5000 | 0.5378 | 108 | 0.5415 | 108 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.4811 | 96 | 0.4794 | 96 | 50-135 | 36-149 | 0 | 0-25 | |
| Endrin | 0.5000 | 0.5335 | 107 | 0.5371 | 107 | 50-135 | 36-149 | 1 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.6733 | 135 | 0.6103 | 122 | 50-135 | 36-149 | 10 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.5057 | 101 | 0.5046 | 101 | 50-135 | 36-149 | 0 | 0-25 | |
| Endosulfan II | 0.5000 | 0.5349 | 107 | 0.5386 | 108 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.5675 | 114 | 0.5887 | 118 | 50-135 | 36-149 | 4 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.5345 | 107 | 0.5392 | 108 | 50-135 | 36-149 | 1 | 0-25 | |
| Methoxychlor | 0.5000 | 0.5191 | 104 | 0.5246 | 105 | 50-135 | 36-149 | 1 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-533-1000 | LCS | Aqueous | GC 31 | 01/24/15 | 02/04/15 10:25 | 150124L09 | | | |
| 099-12-533-1000 | LCSD | Aqueous | GC 31 | 01/24/15 | 02/04/15 10:44 | 150124L09 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Aroclor-1016 | 2.000 | 1.741 | 87 | 1.763 | 88 | 50-135 | 1 | 0-25 | |
| Aroclor-1260 | 2.000 | 1.654 | 83 | 1.659 | 83 | 50-135 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-963-83 | LCS | Aqueous | | GC 26 | 01/26/15 | 02/08/15 12:52 | 150126L03 | | | |
| 099-15-963-83 | LCSD | Aqueous | | GC 26 | 01/26/15 | 02/08/15 13:37 | 150126L03 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.03948 | 99 | 0.04305 | 108 | 30-130 | 13-147 | 9 | 0-30 | |
| Bolstar | 0.04000 | 0.03654 | 91 | 0.03942 | 99 | 30-130 | 13-147 | 8 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.03777 | 94 | 0.04094 | 102 | 30-130 | 13-147 | 8 | 0-30 | |
| Coumaphos | 0.04000 | 0.03960 | 99 | 0.04209 | 105 | 30-130 | 13-147 | 6 | 0-30 | |
| Diazinon | 0.04000 | 0.03975 | 99 | 0.04272 | 107 | 30-130 | 13-147 | 7 | 0-30 | |
| Disulfoton | 0.04000 | 0.03340 | 84 | 0.03727 | 93 | 30-130 | 13-147 | 11 | 0-30 | |
| Ethoprop | 0.04000 | 0.03892 | 97 | 0.04169 | 104 | 30-130 | 13-147 | 7 | 0-30 | |
| Fensulfothion | 0.04000 | 0.04046 | 101 | 0.04286 | 107 | 30-130 | 13-147 | 6 | 0-30 | |
| Fenthion | 0.04000 | 0.03727 | 93 | 0.04031 | 101 | 30-130 | 13-147 | 8 | 0-30 | |
| Merphos | 0.04000 | 0.04891 | 122 | 0.05170 | 129 | 30-130 | 13-147 | 6 | 0-30 | |
| Methyl Parathion | 0.04000 | 0.03953 | 99 | 0.04278 | 107 | 30-130 | 13-147 | 8 | 0-30 | |
| Phorate | 0.04000 | 0.03810 | 95 | 0.04160 | 104 | 30-130 | 13-147 | 9 | 0-30 | |
| Ronnel | 0.04000 | 0.03858 | 96 | 0.04236 | 106 | 30-130 | 13-147 | 9 | 0-30 | |
| Stirophos | 0.04000 | 0.03359 | 84 | 0.03695 | 92 | 30-130 | 13-147 | 10 | 0-30 | |
| Tokuthion | 0.04000 | 0.03651 | 91 | 0.03965 | 99 | 30-130 | 13-147 | 8 | 0-30 | |
| Trichloronate | 0.04000 | 0.03693 | 92 | 0.04050 | 101 | 30-130 | 13-147 | 9 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-034-634 | LCS | Aqueous | GC 40 | 01/26/15 | 01/29/15 19:07 | 150126L04 | | | |
| 095-01-034-634 | LCSD | Aqueous | GC 40 | 01/26/15 | 01/29/15 19:30 | 150126L04 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| 2,4-D | 20.00 | 13.92 | 70 | 12.66 | 63 | 30-130 | 10 | 0-30 | |
| 2,4,5-T | 2.000 | 1.485 | 74 | 1.365 | 68 | 30-130 | 8 | 0-30 | |
| 2,4-DB | 20.00 | 13.54 | 68 | 12.15 | 61 | 30-130 | 11 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-3987 | LCS | Aqueous | | GC/MS CCC | 01/26/15 | 01/29/15 16:05 | 150126L07 | | | |
| 095-01-003-3987 | LCSD | Aqueous | | GC/MS CCC | 01/26/15 | 01/29/15 16:23 | 150126L07 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 200.0 | 197.7 | 99 | 201.5 | 101 | 61-120 | 51-130 | 2 | 0-20 | |
| Acenaphthylene | 200.0 | 201.2 | 101 | 203.8 | 102 | 55-120 | 44-131 | 1 | 0-20 | |
| Butyl Benzyl Phthalate | 200.0 | 193.7 | 97 | 193.1 | 97 | 56-122 | 45-133 | 0 | 0-20 | |
| 4-Chloro-3-Methylphenol | 200.0 | 175.8 | 88 | 177.5 | 89 | 52-120 | 41-131 | 1 | 0-20 | |
| 2-Chlorophenol | 200.0 | 172.9 | 86 | 177.3 | 89 | 47-120 | 35-132 | 2 | 0-20 | |
| 1,4-Dichlorobenzene | 200.0 | 157.5 | 79 | 165.2 | 83 | 36-120 | 22-134 | 5 | 0-20 | |
| Dimethyl Phthalate | 200.0 | 198.4 | 99 | 201.8 | 101 | 60-120 | 50-130 | 2 | 0-20 | |
| 2,4-Dinitrotoluene | 200.0 | 195.0 | 97 | 199.9 | 100 | 61-121 | 51-131 | 3 | 0-20 | |
| Fluorene | 200.0 | 206.3 | 103 | 211.0 | 106 | 67-120 | 58-129 | 2 | 0-20 | |
| N-Nitroso-di-n-propylamine | 200.0 | 174.5 | 87 | 178.5 | 89 | 39-123 | 25-137 | 2 | 0-20 | |
| Naphthalene | 200.0 | 171.5 | 86 | 175.0 | 87 | 54-120 | 43-131 | 2 | 0-20 | |
| 4-Nitrophenol | 200.0 | 86.57 | 43 | 86.38 | 43 | 14-120 | 0-138 | 0 | 0-20 | |
| Pentachlorophenol | 200.0 | 184.2 | 92 | 180.8 | 90 | 31-127 | 15-143 | 2 | 0-20 | |
| Phenol | 200.0 | 83.93 | 42 | 87.00 | 44 | 17-120 | 0-137 | 4 | 0-20 | |
| Pyrene | 200.0 | 193.4 | 97 | 194.6 | 97 | 58-124 | 47-135 | 1 | 0-20 | |
| 1,2,4-Trichlorobenzene | 200.0 | 166.5 | 83 | 171.9 | 86 | 49-120 | 37-132 | 3 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-16246 | LCS | Aqueous | GC/MS O | 01/24/15 | 01/24/15 20:14 | 150124L015 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 52.34 | 105 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 51.73 | 103 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 51.74 | 103 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 52.70 | 105 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 47.41 | 95 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 54.23 | 108 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 48.20 | 96 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 54.91 | 110 | 80-123 | 73-130 | |
| Toluene | 50.00 | 54.79 | 110 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 52.49 | 105 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 40.45 | 81 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 111.2 | 111 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 56.03 | 112 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 50.34 | 101 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 01/24/15
Work Order: 15-01-1511
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

Page 18 of 18

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-16277 | LCS | Aqueous | GC/MS O | 01/30/15 | 01/30/15 12:54 | 150130L010 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 54.43 | 109 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 52.36 | 105 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 53.66 | 107 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 52.60 | 105 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 52.68 | 105 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 53.32 | 107 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 48.58 | 97 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 56.87 | 114 | 80-123 | 73-130 | |
| Toluene | 50.00 | 54.08 | 108 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 52.56 | 105 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 45.58 | 91 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 118.2 | 118 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 60.23 | 120 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 48.98 | 98 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Sample Analysis Summary Report

Work Order: 15-01-1511

Page 1 of 1

| <u>Method</u> | <u>Extraction</u> | <u>Chemist ID</u> | <u>Instrument</u> | <u>Analytical Location</u> |
|-----------------------|-------------------|-------------------|-------------------|----------------------------|
| EPA 1694 (M) Caffeine | EPA 1694 | 262 | LC/TQ 2 | 1 |
| EPA 300.0 | N/A | 650 | IC 15 | 1 |
| EPA 365.1 | N/A | 735 | ACA 1 | 1 |
| EPA 6010B | EPA 3005A Filt. | 771 | ICP 7300 | 1 |
| EPA 6020 | EPA 3005A Filt. | 776 | ICP/MS 04 | 1 |
| EPA 7470A | EPA 7470A Filt. | 915 | Mercury 04 | 1 |
| EPA 8081A | EPA 3510C | 669 | GC 51 | 1 |
| EPA 8082 | EPA 3510C | 944 | GC 31 | 1 |
| EPA 8141A | EPA 3510C | 949 | GC 26 | 1 |
| EPA 8151A | EPA 8151A | 944 | GC 40 | 1 |
| EPA 8260B | EPA 5030C | 867 | GC/MS O | 2 |
| EPA 8270C | EPA 3510C | 923 | GC/MS CCC | 1 |
| SM 2320B | N/A | 885 | PH1/BUR03 | 1 |
| SM 2540 C | N/A | 810 | SC 5 | 1 |
| SM 2540 D | N/A | 689 | N/A | 1 |
| SM 4500 H+ B | N/A | 885 | PH 1 | 1 |
| SM 4500 N Org B | N/A | 685 | BUR05 | 1 |
| SM 5310 B | N/A | 735 | TOC 8 | 1 |

Return to Contents

Location 1: 7440 Lincoln Way, Garden Grove, CA 92841

Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841

Glossary of Terms and Qualifiers

Work Order: 15-01-1511

Page 1 of 1

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|--|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDS or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.



Calscience

7440 Lincoln Way, Garden Grove, CA 92641-1427 • (714) 895-5494
For courier service / sample drop off information, contact us26_sales@eurofinsus.com or call us.

CHAIN OF CUSTODY RECORD

WFO # / LAB USE ONLY
15-01-1511

DATE: 1/23/15
PAGE: 2 OF 3

| | | | |
|---|---------------------------|---|---------------------------------------|
| LABORATORY CLIENT: SWCA Environmental Consultants | | CLIENT PROJECT NAME / NUMBER: EAA 27122 | |
| ADDRESS: 6200 UTSA Blvd. Suite 102 | | PROJECT CONTACT: Philip Pearce | |
| CITY: San Antonio | STATE: TX | ZIP: 78249-1618 | SAMPLER(S) (PRINT): Jennifer Moreland |
| TEL: 210.877.2847 | E-MAIL: P Pearce@swca.com | P.O. NO.: | |

REQUESTED ANALYSES

| | | | |
|---|--------------|---------------|--------------|
| TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD"): | | | |
| <input type="checkbox"/> SAME DAY <input type="checkbox"/> 24 HR <input type="checkbox"/> 48 HR <input type="checkbox"/> 72 HR <input checked="" type="checkbox"/> 10 Days (standard) | | | |
| <input type="checkbox"/> COELT EDF | | LOG CODE: | |
| SPECIAL INSTRUCTIONS: | | | |
| Samples for EPA 6010B, EPA 6020, and EPA 7470 Metals are field filtered EPA 6020 Metals: Zn, Ti, Se, Sb, Pb, Ni, Cu, Cr, Cd, Be, Ba, As, Al, Ag, Fe, Mn Please also analyze each sample for pH. | | | |
| LAB. USE ONLY | SAMPLE ID | SAMPLING DATE | NO. OF CONT. |
| 1 | HCS 210 Peak | 1/22/15 16:28 | 9 |
| 2 | HCS 240 Peak | 1/22/15 16:51 | 9 |
| 3 | HCS 250 Peak | 1/22/15 16:40 | 9 |
| 4 | HCS 260 Peak | 1/22/15 17:03 | 9 |
| 5 | HCS 270 Peak | 1/22/15 16:55 | 9 |

| | | | |
|------------------------------|--------------------------------------|---------------|------------|
| Relinquished by: (Signature) | Received by: (Signature/Affiliation) | Date: 1/23/15 | Time: 5:17 |
| Relinquished by: (Signature) | Received by: (Signature/Affiliation) | Date: 1/24/15 | Time: 1:00 |
| Relinquished by: (Signature) | Received by: (Signature/Affiliation) | Date: | Time: |

ID:SATA (210) 877-2847
NC
TSA BLVD STE 102
TONIO, TX 782491618
STATES US

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ACTWGT: 52.3 LB
CAD: /OFFC1522
DIMS: 24x14x13 IN
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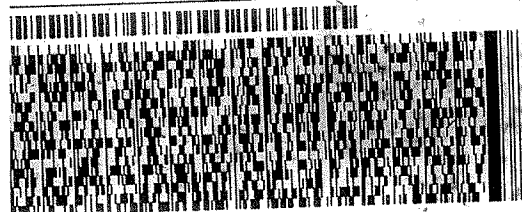
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SWCA INC
6200 UTSA BLVD STE 102
SAN ANTONIO, TX 782491618
UNITED STATES US

SHIP DATE: 23JAN15
ACTWGT: 54.2 LB
CAD: /OFFC1522
DIMS: 24x14x13 IN
BILL SENDER

Page 185 of 200

L SCIENCE
40 LINCOLN WAY

ARDEN GROVE CA 92841
95-5494 REF: DEPT:



5 of 12
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PRIORITY OVERNIGHT

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CA-US SNA

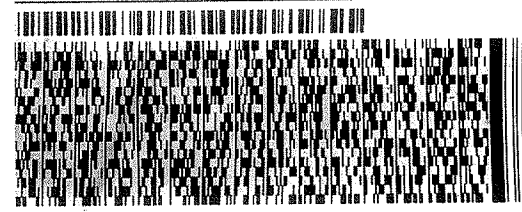


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DIMS: 24x14x13 IN
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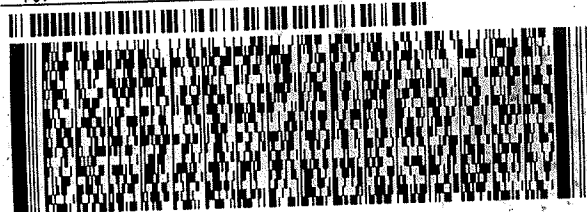
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TO
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7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 885-5494 REF: DEPT:



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Mstr# 8075 3550 1233

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CA-US SNA



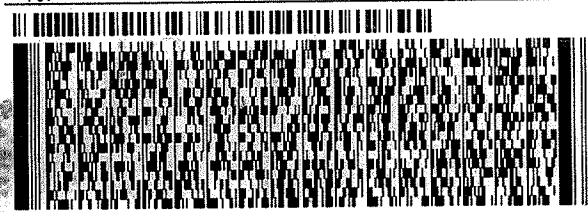
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SWCA INC
6200 UTSA BLVD STE 102
SAN ANTONIO, TX 782491618
UNITED STATES US

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ACTWGT: 54.5 LB
CAD: /OFFC1522
DIMS: 24x14x13 IN
BILL SENDER

TO
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7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 885-5494 REF: DEPT:



8 of 12
MPS# 7801 6985 1211
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Mstr# 8075 3550 1233

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PRIORITY OVERNIGHT

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UNITED STATES US

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UNITED STATES US

DIMS: 24x14x13 IN
BILL SENDER Page 186 of 200

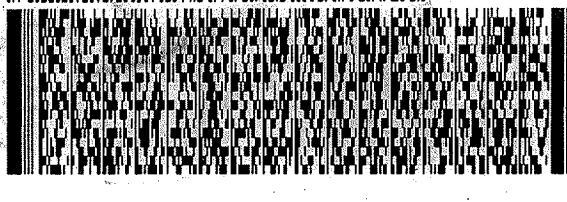
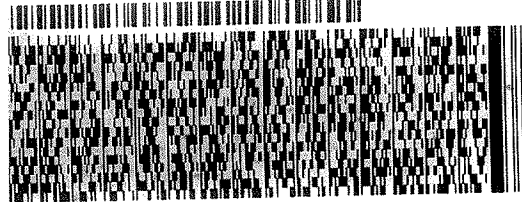
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40 LINCOLN WAY

ARDEN GROVE CA 92841
95-5494 REF: DEPT:

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7440 LINCOLN WAY

GARDEN GROVE CA 92841
(714) 895-5494 REF: DEPT:



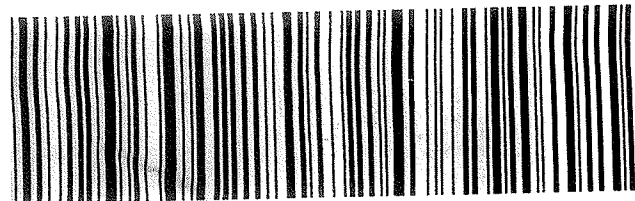
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SATURDAY 12:00P
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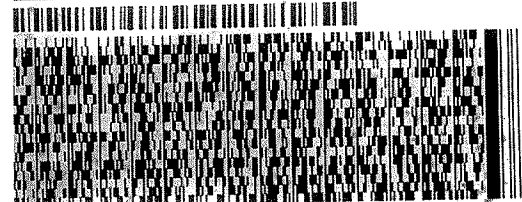
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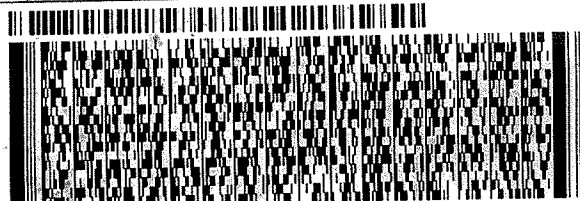


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CA-US SNA

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7440 LINCOLN WAY

GARDEN GROVE CA 92841
(714) 895-5494 REF: DEPT:



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Mstr# 8075 3550 1233

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CA-US SNA



ANTONIO, TX 782491618
UNITED STATES US

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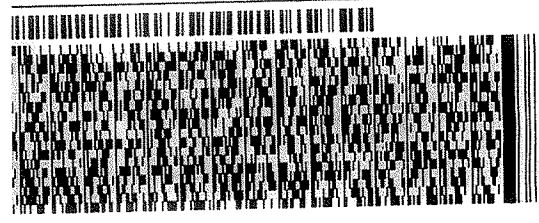
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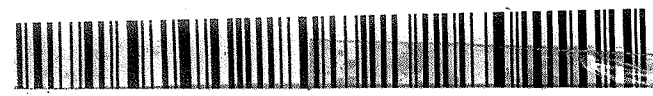
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140 LINCOLN WAY

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CA-US SNA



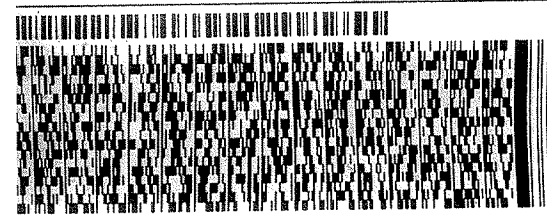
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INC
UTSA BLVD STE 102
ANTONIO, TX 782491618
UNITED STATES US

SHIP DATE: 23JAN15
ACTWGT: 59.3 LB
CAD: 7OFFC1522
DIMS: 24x14x13 IN
BILL SENDER

Part # 156297-433 R12 04/13

AL SCIENCE
140 LINCOLN WAY

GARDEN GROVE CA 92841
895-5494 REF: DEPT:



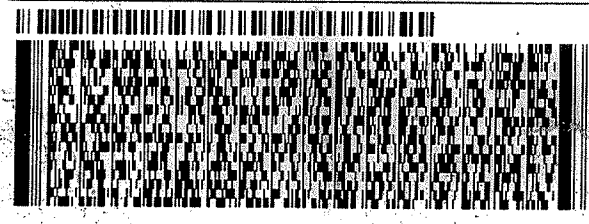
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CA-US SNA



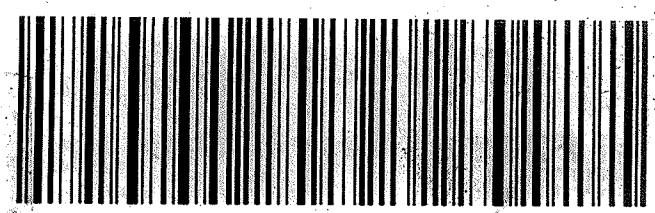
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7440 LINCOLN WAY

GARDEN GROVE CA 92841
(714) 895-5494 REF: DEPT:



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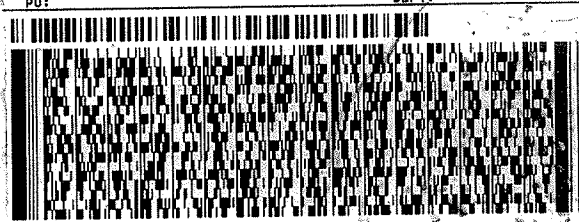


UNITED STATES US

TO

CAL SCIENCE
7440 LINCOLN WAY

GARDEN GROVE CA 92841
(714) 895-5494 REF: DEPT:



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MPS# 7801 6985 1244
0681
Mstr# 8075 3550 1233 0200
X0 APVA 92841
CA-US SNA



Part # 156297-433 R12 04/13

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Calscience

WORK ORDER #: 15-01-1111

SAMPLE RECEIPT FORM

Cooler 1 of 12

CLIENT: SWCA

DATE: 01/24/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 3.1 °C + 0.2 °C (CF) = 3.3 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 802
☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 681

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|--------------------------|-------------------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
|--|--------------------------|-------------------------------------|--------------------------|

| | | | |
|--|--------------------------|-------------------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
|--|--------------------------|-------------------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☐ ☒

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☒ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBz₂na ☐ 100PJ ☐ 100PJna₂ ☒ 5gal cube ☐ _____ ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: 141229B Labeled/Checked by: 681

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 862

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure z₂na: ZnAc₂+NaOH f: Filtered Scanned by: 862

Calscience

WORK ORDER #: 15-01-1611

SAMPLE RECEIPT FORM

Cooler 2 of 12

CLIENT: SWCA

DATE: 01/24/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 3.0 °C + 0.2 °C (CF) = 3.2 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEALS INTACT:
☒ Cooler

☐ _____

☐ No (Not Intact)

☐ Not Present

☐ N/A

Checked by: 802

☐ Sample

☐ _____

☐ No (Not Intact)

☒ Not Present

Checked by: 681

SAMPLE CONDITION:

Yes

No

N/A

Chain-Of-Custody (COC) document(s) received with samples..... ☒
☐
☐

COC document(s) received complete..... ☒
☐
☐
☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

Sampler's name indicated on COC..... ☒
☐
☐

Sample container label(s) consistent with COC..... ☒
☐
☐

Sample container(s) intact and good condition..... ☒
☐
☐

Proper containers and sufficient volume for analyses requested..... ☒
☐
☐

Analyses received within holding time..... ☒
☐
☐

Aqueous samples received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐
☐
☒

Proper preservation noted on COC or sample container..... ☒
☐
☐
☐ Unpreserved vials received for Volatiles analysis

Volatile analysis container(s) free of headspace..... ☒
☐
☐

Tedlar bag(s) free of condensation..... ☐
☐
☒
CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna2 ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna2 ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBznna ☐ 100PJ ☐ 100PJna2 ☒ 5gal cube ☐ _____ ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 681

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope

Reviewed by: 802

Preservative: h: HCL n: HNO3 na2: Na2S2O3 na: NaOH p: H3PO4 s: H2SO4 u: Ultra-pure znna: ZnAc2+NaOH f: Filtered

Scanned by: 802

Calscience

WORK ORDER #: 15-01-1611

SAMPLE RECEIPT FORM

Cooler 3 of 12

CLIENT: SWCA

DATE: 01/24/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 3.3 °C + 0.2 °C (CF) = 3.5 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 802

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 681

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☒

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBznna ☐ 100PJ ☐ 100PJna₂ ☒ 5gal cube ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 681

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope

Reviewed by: 862

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure znna: ZnAc₂+NaOH f: Filtered

Scanned by: 862

Calscience

WORK ORDER #: 15-01-1611

SAMPLE RECEIPT FORM

Cooler 4 of 12

CLIENT: SWCA

DATE: 01/24/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 3.0 °C + 0.2 °C (CF) = 3.2 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEALS INTACT:
☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 802

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 681

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☒

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBznna ☐ 100PJ ☐ 100PJna₂ ☒ 5gal cube ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 681

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 802

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure znna: ZnAc₂+NaOH f: Filtered Scanned by: 802

Calscience

WORK ORDER #: 15-01-11611

SAMPLE RECEIPT FORM

Cooler 5 of 12

CLIENT: SWCA

DATE: 01/24/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 3.1 °C + 0.2 °C (CF) = 3.3 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 802

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 681

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☒

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBznna ☐ 100PJ ☐ 100PJna₂ ☒ 5gal cube ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 681

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 802

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure znna: ZnAc₂+NaOH f: Filtered Scanned by: 802

Calscience

WORK ORDER #: 15-01-1611

SAMPLE RECEIPT FORM

Cooler 6 of 12

CLIENT: SWCA

DATE: 01/24/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 3.4 °C + 0.2 °C (CF) = 3.6 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 802

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 681

SAMPLE CONDITION:

Yes No N/A

Chain-Of-Custody (COC) document(s) received with samples..... ☒ ☐ ☐

COC document(s) received complete..... ☒ ☐ ☐
☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

Sampler's name indicated on COC..... ☒ ☐ ☐

Sample container label(s) consistent with COC..... ☒ ☐ ☐

Sample container(s) intact and good condition..... ☒ ☐ ☐

Proper containers and sufficient volume for analyses requested..... ☒ ☐ ☐

Analyses received within holding time..... ☒ ☐ ☐

Aqueous samples received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☐ ☒

Proper preservation noted on COC or sample container..... ☒ ☐ ☐
☐ Unpreserved vials received for Volatiles analysis

Volatile analysis container(s) free of headspace..... ☒ ☐ ☐

Tedlar bag(s) free of condensation..... ☐ ☐ ☒

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBznna ☐ 100PJ ☐ 100PJna₂ ☒ 5gal cube ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 681

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 802

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure znna: ZnAc₂+NaOH f: Filtered Scanned by: 802

Calscience

WORK ORDER #: 15-01-1511

SAMPLE RECEIPT FORM

Cooler 7 of 12

CLIENT: SWCA

DATE: 01/24/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 3.1 °C + 0.2 °C (CF) = 3.3 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 802

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 681

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☒

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBznna ☐ 100PJ ☐ 100PJna₂ ☒ 5gal cube ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 681

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 802

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure znna: ZnAc₂+NaOH f: Filtered Scanned by: 802

Calscience

WORK ORDER #: 15-01-1611

SAMPLE RECEIPT FORM

Cooler 8 of 12

CLIENT: SWCA

DATE: 01/24/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 3.0 °C + 0.2 °C (CF) = 3.2 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 802

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 681

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☒

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBzanna ☐ 100PJ ☐ 100PJna₂ ☒ 5gal cube ☐ _____ ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 681

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope

Reviewed by: 802

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure zanna: ZnAc₂+NaOH f: Filtered

Scanned by: 802

Calscience

WORK ORDER #: 15-01-1611

SAMPLE RECEIPT FORM

Cooler 9 of 12

CLIENT: SWCA

DATE: 01/24/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 3.3 °C + 0.2 °C (CF) = 3.5 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 802

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 681

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☒

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBz₂na ☐ 100PJ ☐ 100PJna₂ ☒ 5gal cube ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 681

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 802

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure z₂na: ZnAc₂+NaOH f: Filtered Scanned by: 802

Calscience

WORK ORDER #: 15-01-1611

SAMPLE RECEIPT FORM

Cooler 10 of 12

CLIENT: SWCA

DATE: 01/24/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 3.1 °C + 0.2 °C (CF) = 3.3 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 802

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 681

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☒

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBznna ☐ 100PJ ☐ 100PJna₂ ☒ 5gal cube ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 681

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 802

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure znna: ZnAc₂+NaOH f: Filtered Scanned by: 802

Calscience

WORK ORDER #: 15-01-1511

SAMPLE RECEIPT FORM

Cooler 11 of 12

CLIENT: SWCA

DATE: 01/24/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Temperature 3.2°C + 0.2°C (CF) = 3.4°C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 802

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 681

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| <input type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfides <input type="checkbox"/> Dissolved Oxygen..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBznna ☐ 100PJ ☐ 100PJna₂ ☒ 5gal cube ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 681

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 802

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure znna: ZnAc₂+NaOH f: Filtered Scanned by: 802

Calscience

WORK ORDER #: 15-01-1511

SAMPLE RECEIPT FORM

Cooler 12 of 12

CLIENT: SWCA

DATE: 01/24/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 3.1 °C + 0.2 °C (CF) = 3.3 °C ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

☐ Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 802

CUSTODY SEALS INTACT:

☒ Cooler ☐ _____ ☐ No (Not Intact) ☐ Not Present ☐ N/A Checked by: 802

☐ Sample ☐ _____ ☐ No (Not Intact) ☒ Not Present Checked by: 681

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|--------------------------|--------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.

☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.

| | | | |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples received within 15-minute holding time

☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen..... ☐ ☐ ☒

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☐ Unpreserved vials received for Volatiles analysis

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® ☐ TerraCores® ☐ _____

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs

☐ 500AGB ☐ 500AGJ ☐ 500AGJs ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 1PB ☐ 1PBna ☐ 500PB

☐ 250PB ☒ 250PBn ☐ 125PB ☐ 125PBznna ☐ 100PJ ☐ 100PJna₂ ☒ Seal Cube ☐ _____ ☐ _____

Air: ☐ Tedlar® ☐ Canister Other: ☐ _____ Trip Blank Lot#: _____ Labeled/Checked by: 681

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 802

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure znna: ZnAc₂+NaOH f: Filtered Scanned by: 802



Calscience

WORK ORDER #: 15-01-1511

SAMPLE ANOMALY FORM

SAMPLES - CONTAINERS & LABELS:

Comments:

- ☐ Sample(s) NOT RECEIVED but listed on COC
☐ Sample(s) received but NOT LISTED on COC
☐ Holding time expired – list sample ID(s) and test
☐ Insufficient quantities for analysis – list test
☐ Improper container(s) used – list test
☐ Improper preservative used – list test
☐ No preservative noted on COC or label – list test & notify lab
☐ Sample labels illegible – note test/container type
☒ Sample label(s) do not match COC – Note in comments
 - ☒ Sample ID
 - ☒ Date and/or Time Collected
 - ☐ Project Information
 - ☐ # of Container(s)
 - ☐ Analysis☒ Sample container(s) compromised – Note in comments
 - ☐ Water present in sample container
 - ☒ Broken☐ Sample container(s) not labeled
☐ Air sample container(s) compromised – Note in comments
 - ☐ Flat
 - ☐ Very low in volume
 - ☐ Leaking (Not transferred - duplicate bag submitted)
 - ☐ Leaking (transferred into Calscience Tedlar® Bag*)
 - ☐ Leaking (transferred into Client's Tedlar® Bag*)☐ Other: _____

(-4) collection time per label
is 17:10.

(-11) labeled as HC5260 Trail
collection date & time matched.

(-11) 1 of 2 250 CGB W. H₂SO₄
Received broken.

HEADSPACE – Containers with Bubble > 6mm or ¼ inch:

| Sample # | Container ID(s) | # of Vials Received | Sample # | Container ID(s) | # of Vials Received | Sample # | Container ID(s) | # of Cont. received | Analysis |
|----------|-----------------|---------------------|----------|-----------------|---------------------|----------|-----------------|---------------------|----------|
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |

Comments: _____

*Transferred at Client's request.

Initial / Date: 681 01/24/15

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HSM 210 LEAD
Sample Number: AA94367
Sample Matrix: Non Potable Water

Collection Date/Time: 5/5/15 22:10
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94367-A | E. coli | √ | SM 9223B-2004 | 77 | MPN/100 mL | H | 1 | 42654 | 5/6/15 | 12:58 | RSC |
| AA94367-A | E. Coli Holding Time - IDEXX Colilert | | NA | 14.80 | hours | | 0.00 | 42653 | 5/6/15 | 12:58 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42654

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HSM 230 LEAD
Sample Number: AA94368
Sample Matrix: Non Potable Water

Collection Date/Time: 5/5/15 22:20
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94368-A | E. coli | √ | SM 9223B-2004 | 20000 | MPN/100 mL | H | 1 | 42654 | 5/6/15 | 12:58 | RSC |
| AA94368-A | E. Coli Holding Time - IDEXX Colilert | | NA | 14.63 | hours | | 0.00 | 42653 | 5/6/15 | 12:58 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42654

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 231 LEAD
Sample Number: AA94369
Sample Matrix: Non Potable Water

Collection Date/Time: 5/5/15 22:11
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94369-A | E. coli | √ | SM 9223B-2004 | 870 | MPN/100 mL | H | 1 | 42654 | 5/6/15 | 12:58 | RSC |
| AA94369-A | E. Coli Holding Time - IDEXX Colilert | | NA | 14.78 | hours | | 0.00 | 42653 | 5/6/15 | 12:58 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42654

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 240 LEAD
Sample Number: AA94370
Sample Matrix: Non Potable Water

Collection Date/Time: 5/5/15 22:49
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed. Limited sample volume.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94370-A | E. coli | √ | SM 9223B-2004 | 24000 | MPN/100 mL | H | 1 | 42654 | 5/6/15 | 12:58 | RSC |
| AA94370-A | E. Coli Holding Time - IDEXX Colilert | | NA | 14.15 | hours | | 0.00 | 42653 | 5/6/15 | 12:58 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42654

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 250 LEAD
Sample Number: AA94371
Sample Matrix: Non Potable Water

Collection Date/Time: 5/5/15 22:27
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94371-A | E. coli | √ | SM 9223B-2004 | 2400 | MPN/100 mL | H | 1 | 42654 | 5/6/15 | 12:58 | RSC |
| AA94371-A | E. Coli Holding Time - IDEXX Colilert | | NA | 14.52 | hours | | 0.00 | 42653 | 5/6/15 | 12:58 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42654

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 260 LEAD
Sample Number: AA94372
Sample Matrix: Non Potable Water

Collection Date/Time: 5/5/15 22:46
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94372-A | E. coli | √ | SM 9223B-2004 | 2400 | MPN/100 mL | H | 1 | 42654 | 5/6/15 | 12:58 | RSC |
| AA94372-A | E. Coli Holding Time - IDEXX Colilert | | NA | 14.20 | hours | | 0.00 | 42653 | 5/6/15 | 12:58 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42654

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 270 LEAD
Sample Number: AA94373
Sample Matrix: Non Potable Water

Collection Date/Time: 5/5/15 23:06
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94373-A | E. coli | √ | SM 9223B-2004 | 1400 | MPN/100 mL | H | 1 | 42654 | 5/6/15 | 12:58 | RSC |
| AA94373-A | E. Coli Holding Time - IDEXX Colilert | | NA | 13.87 | hours | | 0.00 | 42653 | 5/6/15 | 12:58 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42654

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 210 PEAK
Sample Number: AA94374
Sample Matrix: Non Potable Water

Collection Date/Time: 5/5/15 23:24
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94374-A | E. coli | √ | SM 9223B-2004 | 66 | MPN/100 mL | H | 1 | 42654 | 5/6/15 | 12:58 | RSC |
| AA94374-A | E. Coli Holding Time - IDEXX Colilert | | NA | 13.57 | hours | | 0.00 | 42653 | 5/6/15 | 12:58 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42654

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 230 PEAK
Sample Number: AA94375
Sample Matrix: Non Potable Water

Collection Date/Time: 5/5/15 23:35
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94375-A | E. coli | √ | SM 9223B-2004 | 14000 | MPN/100 mL | H | 1 | 42654 | 5/6/15 | 12:58 | RSC |
| AA94375-A | E. Coli Holding Time - IDEXX Colilert | | NA | 13.38 | hours | | 0.00 | 42653 | 5/6/15 | 12:58 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42654

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 231 PEAK
Sample Number: AA94376
Sample Matrix: Non Potable Water

Collection Date/Time: 5/5/15 22:11
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94376-A | E. coli | √ | SM 9223B-2004 | 770 | MPN/100 mL | H | 1 | 42656 | 5/6/15 | 13:17 | RSC |
| AA94376-A | E. Coli Holding Time - IDEXX Colilert | | NA | 15.10 | hours | | 0.00 | 42655 | 5/6/15 | 13:17 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42656

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 240 PEAK
Sample Number: AA94377
Sample Matrix: Non Potable Water

Collection Date/Time: 5/6/15 00:05
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94377-A | E. coli | √ | SM 9223B-2004 | 290 | MPN/100 mL | H | 1 | 42656 | 5/6/15 | 13:17 | RSC |
| AA94377-A | E. Coli Holding Time - IDEXX Colilert | | NA | 13.20 | hours | | 0.00 | 42655 | 5/6/15 | 13:17 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42656

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 250 PEAK
Sample Number: AA94378
Sample Matrix: Non Potable Water

Collection Date/Time: 5/5/15 23:30
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94378-A | E. coli | √ | SM 9223B-2004 | 2200 | MPN/100 mL | H | 1 | 42656 | 5/6/15 | 13:17 | RSC |
| AA94378-A | E. Coli Holding Time - IDEXX Colilert | | NA | 13.78 | hours | | 0.00 | 42655 | 5/6/15 | 13:17 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42656

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 260 PEAK
Sample Number: AA94379
Sample Matrix: Non Potable Water

Collection Date/Time: 5/5/15 23:49
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94379-A | E. coli | √ | SM 9223B-2004 | 2100 | MPN/100 mL | H | 1 | 42656 | 5/6/15 | 13:17 | RSC |
| AA94379-A | E. Coli Holding Time - IDEXX Colilert | | NA | 13.47 | hours | | 0.00 | 42655 | 5/6/15 | 13:17 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42656

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 270 PEAK
Sample Number: AA94380
Sample Matrix: Non Potable Water

Collection Date/Time: 5/6/15 00:06
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94380-A | E. coli | √ | SM 9223B-2004 | 5800 | MPN/100 mL | H | 1 | 42656 | 5/6/15 | 13:17 | RSC |
| AA94380-A | E. Coli Holding Time - IDEXX Colilert | | NA | 13.18 | hours | | 0.00 | 42655 | 5/6/15 | 13:17 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42656

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 210 TRAIL
Sample Number: AA94381
Sample Matrix: Non Potable Water

Collection Date/Time: 5/6/15 01:30
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94381-A | E. coli | √ | SM 9223B-2004 | 270 | MPN/100 mL | H | 1 | 42656 | 5/6/15 | 13:17 | RSC |
| AA94381-A | E. Coli Holding Time - IDEXX Colilert | | NA | 11.78 | hours | | 0.00 | 42655 | 5/6/15 | 13:17 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42656

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: FDHSM 210 TRAIL
Sample Number: AA94382
Sample Matrix: Non Potable Water

Collection Date/Time: 5/6/15 01:30
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Hold time exceeded, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94382-A | E. coli | √ | SM 9223B-2004 | 330 | MPN/100 mL | H | 1 | 42656 | 5/6/15 | 13:17 | RSC |
| AA94382-A | E. Coli Holding Time - IDEXX Colilert | | NA | 11.78 | hours | | 0.00 | 42655 | 5/6/15 | 13:17 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42656

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 230 TRAIL
Sample Number: AA94383
Sample Matrix: Non Potable Water

Collection Date/Time: 5/6/15 02:15
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Analysis Comments: AA94383-A E. coli

Customer submitted samples with a very limited hold time. Analyst set up samples for analysis as soon as possible. Customer requested the continuation of the analysis.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94383-A | E. coli | √ | SM 9223B-2004 | 5200 | MPN/100 mL | H* | 1 | 42656 | 5/6/15 | 13:17 | RSC |
| AA94383-A | E. Coli Holding Time - IDEXX Colilert | | NA | 11.03 | hours | | 0.00 | 42655 | 5/6/15 | 13:17 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42656

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: FDHSM 230 TRAIL
Sample Number: AA94384
Sample Matrix: Non Potable Water

Collection Date/Time: 5/6/15 02:15
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Analysis Comments: AA94384-A E. coli

Customer submitted samples with a very limited hold time. Analyst set up samples for analysis as soon as possible. Customer requested the continuation of the analysis.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94384-A | E. coli | √ | SM 9223B-2004 | 4600 | MPN/100 mL | H* | 1 | 42656 | 5/6/15 | 13:17 | RSC |
| AA94384-A | E. Coli Holding Time - IDEXX Colilert | | NA | 11.03 | hours | | 0.00 | 42655 | 5/6/15 | 13:17 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42656

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 231 TRAIL
Sample Number: AA94385
Sample Matrix: Non Potable Water

Collection Date/Time: 5/6/15 02:40
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Analysis Comments: AA94385-A E. coli

Customer submitted samples with a very limited hold time. Analyst set up samples for analysis as soon as possible. Customer requested the continuation of the analysis.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94385-A | E. coli | √ | SM 9223B-2004 | 280 | MPN/100 mL | H* | 1 | 42658 | 5/6/15 | 13:34 | RSC |
| AA94385-A | E. Coli Holding Time - IDEXX Colilert | | NA | 10.90 | hours | | 0.00 | 42657 | 5/6/15 | 13:34 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42658

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: FDHSM 231 TRAIL
Sample Number: AA94386
Sample Matrix: Non Potable Water

Collection Date/Time: 5/6/15 02:40
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Analysis Comments: AA94386-A E. coli

Customer submitted samples with a very limited hold time. Analyst set up samples for analysis as soon as possible. Customer requested the continuation of the analysis.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94386-A | E. coli | √ | SM 9223B-2004 | 330 | MPN/100 mL | H* | 1 | 42658 | 5/6/15 | 13:34 | RSC |
| AA94386-A | E. Coli Holding Time - IDEXX Colilert | | NA | 10.90 | hours | | 0.00 | 42657 | 5/6/15 | 13:34 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42658

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 240 TRAIL
Sample Number: AA94387
Sample Matrix: Non Potable Water

Collection Date/Time: 5/6/15 03:05
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Analysis Comments: AA94387-A E. coli

Customer submitted samples with a very limited hold time. Analyst set up samples for analysis as soon as possible. Customer requested the continuation of the analysis.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94387-A | E. coli | √ | SM 9223B-2004 | 210 | MPN/100 mL | H* | 1 | 42658 | 5/6/15 | 13:34 | RSC |
| AA94387-A | E. Coli Holding Time - IDEXX Colilert | | NA | 10.48 | hours | | 0.00 | 42657 | 5/6/15 | 13:34 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42658

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 250 TRAIL
Sample Number: AA94388
Sample Matrix: Non Potable Water

Collection Date/Time: 5/6/15 03:58
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Analysis Comments: AA94388-A E. coli

Customer submitted samples with a very limited hold time. Analyst set up samples for analysis as soon as possible. Customer requested the continuation of the analysis.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94388-A | E. coli | √ | SM 9223B-2004 | 2500 | MPN/100 mL | H* | 1 | 42658 | 5/6/15 | 13:34 | RSC |
| AA94388-A | E. Coli Holding Time - IDEXX Colilert | | NA | 9.60 | hours | | 0.00 | 42657 | 5/6/15 | 13:34 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42658

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM 260 TRAIL
Sample Number: AA94389
Sample Matrix: Non Potable Water

Collection Date/Time: 5/6/15 04:20
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Analysis Comments: AA94389-A E. coli

Customer submitted samples with a very limited hold time. Analyst set up samples for analysis as soon as possible. Customer requested the continuation of the analysis.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94389-A | E. coli | √ | SM 9223B-2004 | 1100 | MPN/100 mL | H* | 1 | 42658 | 5/6/15 | 13:34 | RSC |
| AA94389-A | E. Coli Holding Time - IDEXX Colilert | | NA | 9.23 | hours | | 0.00 | 42657 | 5/6/15 | 13:34 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42658

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HSM 270 TRAIL
Sample Number: AA94390
Sample Matrix: Non Potable Water

Collection Date/Time: 5/6/15 04:40
Receipt Date/Time: 5/6/15 09:35

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact David Hernandez, Laboratory Supervisor, at (210) 302-3669

Analysis identified with a "✓" complies with NELAP requirements unless otherwise specified in the case narrative.

Analysis Comments: AA94390-A E. coli

Customer submitted samples with a very limited hold time. Analyst set up samples for analysis as soon as possible. Customer requested the continuation of the analysis.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA94390-A | E. coli | ✓ | SM 9223B-2004 | 1700 | MPN/100 mL | H* | 1 | 42658 | 5/6/15 | 13:34 | RSC |
| AA94390-A | E. Coli Holding Time - IDEXX Colilert | | NA | 8.90 | hours | | 0.00 | 42657 | 5/6/15 | 13:34 | RSC |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



May 11, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-42658

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Jeanette Hernandez
Water Quality Planner / QAO

5/11/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable



WORK ORDER NUMBER: 15-05-0402

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: SWCA Environmental Consultants

Client Project Name: EAA 27122

Attention: Philip Pearce
6200 UTSA Blvd.
Suite 102
San Antonio, TX 78249-1618

A handwritten signature in black ink, appearing to read "Don Burley".

Approved for release on 06/01/2015 by:
Don Burley
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

Contents

Client Project Name: EAA 27122
Work Order Number: 15-05-0402

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Work Order Narrative

Work Order: 15-05-0402

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Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 05/07/15. They were assigned to Work Order 15-05-0402.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.



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Sample Summary

| | | |
|--|-----------------------|----------------|
| Client: SWCA Environmental Consultants | Work Order: | 15-05-0402 |
| 6200 UTSA Blvd., Suite 102 | Project Name: | EAA 27122 |
| San Antonio, TX 78249-1618 | PO Number: | |
| | Date/Time Received: | 05/07/15 10:30 |
| | Number of Containers: | 217 |
| Attn: Philip Pearce | | |

| Sample Identification | Lab Number | Collection Date and Time | Number of Containers | Matrix |
|-----------------------|---------------|--------------------------|----------------------|---------|
| HSM 210 Lead | 15-05-0402-1 | 05/05/15 22:10 | 9 | Aqueous |
| HSM 230 Lead | 15-05-0402-2 | 05/05/15 22:20 | 9 | Aqueous |
| HSM 231 Lead | 15-05-0402-3 | 05/05/15 22:11 | 9 | Aqueous |
| HSM 240 Lead | 15-05-0402-4 | 05/05/15 22:49 | 9 | Aqueous |
| HSM 250 Lead | 15-05-0402-5 | 05/05/15 22:27 | 9 | Aqueous |
| HSM 260 Lead | 15-05-0402-6 | 05/05/15 22:46 | 9 | Aqueous |
| HSM 270 Lead | 15-05-0402-7 | 05/05/15 23:06 | 9 | Aqueous |
| HSM 210 Peak | 15-05-0402-8 | 05/05/15 23:24 | 9 | Aqueous |
| HSM 230 Peak | 15-05-0402-9 | 05/05/15 23:35 | 9 | Aqueous |
| HSM 231 Peak | 15-05-0402-10 | 05/05/15 22:11 | 9 | Aqueous |
| HSM 240 Peak | 15-05-0402-11 | 05/06/15 00:05 | 9 | Aqueous |
| HSM 250 Peak | 15-05-0402-12 | 05/05/15 23:30 | 9 | Aqueous |
| HSM 260 Peak | 15-05-0402-13 | 05/05/15 23:49 | 9 | Aqueous |
| HSM 270 Peak | 15-05-0402-14 | 05/06/15 00:06 | 9 | Aqueous |
| TB06 | 15-05-0402-15 | 05/06/15 10:11 | 1 | Aqueous |
| HSM 210 Trail | 15-05-0402-16 | 05/06/15 01:30 | 9 | Aqueous |
| HSM 230 Trail | 15-05-0402-17 | 05/06/15 02:15 | 9 | Aqueous |
| HSM 231 Trail | 15-05-0402-18 | 05/06/15 02:40 | 9 | Aqueous |
| HSM 240 Trail | 15-05-0402-19 | 05/06/15 03:05 | 9 | Aqueous |
| HSM 250 Trail | 15-05-0402-20 | 05/06/15 03:58 | 9 | Aqueous |
| HSM 260 Trail | 15-05-0402-21 | 05/06/15 04:20 | 9 | Aqueous |
| HSM 270 Trail | 15-05-0402-22 | 05/06/15 04:40 | 9 | Aqueous |
| FDHSM 210 Trail | 15-05-0402-23 | 05/06/15 01:30 | 9 | Aqueous |
| FDHSM 230 Trail | 15-05-0402-24 | 05/06/15 02:15 | 9 | Aqueous |
| FDHSM 231 Trail | 15-05-0402-25 | 05/06/15 02:40 | 9 | Aqueous |

[Return to Contents](#)



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

Page 1 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| HSM 210 Lead | 15-05-0402-1-I | 05/05/15 22:10 | Aqueous | IC 15 | N/A | 05/07/15 14:53 | 150507L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.21 | 0.10 | 0.025 | 1.00 | |
| Chloride | 26 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.15 | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | 0.38 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 36 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 230 Lead | 15-05-0402-2-I | 05/05/15 22:20 | Aqueous | IC 15 | N/A | 05/07/15 15:09 | 150507L01 |
|---------------------|-----------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.096 | 0.10 | 0.025 | 1.00 | J |
| Chloride | 7.2 | 1.0 | 0.12 | 1.00 | |
| Bromide | ND | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | 0.79 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 12 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 231 Lead | 15-05-0402-3-I | 05/05/15 22:11 | Aqueous | IC 15 | N/A | 05/07/15 15:26 | 150507L01 |
|---------------------|-----------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.17 | 0.10 | 0.025 | 1.00 | |
| Chloride | 16 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.073 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.0 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 24 | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| HSM 240 Lead | 15-05-0402-4-I | 05/05/15 22:49 | Aqueous | IC 15 | N/A | 05/07/15 15:43 | 150507L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.15 | 0.10 | 0.025 | 1.00 | |
| Chloride | 11 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.045 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 0.68 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 16 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 250 Lead | 15-05-0402-5-I | 05/05/15 22:27 | Aqueous | IC 15 | N/A | 05/07/15 16:00 | 150507L01 |
|---------------------|-----------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.18 | 0.10 | 0.025 | 1.00 | |
| Chloride | 17 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.081 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.0 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 25 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 260 Lead | 15-05-0402-6-I | 05/05/15 22:46 | Aqueous | IC 15 | N/A | 05/07/15 16:17 | 150507L01 |
|---------------------|-----------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.18 | 0.10 | 0.025 | 1.00 | |
| Chloride | 18 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.091 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.1 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 27 | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| HSM 270 Lead | 15-05-0402-7-I | 05/05/15 23:06 | Aqueous | IC 15 | N/A | 05/07/15 16:34 | 150507L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.17 | 0.10 | 0.025 | 1.00 | |
| Chloride | 18 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.092 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.1 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 27 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 210 Peak | 15-05-0402-8-I | 05/05/15 23:24 | Aqueous | IC 15 | N/A | 05/07/15 16:50 | 150507L01 |
|---------------------|-----------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.21 | 0.10 | 0.025 | 1.00 | |
| Chloride | 26 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.14 | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | 0.38 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 36 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 230 Peak | 15-05-0402-9-I | 05/05/15 23:35 | Aqueous | IC 15 | N/A | 05/07/15 17:59 | 150507L01 |
|---------------------|-----------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.16 | 0.10 | 0.025 | 1.00 | |
| Chloride | 17 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.082 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.4 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 31 | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| HSM 231 Peak | 15-05-0402-10-I | 05/05/15 22:11 | Aqueous | IC 15 | N/A | 05/07/15 18:16 | 150507L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.18 | 0.10 | 0.025 | 1.00 | |
| Chloride | 18 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.091 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.1 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 28 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 240 Peak | 15-05-0402-11-I | 05/06/15 00:05 | Aqueous | IC 15 | N/A | 05/07/15 18:32 | 150507L01 |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.18 | 0.10 | 0.025 | 1.00 | |
| Chloride | 19 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.088 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.1 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 28 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 250 Peak | 15-05-0402-12-I | 05/05/15 23:30 | Aqueous | IC 15 | N/A | 05/07/15 18:49 | 150507L01 |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.22 | 0.10 | 0.025 | 1.00 | |
| Chloride | 20 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.11 | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | 1.0 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 29 | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| HSM 260 Peak | 15-05-0402-13-I | 05/05/15 23:49 | Aqueous | IC 15 | N/A | 05/07/15 19:06 | 150507L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.18 | 0.10 | 0.025 | 1.00 | |
| Chloride | 17 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.092 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.1 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 25 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 270 Peak | 15-05-0402-14-I | 05/06/15 00:06 | Aqueous | IC 15 | N/A | 05/07/15 19:23 | 150507L01 |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.18 | 0.10 | 0.025 | 1.00 | |
| Chloride | 17 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.073 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.0 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 25 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 210 Trail | 15-05-0402-16-I | 05/06/15 01:30 | Aqueous | IC 15 | N/A | 05/07/15 19:40 | 150507L01 |
|----------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.22 | 0.10 | 0.025 | 1.00 | |
| Chloride | 26 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.14 | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | 0.41 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 35 | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Trail | 15-05-0402-17-I | 05/06/15 02:15 | Aqueous | IC 15 | N/A | 05/07/15 19:57 | 150507L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.16 | 0.10 | 0.025 | 1.00 | |
| Chloride | 17 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.076 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.5 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 27 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|-------|-----|----------------|-----------|
| HSM 231 Trail | 15-05-0402-18-I | 05/06/15 02:40 | Aqueous | IC 15 | N/A | 05/07/15 20:14 | 150507L01 |
|---------------|-----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.19 | 0.10 | 0.025 | 1.00 | |
| Chloride | 19 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.094 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.2 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 28 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|-------|-----|----------------|-----------|
| HSM 240 Trail | 15-05-0402-19-I | 05/06/15 03:05 | Aqueous | IC 15 | N/A | 05/07/15 21:38 | 150507L02 |
|---------------|-----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.20 | 0.10 | 0.025 | 1.00 | |
| Chloride | 19 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.092 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.2 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 28 | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-05-0402-20-I | 05/06/15 03:58 | Aqueous | IC 15 | N/A | 05/07/15 21:55 | 150507L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.19 | 0.10 | 0.025 | 1.00 | |
| Chloride | 18 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.089 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.1 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 27 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|-------|-----|----------------|-----------|
| HSM 260 Trail | 15-05-0402-21-I | 05/06/15 04:20 | Aqueous | IC 15 | N/A | 05/07/15 22:12 | 150507L02 |
|---------------|-----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.17 | 0.10 | 0.025 | 1.00 | |
| Chloride | 18 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.097 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.1 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 27 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|-------|-----|----------------|-----------|
| HSM 270 Trail | 15-05-0402-22-I | 05/06/15 04:40 | Aqueous | IC 15 | N/A | 05/07/15 22:29 | 150507L02 |
|---------------|-----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.18 | 0.10 | 0.025 | 1.00 | |
| Chloride | 18 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.099 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.1 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 27 | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| FDHSM 210 Trail | 15-05-0402-23-I | 05/06/15 01:30 | Aqueous | IC 15 | N/A | 05/07/15 22:45 | 150507L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.21 | 0.10 | 0.025 | 1.00 | |
| Chloride | 25 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.16 | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | 0.41 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 35 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|------------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| FDHSM 230 Trail | 15-05-0402-24-I | 05/06/15 02:15 | Aqueous | IC 15 | N/A | 05/07/15 23:02 | 150507L02 |
|------------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.16 | 0.10 | 0.025 | 1.00 | |
| Chloride | 17 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.079 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.5 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 27 | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|------------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| FDHSM 231 Trail | 15-05-0402-25-I | 05/06/15 02:40 | Aqueous | IC 15 | N/A | 05/07/15 23:19 | 150507L02 |
|------------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.19 | 0.10 | 0.025 | 1.00 | |
| Chloride | 19 | 1.0 | 0.12 | 1.00 | |
| Bromide | 0.097 | 0.10 | 0.037 | 1.00 | J |
| Nitrate (as N) | 1.2 | 0.10 | 0.025 | 1.00 | |
| Sulfate | 28 | 1.0 | 0.19 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-906-5739 | N/A | Aqueous | IC 15 | N/A | 05/07/15 11:05 | 150507L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | ND | 0.10 | 0.025 | 1.00 | |
| Chloride | ND | 1.0 | 0.12 | 1.00 | |
| Bromide | ND | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | ND | 0.10 | 0.025 | 1.00 | |
| Sulfate | ND | 1.0 | 0.19 | 1.00 | |

| | | | | | | | |
|--------------|-----------------|-----|---------|-------|-----|----------------|-----------|
| Method Blank | 099-12-906-5725 | N/A | Aqueous | IC 15 | N/A | 05/07/15 21:04 | 150507L02 |
|--------------|-----------------|-----|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | ND | 0.10 | 0.025 | 1.00 | |
| Chloride | ND | 1.0 | 0.12 | 1.00 | |
| Bromide | ND | 0.10 | 0.037 | 1.00 | |
| Nitrate (as N) | ND | 0.10 | 0.025 | 1.00 | |
| Sulfate | ND | 1.0 | 0.19 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

Page 1 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 210 Lead | 15-05-0402-1-D | 05/05/15 22:10 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:16 | 150507LA5 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 93.1 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 20.6 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.21 | 0.500 | 0.103 | 1.00 | |
| Sodium | 17.3 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.742 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.34 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 230 Lead | 15-05-0402-2-D | 05/05/15 22:20 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:18 | 150507LA5 |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 40.1 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 4.77 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 3.24 | 0.500 | 0.103 | 1.00 | |
| Sodium | 6.32 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.161 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 2.01 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 231 Lead | 15-05-0402-3-D | 05/05/15 22:11 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:20 | 150507LA5 |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 83.0 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 15.2 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.85 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.507 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.91 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 240 Lead | 15-05-0402-4-D | 05/05/15 22:49 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:21 | 150507LA5 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 52.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 9.34 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.62 | 0.500 | 0.103 | 1.00 | |
| Sodium | 9.39 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.321 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 2.46 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 250 Lead | 15-05-0402-5-D | 05/05/15 22:27 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:23 | 150507LA5 |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 86.4 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 15.6 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.67 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.533 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.94 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 260 Lead | 15-05-0402-6-D | 05/05/15 22:46 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:25 | 150507LA5 |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 92.5 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.1 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.60 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.7 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.578 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.26 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

Page 3 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 270 Lead | 15-05-0402-7-D | 05/05/15 23:06 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:31 | 150507LA5 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 91.3 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.1 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.13 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.572 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.31 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 210 Peak | 15-05-0402-8-D | 05/05/15 23:24 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:33 | 150507LA5 |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 87.4 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 19.8 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.71 | 0.500 | 0.103 | 1.00 | |
| Sodium | 16.0 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.686 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.20 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 230 Peak | 15-05-0402-9-D | 05/05/15 23:35 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:35 | 150507LA5 |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 85.9 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 10.9 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.85 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.6 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.380 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.91 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 231 Peak | 15-05-0402-10-D | 05/05/15 22:11 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:36 | 150507LA5 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 97.6 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.8 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.60 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.598 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.46 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 240 Peak | 15-05-0402-11-D | 05/06/15 00:05 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:38 | 150507LA5 |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 95.6 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.8 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.86 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.5 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.598 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.43 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 250 Peak | 15-05-0402-12-D | 05/05/15 23:30 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:40 | 150507LA5 |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 90.6 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 16.7 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.68 | 0.500 | 0.103 | 1.00 | |
| Sodium | 15.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.574 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.28 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

Page 5 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 260 Peak | 15-05-0402-13-D | 05/05/15 23:49 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:42 | 150507LA5 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 87.5 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 16.4 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.06 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.9 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.549 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.18 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 270 Peak | 15-05-0402-14-D | 05/06/15 00:06 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:43 | 150507LA5 |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 80.9 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 14.7 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.10 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.507 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 3.73 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 210 Trail | 15-05-0402-16-D | 05/06/15 01:30 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:45 | 150507LA5 |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 90.1 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 20.3 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.29 | 0.500 | 0.103 | 1.00 | |
| Sodium | 16.5 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.723 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.45 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Trail | 15-05-0402-17-D | 05/06/15 02:15 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:47 | 150507LA5 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 83.8 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 12.1 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.52 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.404 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.01 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|----------|----------|----------------|-----------|
| HSM 231 Trail | 15-05-0402-18-D | 05/06/15 02:40 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:53 | 150507LA5 |
|---------------|-----------------|----------------|---------|----------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 98.7 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 18.4 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.53 | 0.500 | 0.103 | 1.00 | |
| Sodium | 14.0 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.608 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.64 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|----------|----------|----------------|-----------|
| HSM 240 Trail | 15-05-0402-19-D | 05/06/15 03:05 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:55 | 150507LA5 |
|---------------|-----------------|----------------|---------|----------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 90.9 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.0 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.32 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.565 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.22 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-05-0402-20-D | 05/06/15 03:58 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:57 | 150507LA5 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 93.9 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.1 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.87 | 0.500 | 0.103 | 1.00 | |
| Sodium | 14.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.585 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.31 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|----------|----------|----------------|-----------|
| HSM 260 Trail | 15-05-0402-21-D | 05/06/15 04:20 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:58 | 150507LA5 |
|---------------|-----------------|----------------|---------|----------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 96.4 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.9 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.68 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.587 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.50 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|----------|----------|----------------|------------|
| HSM 270 Trail | 15-05-0402-22-D | 05/06/15 04:40 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 22:00 | 150507LA6A |
|---------------|-----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 93.1 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.3 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.75 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.569 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.41 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| FDHSM 210 Trail | 15-05-0402-23-D | 05/06/15 01:30 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 22:02 | 150507LA6A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 91.7 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 20.1 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 2.06 | 0.500 | 0.103 | 1.00 | |
| Sodium | 16.5 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.715 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.41 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|------------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| FDHSM 230 Trail | 15-05-0402-24-D | 05/06/15 02:15 | Aqueous | ICP 7300 | 05/07/15 | 05/08/15 19:05 | 150507LA6A |
|------------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 86.8 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 12.5 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 2.05 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.9 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.411 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.11 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|------------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| FDHSM 231 Trail | 15-05-0402-25-D | 05/06/15 02:40 | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 22:05 | 150507LA6A |
|------------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 97.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 18.4 | 0.100 | 0.00336 | 1.00 | B |
| Potassium | 1.48 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.5 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.597 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.65 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-683-1264 | N/A | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 20:59 | 150507LA5 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | ND | 0.100 | 0.0118 | 1.00 | |
| Magnesium | ND | 0.100 | 0.00336 | 1.00 | |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|-----------------|-----|---------|----------|----------|----------------|------------|
| Method Blank | 099-15-683-1265 | N/A | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:01 | 150507LA6A |
|--------------|-----------------|-----|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|--------|---------|------|------------|
| Calcium | ND | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 0.00476 | 0.100 | 0.00336 | 1.00 | J |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Lead | 15-05-0402-1-D | 05/05/15 22:10 | Aqueous | ICP/MS 04 | 05/07/15 | 05/11/15 21:21 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000130 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00112 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0396 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000802 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | 0.000112 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00206 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0117 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00539 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0556 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0150 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 2 of 26

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Lead | 15-05-0402-2-D | 05/05/15 22:20 | Aqueous | ICP/MS 04 | 05/07/15 | 05/11/15 21:25 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000679 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00192 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0230 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.00158 | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00565 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000566 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00182 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0540 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.102 | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.151 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0132 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|------------------|-----------------|-----------------------|-------------------|
| HSM 231 Lead | 15-05-0402-3-D | 05/05/15 22:11 | Aqueous | ICP/MS 04 | 05/07/15 | 05/11/15 22:17 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Antimony | 0.000297 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.000935 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | 0.0355 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000550 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00284 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000276 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00202 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000445 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0248 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0298 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0846 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00472 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Lead | 15-05-0402-4-D | 05/05/15 22:49 | Aqueous | ICP/MS 04 | 05/07/15 | 05/11/15 22:21 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000745 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00120 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0290 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.00123 | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00355 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000370 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00181 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000229 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0240 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0156 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0937 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0175 | 0.00100 | 0.000139 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Lead | 15-05-0402-5-D | 05/05/15 22:27 | Aqueous | ICP/MS 04 | 05/07/15 | 05/11/15 22:25 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000288 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.000936 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | 0.0398 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000411 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00265 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000148 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00267 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000335 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0159 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0113 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0880 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00595 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Lead | 15-05-0402-6-D | 05/05/15 22:46 | Aqueous | ICP/MS 04 | 05/07/15 | 05/11/15 22:29 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000147 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.000899 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | 0.0415 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00207 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000159 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00256 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000377 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0159 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00800 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0754 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00445 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Lead | 15-05-0402-7-D | 05/05/15 23:06 | Aqueous | ICP/MS 04 | 05/07/15 | 05/11/15 22:33 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000147 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.000616 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | 0.0405 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00148 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000122 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00254 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000395 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0158 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00784 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0708 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00432 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Peak | 15-05-0402-8-D | 05/05/15 23:24 | Aqueous | ICP/MS 04 | 05/07/15 | 05/11/15 22:37 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000112 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00116 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0381 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000793 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00219 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000250 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00991 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00555 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0707 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0154 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Peak | 15-05-0402-9-D | 05/05/15 23:35 | Aqueous | ICP/MS 04 | 05/07/15 | 05/11/15 22:41 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000457 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00128 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0403 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000721 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00366 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000162 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00238 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000370 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0249 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0124 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.148 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0114 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 231 Peak | 15-05-0402-10-D | 05/05/15 22:11 | Aqueous | ICP/MS 04 | 05/07/15 | 05/11/15 22:45 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000176 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00128 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0408 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00131 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000128 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00235 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000324 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0144 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0108 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.109 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00258 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Peak | 15-05-0402-11-D | 05/06/15 00:05 | Aqueous | ICP/MS 04 | 05/07/15 | 05/11/15 22:49 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000177 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00153 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0409 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00109 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000110 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00224 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000305 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0179 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0123 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0821 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00186 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Peak | 15-05-0402-12-D | 05/05/15 23:30 | Aqueous | ICP/MS 04 | 05/07/15 | 05/11/15 22:53 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000269 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.00121 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0433 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00139 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000238 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00229 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000340 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0345 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00858 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.104 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0109 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Peak | 15-05-0402-13-D | 05/05/15 23:49 | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 16:59 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000219 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0359 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00118 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000164 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00237 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000214 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0225 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0308 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0884 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00520 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Peak | 15-05-0402-14-D | 05/06/15 00:06 | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 17:03 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000137 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0346 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00159 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000119 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00214 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000405 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0111 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00994 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0781 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00467 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-05-0402-16-D | 05/06/15 01:30 | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 17:07 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0362 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000804 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00239 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000394 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0106 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00574 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0724 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0184 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Trail | 15-05-0402-17-D | 05/06/15 02:15 | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 17:11 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000248 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0359 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000560 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00229 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000122 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00234 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000308 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0308 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0119 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0929 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00719 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 231 Trail | 15-05-0402-18-D | 05/06/15 02:40 | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 17:15 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0382 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000402 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.000596 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00256 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000386 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0180 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00844 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0695 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00172 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-05-0402-19-D | 05/06/15 03:05 | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 17:55 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0356 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000909 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00245 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000279 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00931 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00686 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0921 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00159 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-05-0402-20-D | 05/06/15 03:58 | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 17:59 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0373 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000766 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00261 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000369 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0123 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00740 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0719 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00241 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Trail | 15-05-0402-21-D | 05/06/15 04:20 | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 18:03 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0380 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00133 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000119 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00268 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000250 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0168 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00937 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.145 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00383 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Trail | 15-05-0402-22-D | 05/06/15 04:40 | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 01:00 | 150507LA6F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000132 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.000754 | 0.00100 | 0.000386 | 1.00 | B,J |
| Barium | 0.0421 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00116 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000125 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00234 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000295 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0214 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00632 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0778 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00423 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 210 Trail | 15-05-0402-23-D | 05/06/15 01:30 | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 00:48 | 150507LA6F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | 0.00124 | 0.00100 | 0.000386 | 1.00 | B |
| Barium | 0.0397 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000945 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00207 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000263 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0110 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00347 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0708 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0203 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 230 Trail | 15-05-0402-24-D | 05/06/15 02:15 | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 00:52 | 150507LA6F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000244 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | 0.000733 | 0.00100 | 0.000386 | 1.00 | B,J |
| Barium | 0.0387 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000634 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00275 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000124 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00235 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000317 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0278 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0117 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0758 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00682 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 231 Trail | 15-05-0402-25-D | 05/06/15 02:40 | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 00:56 | 150507LA6F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | 0.000941 | 0.00100 | 0.000386 | 1.00 | B,J |
| Barium | 0.0430 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00105 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000146 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00225 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000319 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0361 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0860 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00175 | 0.00100 | 0.000139 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-820 | N/A | Aqueous | ICP/MS 04 | 05/07/15 | 05/09/15 02:52 | 150507LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-821 | N/A | Aqueous | ICP/MS 04 | 05/07/15 | 05/09/15 02:56 | 150507LA6F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | 0.000469 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

Page 1 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 210 Lead | 15-05-0402-1-D | 05/05/15 22:10 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:18 | 150511L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 230 Lead | 15-05-0402-2-D | 05/05/15 22:20 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:20 | 150511L04 |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 231 Lead | 15-05-0402-3-D | 05/05/15 22:11 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:22 | 150511L04 |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 240 Lead | 15-05-0402-4-D | 05/05/15 22:49 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:24 | 150511L04 |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 250 Lead | 15-05-0402-5-D | 05/05/15 22:27 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:27 | 150511L04 |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 260 Lead | 15-05-0402-6-D | 05/05/15 22:46 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:29 | 150511L04 |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

Page 2 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 270 Lead | 15-05-0402-7-D | 05/05/15 23:06 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:31 | 150511L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 210 Peak | 15-05-0402-8-D | 05/05/15 23:24 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:33 | 150511L04 |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 230 Peak | 15-05-0402-9-D | 05/05/15 23:35 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:41 | 150511L04 |
|---------------------|-----------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 231 Peak | 15-05-0402-10-D | 05/05/15 22:11 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:43 | 150511L04 |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 240 Peak | 15-05-0402-11-D | 05/06/15 00:05 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:46 | 150511L04 |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 250 Peak | 15-05-0402-12-D | 05/05/15 23:30 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:48 | 150511L04 |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 260 Peak | 15-05-0402-13-D | 05/05/15 23:49 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:50 | 150511L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 270 Peak | 15-05-0402-14-D | 05/06/15 00:06 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:52 | 150511L04 |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 210 Trail | 15-05-0402-16-D | 05/06/15 01:30 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:54 | 150511L04 |
|----------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 230 Trail | 15-05-0402-17-D | 05/06/15 02:15 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:57 | 150511L04 |
|----------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 231 Trail | 15-05-0402-18-D | 05/06/15 02:40 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:59 | 150511L04 |
|----------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|
| HSM 240 Trail | 15-05-0402-19-D | 05/06/15 03:05 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 19:01 | 150511L04 |
|----------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-05-0402-20-D | 05/06/15 03:58 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 19:08 | 150511L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|------------|----------|----------------|-----------|
| HSM 260 Trail | 15-05-0402-21-D | 05/06/15 04:20 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 19:10 | 150511L04 |
|---------------|-----------------|----------------|---------|------------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|------------|----------|----------------|-----------|
| HSM 270 Trail | 15-05-0402-22-D | 05/06/15 04:40 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 19:12 | 150511L05 |
|---------------|-----------------|----------------|---------|------------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|-----------|
| FDHSM 210 Trail | 15-05-0402-23-D | 05/06/15 01:30 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 19:14 | 150511L05 |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|-----------|
| FDHSM 230 Trail | 15-05-0402-24-D | 05/06/15 02:15 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 19:17 | 150511L05 |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|-----------|
| FDHSM 231 Trail | 15-05-0402-25-D | 05/06/15 02:40 | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 19:19 | 150511L05 |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-763-554 | N/A | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 17:55 | 150511L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|-----|---------|------------|----------|----------------|-----------|
| Method Blank | 099-15-763-555 | N/A | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 17:58 | 150511L05 |
|--------------|----------------|-----|---------|------------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|-----------------------|-----------------------|----------------|-----------------------|-----------------|-----------------------|-------------------|
| HSM 210 Lead | 15-05-0402-1-I | 05/05/15 22:10 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 13:18 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 10 | | 9.8 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 43 | | 31-200 | | | |
| HSM 230 Lead | 15-05-0402-2-I | 05/05/15 22:20 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 18:25 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 4800 | | 200 | | 20.0 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 47 | | 31-200 | | | |
| HSM 231 Lead | 15-05-0402-3-I | 05/05/15 22:11 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 18:33 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 370 | | 190 | | 20.0 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 71 | | 31-200 | | | |
| HSM 240 Lead | 15-05-0402-4-I | 05/05/15 22:49 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 18:42 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 3100 | | 200 | | 20.0 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 62 | | 31-200 | | | |
| HSM 250 Lead | 15-05-0402-5-I | 05/05/15 22:27 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 18:50 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 350 | | 97 | | 10.0 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 66 | | 31-200 | | | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|------------------------|-----------------------|----------------|-----------------------|-----------------|-----------------------|-------------------|
| HSM 260 Lead | 15-05-0402-6-I | 05/05/15 22:46 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 14:00 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 230 | | 9.9 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 40 | | 31-200 | | | |
| HSM 270 Lead | 15-05-0402-7-I | 05/05/15 23:06 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 14:08 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 180 | | 9.9 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 40 | | 31-200 | | | |
| HSM 210 Peak | 15-05-0402-8-I | 05/05/15 23:24 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 14:17 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 17 | | 9.7 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 39 | | 31-200 | | | |
| HSM 230 Peak | 15-05-0402-9-I | 05/05/15 23:35 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 19:01 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 830 | | 98 | | 10.0 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 57 | | 31-200 | | | |
| HSM 231 Peak | 15-05-0402-10-I | 05/05/15 22:11 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 14:33 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 83 | | 9.8 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 40 | | 31-200 | | | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|------------------------|-----------------------|----------------|-----------------------|-----------------|-----------------------|-------------------|
| HSM 240 Peak | 15-05-0402-11-I | 05/06/15 00:05 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 14:42 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 44 | | 9.9 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 37 | | 31-200 | | | |
| HSM 250 Peak | 15-05-0402-12-I | 05/05/15 23:30 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 14:50 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 230 | | 9.4 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 40 | | 31-200 | | | |
| HSM 260 Peak | 15-05-0402-13-I | 05/05/15 23:49 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 19:10 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 870 | | 48 | | 5.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 51 | | 31-200 | | | |
| HSM 270 Peak | 15-05-0402-14-I | 05/06/15 00:06 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 19:18 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 450 | | 49 | | 5.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 57 | | 31-200 | | | |
| HSM 210 Trail | 15-05-0402-16-I | 05/06/15 01:30 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 15:15 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 13 | | 9.9 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 46 | | 31-200 | | | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|------------------------|-----------------------|-----------------------|----------------|-------------------|-----------------------|-------------------|
| HSM 230 Trail | 15-05-0402-17-I | 05/06/15 02:15 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 19:27 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | <u>RL</u> | | <u>DF</u> | | <u>Qualifiers</u> |
| Caffeine | | 290 | 49 | | 5.00 | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | <u>Control Limits</u> | | <u>Qualifiers</u> | | |
| Caffeine-C13 (Surrogate) | | 58 | 31-200 | | | | |
| HSM 231 Trail | 15-05-0402-18-I | 05/06/15 02:40 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 15:32 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | <u>RL</u> | | <u>DF</u> | | <u>Qualifiers</u> |
| Caffeine | | 25 | 9.6 | | 1.00 | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | <u>Control Limits</u> | | <u>Qualifiers</u> | | |
| Caffeine-C13 (Surrogate) | | 40 | 31-200 | | | | |
| HSM 240 Trail | 15-05-0402-19-I | 05/06/15 03:05 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 15:40 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | <u>RL</u> | | <u>DF</u> | | <u>Qualifiers</u> |
| Caffeine | | 25 | 9.7 | | 1.00 | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | <u>Control Limits</u> | | <u>Qualifiers</u> | | |
| Caffeine-C13 (Surrogate) | | 48 | 31-200 | | | | |
| HSM 250 Trail | 15-05-0402-20-I | 05/06/15 03:58 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 15:48 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | <u>RL</u> | | <u>DF</u> | | <u>Qualifiers</u> |
| Caffeine | | 130 | 9.6 | | 1.00 | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | <u>Control Limits</u> | | <u>Qualifiers</u> | | |
| Caffeine-C13 (Surrogate) | | 44 | 31-200 | | | | |
| HSM 260 Trail | 15-05-0402-21-I | 05/06/15 04:20 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 15:59 | 150512L22 |
| <u>Parameter</u> | | <u>Result</u> | <u>RL</u> | | <u>DF</u> | | <u>Qualifiers</u> |
| Caffeine | | 110 | 9.8 | | 1.00 | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | <u>Control Limits</u> | | <u>Qualifiers</u> | | |
| Caffeine-C13 (Surrogate) | | 46 | 31-200 | | | | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|------------------------|-----------------------|----------------|-----------------------|-----------------|-----------------------|-------------------|
| HSM 270 Trail | 15-05-0402-22-I | 05/06/15 04:40 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 16:08 | 150512L22 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 190 | | 9.7 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 47 | | 31-200 | | | |
| FDHSM 210 Trail | 15-05-0402-23-I | 05/06/15 01:30 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 16:16 | 150512L22 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 14 | | 9.9 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 35 | | 31-200 | | | |
| FDHSM 230 Trail | 15-05-0402-24-I | 05/06/15 02:15 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 19:35 | 150512L22 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 340 | | 20 | | 2.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 47 | | 31-200 | | | |
| FDHSM 231 Trail | 15-05-0402-25-I | 05/06/15 02:40 | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 16:42 | 150512L22 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | 25 | | 9.8 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 46 | | 31-200 | | | |
| Method Blank | 099-16-376-13 | N/A | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 12:36 | 150512L21 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | | <u>DF</u> | <u>Qualifiers</u> |
| Caffeine | | ND | | 10 | | 1.00 | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 76 | | 31-200 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine
Units: ng/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------|-------------------|---------------------|---------|-----------------------|---------------|--------------------|-------------------|
| Method Blank | 099-16-376-14 | N/A | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 17:18 | 150512L22 |
| <u>Parameter</u> | | <u>Result</u> | | <u>RL</u> | <u>DF</u> | | <u>Qualifiers</u> |
| Caffeine | | ND | | 10 | 1.00 | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> | | <u>Control Limits</u> | | <u>Qualifiers</u> | |
| Caffeine-C13 (Surrogate) | | 77 | | 31-200 | | | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 210 Lead | 15-05-0402-1-I | 05/05/15 22:10 | Aqueous | GC 44 | 05/08/15 | 05/11/15 12:48 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.098 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.098 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.098 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.098 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.098 | 0.028 | 1.00 | |
| Aldrin | ND | 0.098 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.098 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.098 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.098 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.098 | 0.026 | 1.00 | |
| Endrin | ND | 0.098 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.098 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.098 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.098 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.098 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.098 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.098 | 0.025 | 1.00 | |
| Chlordane | ND | 0.98 | 0.32 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.58 | 1.00 | |
| Endrin Ketone | ND | 0.098 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 57 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 56 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Lead | 15-05-0402-2-I | 05/05/15 22:20 | Aqueous | GC 44 | 05/08/15 | 05/11/15 13:03 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.098 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.098 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.098 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.098 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.098 | 0.028 | 1.00 | |
| Aldrin | ND | 0.098 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.098 | 0.025 | 1.00 | |
| Endosulfan I | 0.048 | 0.098 | 0.027 | 1.00 | J |
| Dieldrin | ND | 0.098 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.098 | 0.026 | 1.00 | |
| Endrin | ND | 0.098 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.098 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.098 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.098 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.098 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.098 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.098 | 0.025 | 1.00 | |
| Chlordane | ND | 0.98 | 0.32 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.58 | 1.00 | |
| Endrin Ketone | ND | 0.098 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 99 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 102 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Lead | 15-05-0402-3-I | 05/05/15 22:11 | Aqueous | GC 44 | 05/08/15 | 05/11/15 13:17 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 83 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 91 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Lead | 15-05-0402-4-F | 05/05/15 22:49 | Aqueous | GC 44 | 05/08/15 | 05/11/15 13:31 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | 0.045 | 0.10 | 0.028 | 1.00 | J |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 90 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 95 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 250 Lead | 15-05-0402-5-I | 05/05/15 22:27 | Aqueous | GC 44 | 05/08/15 | 05/11/15 13:45 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 86 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 94 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Lead | 15-05-0402-6-I | 05/05/15 22:46 | Aqueous | GC 44 | 05/08/15 | 05/11/15 14:00 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.097 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.097 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.097 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.097 | 0.028 | 1.00 | |
| Aldrin | ND | 0.097 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.097 | 0.024 | 1.00 | |
| Endosulfan I | ND | 0.097 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.097 | 0.028 | 1.00 | |
| 4,4'-DDE | ND | 0.097 | 0.026 | 1.00 | |
| Endrin | ND | 0.097 | 0.030 | 1.00 | |
| Endrin Aldehyde | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan II | ND | 0.097 | 0.026 | 1.00 | |
| 4,4'-DDT | ND | 0.097 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.097 | 0.028 | 1.00 | |
| Methoxychlor | ND | 0.097 | 0.024 | 1.00 | |
| Chlordane | ND | 0.97 | 0.32 | 1.00 | |
| Toxaphene | ND | 1.9 | 0.57 | 1.00 | |
| Endrin Ketone | ND | 0.097 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 87 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 93 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 270 Lead | 15-05-0402-7-I | 05/05/15 23:06 | Aqueous | GC 44 | 05/08/15 | 05/11/15 14:14 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 87 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 92 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 210 Peak | 15-05-0402-8-I | 05/05/15 23:24 | Aqueous | GC 44 | 05/08/15 | 05/11/15 16:17 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 80 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 88 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 230 Peak | 15-05-0402-9-I | 05/05/15 23:35 | Aqueous | GC 44 | 05/08/15 | 05/11/15 16:32 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 75 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 90 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Peak | 15-05-0402-10-I | 05/05/15 22:11 | Aqueous | GC 44 | 05/08/15 | 05/11/15 16:46 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.096 | 0.027 | 1.00 | |
| Gamma-BHC | ND | 0.096 | 0.029 | 1.00 | |
| Beta-BHC | ND | 0.096 | 0.029 | 1.00 | |
| Heptachlor | ND | 0.096 | 0.025 | 1.00 | |
| Delta-BHC | ND | 0.096 | 0.027 | 1.00 | |
| Aldrin | ND | 0.096 | 0.026 | 1.00 | |
| Heptachlor Epoxide | ND | 0.096 | 0.024 | 1.00 | |
| Endosulfan I | ND | 0.096 | 0.027 | 1.00 | |
| Dieldrin | ND | 0.096 | 0.027 | 1.00 | |
| 4,4'-DDE | ND | 0.096 | 0.026 | 1.00 | |
| Endrin | ND | 0.096 | 0.029 | 1.00 | |
| Endrin Aldehyde | ND | 0.096 | 0.025 | 1.00 | |
| 4,4'-DDD | ND | 0.096 | 0.026 | 1.00 | |
| Endosulfan II | ND | 0.096 | 0.026 | 1.00 | |
| 4,4'-DDT | ND | 0.096 | 0.026 | 1.00 | |
| Endosulfan Sulfate | ND | 0.096 | 0.028 | 1.00 | |
| Methoxychlor | ND | 0.096 | 0.024 | 1.00 | |
| Chlordane | ND | 0.96 | 0.32 | 1.00 | |
| Toxaphene | ND | 1.9 | 0.57 | 1.00 | |
| Endrin Ketone | ND | 0.096 | 0.023 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 84 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 93 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Peak | 15-05-0402-11-I | 05/06/15 00:05 | Aqueous | GC 44 | 05/08/15 | 05/11/15 17:07 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 0 | 50-135 | 2,6 |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 250 Peak | 15-05-0402-12-I | 05/05/15 23:30 | Aqueous | GC 44 | 05/08/15 | 05/11/15 17:22 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 80 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 93 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 260 Peak | 15-05-0402-13-I | 05/05/15 23:49 | Aqueous | GC 44 | 05/08/15 | 05/11/15 17:36 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 55 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 61 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 270 Peak | 15-05-0402-14-I | 05/06/15 00:06 | Aqueous | GC 44 | 05/08/15 | 05/11/15 17:50 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 79 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-05-0402-16-I | 05/06/15 01:30 | Aqueous | GC 44 | 05/08/15 | 05/11/15 18:04 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 77 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Trail | 15-05-0402-17-I | 05/06/15 02:15 | Aqueous | GC 44 | 05/08/15 | 05/11/15 18:19 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | 0.038 | 0.10 | 0.027 | 1.00 | J |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 79 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 91 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 231 Trail | 15-05-0402-18-I | 05/06/15 02:40 | Aqueous | GC 44 | 05/08/15 | 05/11/15 18:33 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 71 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 78 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-05-0402-19-I | 05/06/15 03:05 | Aqueous | GC 44 | 05/08/15 | 05/11/15 18:47 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 64 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 71 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-05-0402-20-I | 05/06/15 03:58 | Aqueous | GC 44 | 05/08/15 | 05/11/15 19:02 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 67 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 70 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Trail | 15-05-0402-21-I | 05/06/15 04:20 | Aqueous | GC 44 | 05/08/15 | 05/11/15 19:16 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 83 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 95 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Trail | 15-05-0402-22-I | 05/06/15 04:40 | Aqueous | GC 44 | 05/09/15 | 05/11/15 20:13 | 150509L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 75 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 92 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| FDHSM 210 Trail | 15-05-0402-23-I | 05/06/15 01:30 | Aqueous | GC 44 | 05/09/15 | 05/11/15 20:27 | 150509L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 60 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 75 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 230 Trail | 15-05-0402-24-I | 05/06/15 02:15 | Aqueous | GC 44 | 05/09/15 | 05/11/15 20:42 | 150509L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 77 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 103 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 231 Trail | 15-05-0402-25-I | 05/06/15 02:40 | Aqueous | GC 44 | 05/09/15 | 05/11/15 20:56 | 150509L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 90 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 96 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-803 | N/A | Aqueous | GC 44 | 05/08/15 | 05/11/15 12:34 | 150508L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 83 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-804 | N/A | Aqueous | GC 44 | 05/09/15 | 05/11/15 19:59 | 150509L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 84 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 108 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 210 Lead | 15-05-0402-1-I | 05/05/15 22:10 | Aqueous | GC 31 | 05/08/15 | 05/11/15 18:57 | 150508L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 0.98 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.98 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 0.98 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.98 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 0.98 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.98 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.98 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.98 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 80 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 60 | 50-135 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Lead | 15-05-0402-2-I | 05/05/15 22:20 | Aqueous | GC 31 | 05/08/15 | 05/12/15 15:45 | 150508L15 |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 0.98 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.98 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 0.98 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.98 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 0.98 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.98 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.98 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.98 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 90 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 91 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Lead | 15-05-0402-3-I | 05/05/15 22:11 | Aqueous | GC 31 | 05/08/15 | 05/12/15 16:04 | 150508L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 91 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 85 | 50-135 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Lead | 15-05-0402-4-I | 05/05/15 22:49 | Aqueous | GC 31 | 05/08/15 | 05/12/15 16:42 | 150508L15 |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 89 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 74 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 250 Lead | 15-05-0402-5-I | 05/05/15 22:27 | Aqueous | GC 31 | 05/08/15 | 05/12/15 17:20 | 150508L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 0.99 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.99 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 0.99 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 0.99 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 0.99 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.99 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.99 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.99 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 102 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 87 | 50-135 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Lead | 15-05-0402-6-I | 05/05/15 22:46 | Aqueous | GC 31 | 05/08/15 | 05/12/15 17:40 | 150508L15 |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 0.97 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.97 | 0.27 | 1.00 | |
| Aroclor-1232 | ND | 0.97 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.97 | 0.17 | 1.00 | |
| Aroclor-1248 | ND | 0.97 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.97 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.97 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.97 | 0.25 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 105 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 89 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 270 Lead | 15-05-0402-7-I | 05/05/15 23:06 | Aqueous | GC 31 | 05/08/15 | 05/12/15 17:59 | 150508L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 0.99 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 0.99 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 0.99 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 0.99 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 0.99 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 0.99 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.99 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 0.99 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 103 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 87 | 50-135 | |

| | | | | | | | |
|---------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 210 Peak | 15-05-0402-8-I | 05/05/15 23:24 | Aqueous | GC 31 | 05/08/15 | 05/12/15 18:18 | 150508L15 |
|---------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 103 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 83 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Peak | 15-05-0402-9-I | 05/05/15 23:35 | Aqueous | GC 31 | 05/08/15 | 05/12/15 18:37 | 150508L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 98 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Peak | 15-05-0402-10-I | 05/05/15 22:11 | Aqueous | GC 31 | 05/08/15 | 05/12/15 18:56 | 150508L15 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 0.96 | 0.28 | 1.00 | |
| Aroclor-1221 | ND | 0.96 | 0.27 | 1.00 | |
| Aroclor-1232 | ND | 0.96 | 0.24 | 1.00 | |
| Aroclor-1242 | ND | 0.96 | 0.17 | 1.00 | |
| Aroclor-1248 | ND | 0.96 | 0.19 | 1.00 | |
| Aroclor-1254 | ND | 0.96 | 0.22 | 1.00 | |
| Aroclor-1260 | ND | 0.96 | 0.25 | 1.00 | |
| Aroclor-1262 | ND | 0.96 | 0.25 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 101 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 84 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Peak | 15-05-0402-11-I | 05/06/15 00:05 | Aqueous | GC 31 | 05/08/15 | 05/12/15 19:15 | 150508L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 100 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 82 | 50-135 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 250 Peak | 15-05-0402-12-I | 05/05/15 23:30 | Aqueous | GC 31 | 05/08/15 | 05/12/15 19:34 | 150508L15 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 98 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 83 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Peak | 15-05-0402-13-I | 05/05/15 23:49 | Aqueous | GC 31 | 05/08/15 | 05/12/15 19:53 | 150508L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 67 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 54 | 50-135 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 270 Peak | 15-05-0402-14-I | 05/06/15 00:06 | Aqueous | GC 31 | 05/08/15 | 05/12/15 20:12 | 150508L15 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 90 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 76 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-05-0402-16-I | 05/06/15 01:30 | Aqueous | GC 31 | 05/08/15 | 05/12/15 20:31 | 150508L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 96 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 77 | 50-135 | |

| HSM 230 Trail | 15-05-0402-17-I | 05/06/15 02:15 | Aqueous | GC 31 | 05/08/15 | 05/12/15 20:50 | 150508L15 |
|---------------|-----------------|----------------|---------|-------|----------|----------------|-----------|
|---------------|-----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 98 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 82 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Trail | 15-05-0402-18-I | 05/06/15 02:40 | Aqueous | GC 31 | 05/08/15 | 05/12/15 21:09 | 150508L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 91 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 72 | 50-135 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Trail | 15-05-0402-19-I | 05/06/15 03:05 | Aqueous | GC 31 | 05/08/15 | 05/12/15 21:28 | 150508L15 |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 84 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 64 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-05-0402-20-I | 05/06/15 03:58 | Aqueous | GC 31 | 05/08/15 | 05/12/15 21:47 | 150508L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 80 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 63 | 50-135 | |

| HSM 260 Trail | 15-05-0402-21-I | 05/06/15 04:20 | Aqueous | GC 31 | 05/08/15 | 05/12/15 22:06 | 150508L15 |
|---------------|-----------------|----------------|---------|-------|----------|----------------|-----------|
|---------------|-----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 106 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 89 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Trail | 15-05-0402-22-I | 05/06/15 04:40 | Aqueous | GC 31 | 05/09/15 | 05/12/15 22:25 | 150509L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 102 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 88 | 50-135 | |

| FDHSM 210 Trail | 15-05-0402-23-I | 05/06/15 01:30 | Aqueous | GC 31 | 05/09/15 | 05/12/15 22:44 | 150509L12 |
|-----------------|-----------------|-------------------|---------|-------|----------|-------------------|-----------|
|-----------------|-----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 79 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 69 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 230 Trail | 15-05-0402-24-I | 05/06/15 02:15 | Aqueous | GC 31 | 05/09/15 | 05/12/15 23:03 | 150509L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 107 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 95 | 50-135 | |

| | | | | | | | |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 231 Trail | 15-05-0402-25-I | 05/06/15 02:40 | Aqueous | GC 31 | 05/09/15 | 05/12/15 23:22 | 150509L12 |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 110 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 91 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

Page 13 of 13

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-533-1035 | N/A | Aqueous | GC 31 | 05/08/15 | 05/11/15 18:38 | 150508L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 105 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 83 | 50-135 | |

| Method Blank | 099-12-533-1034 | N/A | Aqueous | GC 31 | 05/09/15 | 05/11/15 17:21 | 150509L12 |
|--------------|-----------------|-----|---------|-------|----------|----------------|-----------|
|--------------|-----------------|-----|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 105 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 98 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 210 Lead | 15-05-0402-1-I | 05/05/15 22:10 | Aqueous | GC 26 | 05/09/15 | 05/12/15 13:25 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 130 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Lead | 15-05-0402-2-I | 05/05/15 22:20 | Aqueous | GC 26 | 05/09/15 | 05/12/15 14:10 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 140 | 30-130 | 2,7 |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Lead | 15-05-0402-3-I | 05/05/15 22:11 | Aqueous | GC 26 | 05/09/15 | 05/12/15 14:54 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 122 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Lead | 15-05-0402-4-I | 05/05/15 22:49 | Aqueous | GC 26 | 05/09/15 | 05/12/15 15:38 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 129 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 250 Lead | 15-05-0402-5-I | 05/05/15 22:27 | Aqueous | GC 26 | 05/09/15 | 05/12/15 16:22 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 130 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

Page 6 of 26

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 260 Lead | 15-05-0402-6-I | 05/05/15 22:46 | Aqueous | GC 26 | 05/09/15 | 05/12/15 17:50 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 114 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 270 Lead | 15-05-0402-7-I | 05/05/15 23:06 | Aqueous | GC 26 | 05/09/15 | 05/12/15 18:34 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 121 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 210 Peak | 15-05-0402-8-I | 05/05/15 23:24 | Aqueous | GC 26 | 05/09/15 | 05/12/15 19:19 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 113 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 230 Peak | 15-05-0402-9-I | 05/05/15 23:35 | Aqueous | GC 26 | 05/09/15 | 05/12/15 20:03 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 123 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Peak | 15-05-0402-10-I | 05/05/15 22:11 | Aqueous | GC 26 | 05/09/15 | 05/12/15 20:47 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 117 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Peak | 15-05-0402-11-I | 05/06/15 00:05 | Aqueous | GC 26 | 05/09/15 | 05/12/15 21:31 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 114 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 250 Peak | 15-05-0402-12-I | 05/05/15 23:30 | Aqueous | GC 26 | 05/09/15 | 05/12/15 22:15 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 114 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 260 Peak | 15-05-0402-13-I | 05/05/15 23:49 | Aqueous | GC 26 | 05/09/15 | 05/12/15 22:59 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 101 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 270 Peak | 15-05-0402-14-I | 05/06/15 00:06 | Aqueous | GC 26 | 05/09/15 | 05/13/15 00:27 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 91 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-05-0402-16-I | 05/06/15 01:30 | Aqueous | GC 26 | 05/09/15 | 05/13/15 01:11 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 92 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Trail | 15-05-0402-17-I | 05/06/15 02:15 | Aqueous | GC 26 | 05/09/15 | 05/13/15 01:55 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 98 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 231 Trail | 15-05-0402-18-I | 05/06/15 02:40 | Aqueous | GC 26 | 05/09/15 | 05/13/15 02:40 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 96 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-05-0402-19-I | 05/06/15 03:05 | Aqueous | GC 26 | 05/09/15 | 05/13/15 03:24 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 96 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-05-0402-20-I | 05/06/15 03:58 | Aqueous | GC 26 | 05/09/15 | 05/13/15 04:08 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 98 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Trail | 15-05-0402-21-I | 05/06/15 04:20 | Aqueous | GC 26 | 05/09/15 | 05/13/15 04:52 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 100 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Trail | 15-05-0402-22-I | 05/06/15 04:40 | Aqueous | GC 26 | 05/09/15 | 05/13/15 07:48 | 150509L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 91 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 210 Trail | 15-05-0402-23-I | 05/06/15 01:30 | Aqueous | GC 26 | 05/09/15 | 05/13/15 08:33 | 150509L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 91 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 230 Trail | 15-05-0402-24-I | 05/06/15 02:15 | Aqueous | GC 26 | 05/09/15 | 05/13/15 09:17 | 150509L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 83 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| FDHSM 231 Trail | 15-05-0402-25-I | 05/06/15 02:40 | Aqueous | GC 26 | 05/09/15 | 05/13/15 09:58 | 150509L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 90 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-93 | N/A | Aqueous | GC 26 | 05/09/15 | 05/12/15 11:13 | 150509L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 106 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-94 | N/A | Aqueous | GC 26 | 05/09/15 | 05/13/15 05:36 | 150509L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 104 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 210 Lead | 15-05-0402-1-I | 05/05/15 22:10 | Aqueous | GC 40 | 05/08/15 | 05/15/15 05:45 | 150508L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 67 | 0-123 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Lead | 15-05-0402-2-I | 05/05/15 22:20 | Aqueous | GC 40 | 05/08/15 | 05/15/15 06:08 | 150508L12 |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 93 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Lead | 15-05-0402-3-I | 05/05/15 22:11 | Aqueous | GC 40 | 05/08/15 | 05/15/15 06:31 | 150508L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 83 | 0-123 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Lead | 15-05-0402-4-I | 05/05/15 22:49 | Aqueous | GC 40 | 05/08/15 | 05/15/15 06:55 | 150508L12 |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 79 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 250 Lead | 15-05-0402-5-I | 05/05/15 22:27 | Aqueous | GC 40 | 05/08/15 | 05/15/15 07:18 | 150508L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 77 | 0-123 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Lead | 15-05-0402-6-I | 05/05/15 22:46 | Aqueous | GC 40 | 05/08/15 | 05/15/15 07:41 | 150508L12 |
|---------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 86 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 270 Lead | 15-05-0402-7-I | 05/05/15 23:06 | Aqueous | GC 40 | 05/08/15 | 05/15/15 08:04 | 150508L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 86 | 0-123 | |

| | | | | | | | |
|---------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 210 Peak | 15-05-0402-8-I | 05/05/15 23:24 | Aqueous | GC 40 | 05/08/15 | 05/15/15 08:27 | 150508L12 |
|---------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 77 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Peak | 15-05-0402-9-I | 05/05/15 23:35 | Aqueous | GC 40 | 05/08/15 | 05/15/15 08:51 | 150508L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 80 | 0-123 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Peak | 15-05-0402-10-I | 05/05/15 22:11 | Aqueous | GC 40 | 05/08/15 | 05/15/15 09:14 | 150508L12 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 95 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Peak | 15-05-0402-11-I | 05/06/15 00:05 | Aqueous | GC 40 | 05/08/15 | 05/15/15 09:37 | 150508L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 75 | 0-123 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 250 Peak | 15-05-0402-12-I | 05/05/15 23:30 | Aqueous | GC 40 | 05/08/15 | 05/15/15 10:00 | 150508L12 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 52 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Peak | 15-05-0402-13-I | 05/05/15 23:49 | Aqueous | GC 40 | 05/08/15 | 05/15/15 10:47 | 150508L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 82 | 0-123 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 270 Peak | 15-05-0402-14-I | 05/06/15 00:06 | Aqueous | GC 40 | 05/08/15 | 05/15/15 11:10 | 150508L12 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 87 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

Page 8 of 13

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-05-0402-16-I | 05/06/15 01:30 | Aqueous | GC 40 | 05/08/15 | 05/15/15 11:33 | 150508L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 82 | 0-123 | |

| HSM 230 Trail | 15-05-0402-17-I | 05/06/15 02:15 | Aqueous | GC 40 | 05/08/15 | 05/15/15 11:56 | 150508L12 |
|---------------|-----------------|-------------------|---------|-------|----------|-------------------|-----------|
|---------------|-----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 76 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Trail | 15-05-0402-18-I | 05/06/15 02:40 | Aqueous | GC 40 | 05/08/15 | 05/15/15 12:19 | 150508L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 89 | 0-123 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Trail | 15-05-0402-19-I | 05/06/15 03:05 | Aqueous | GC 40 | 05/08/15 | 05/15/15 12:42 | 150508L12 |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 83 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 250 Trail | 15-05-0402-20-I | 05/06/15 03:58 | Aqueous | GC 40 | 05/08/15 | 05/15/15 13:06 | 150508L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 82 | 0-123 | |

| | | | | | | | |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 260 Trail | 15-05-0402-21-I | 05/06/15 04:20 | Aqueous | GC 40 | 05/08/15 | 05/15/15 13:29 | 150508L12 |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 88 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Trail | 15-05-0402-22-I | 05/06/15 04:40 | Aqueous | GC 40 | 05/08/15 | 05/15/15 13:52 | 150508L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 84 | 0-123 | |

| | | | | | | | |
|-----------------|-----------------|-------------------|---------|-------|----------|-------------------|-----------|
| FDHSM 210 Trail | 15-05-0402-23-I | 05/06/15 01:30 | Aqueous | GC 40 | 05/08/15 | 05/15/15 14:15 | 150508L13 |
|-----------------|-----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 85 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 230 Trail | 15-05-0402-24-I | 05/06/15 02:15 | Aqueous | GC 40 | 05/08/15 | 05/15/15 14:38 | 150508L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 71 | 0-123 | |

| | | | | | | | |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 231 Trail | 15-05-0402-25-I | 05/06/15 02:40 | Aqueous | GC 40 | 05/08/15 | 05/15/15 15:01 | 150508L13 |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 85 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-034-644 | N/A | Aqueous | GC 40 | 05/08/15 | 05/15/15 01:30 | 150508L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 101 | 0-123 | |

| | | | | | | | |
|--------------|----------------|-----|---------|-------|----------|----------------|-----------|
| Method Blank | 095-01-034-645 | N/A | Aqueous | GC 40 | 05/08/15 | 05/15/15 03:03 | 150508L13 |
|--------------|----------------|-----|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 100 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 1 of 75

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|------------------|-----------------|-----------------------|------------------|
| HSM 210 Lead | 15-05-0402-1-H | 05/05/15 22:10 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 14:06 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 88 | 33-120 | |
| 2-Fluorophenol | 81 | 24-120 | |
| Nitrobenzene-d5 | 78 | 38-120 | |
| p-Terphenyl-d14 | 87 | 41-137 | |
| Phenol-d6 | 82 | 16-120 | |
| 2,4,6-Tribromophenol | 78 | 27-159 | |



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|------------------|-----------------|-----------------------|------------------|
| HSM 230 Lead | 15-05-0402-2-H | 05/05/15 22:20 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 14:40 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 11 | 3.0 | 1.00 | |
| Acenaphthylene | ND | 11 | 3.1 | 1.00 | |
| Aniline | ND | 11 | 1.6 | 1.00 | |
| Anthracene | ND | 11 | 3.2 | 1.00 | |
| Azobenzene | ND | 11 | 2.8 | 1.00 | |
| Benzidine | ND | 53 | 6.9 | 1.00 | |
| Benzo (a) Anthracene | ND | 11 | 5.0 | 1.00 | |
| Benzo (a) Pyrene | ND | 11 | 2.6 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 11 | 2.4 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 11 | 2.7 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 11 | 3.4 | 1.00 | |
| Benzoic Acid | ND | 53 | 13 | 1.00 | |
| Benzyl Alcohol | ND | 11 | 2.3 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 11 | 2.7 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 27 | 2.6 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 11 | 3.4 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 11 | 3.4 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 11 | 2.9 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 11 | 2.6 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 11 | 2.5 | 1.00 | |
| 4-Chloroaniline | ND | 11 | 2.1 | 1.00 | |
| 2-Chloronaphthalene | ND | 11 | 2.9 | 1.00 | |
| 2-Chlorophenol | ND | 11 | 2.5 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 11 | 2.8 | 1.00 | |
| Chrysene | ND | 11 | 3.0 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 11 | 3.1 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 11 | 2.6 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 11 | 2.7 | 1.00 | |
| Dibenzofuran | ND | 11 | 3.0 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 11 | 3.2 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 11 | 3.3 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 11 | 3.1 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 27 | 2.8 | 1.00 | |
| 2,4-Dichlorophenol | ND | 11 | 2.6 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 11 | 2.9 | 1.00 | |
| Dimethyl Phthalate | ND | 11 | 2.8 | 1.00 | |
| 2,4-Dimethylphenol | ND | 11 | 2.6 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 53 | 15 | 1.00 | |
| 2,4-Dinitrophenol | ND | 53 | 14 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 11 | 2.5 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 11 | 2.5 | 1.00 | |
| Fluoranthene | ND | 11 | 3.3 | 1.00 | |
| Fluorene | ND | 11 | 2.9 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 11 | 3.1 | 1.00 | |
| Hexachlorobenzene | ND | 11 | 3.3 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 27 | 7.4 | 1.00 | |
| Hexachloroethane | ND | 11 | 3.2 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 11 | 2.3 | 1.00 | |
| Isophorone | ND | 11 | 2.7 | 1.00 | |
| 2-Methylnaphthalene | ND | 11 | 3.0 | 1.00 | |
| 1-Methylnaphthalene | ND | 11 | 3.0 | 1.00 | |
| 2-Methylphenol | ND | 11 | 2.2 | 1.00 | |
| 3/4-Methylphenol | ND | 11 | 2.3 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 11 | 2.5 | 1.00 | |
| N-Nitrosodimethylamine | ND | 11 | 3.4 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 11 | 2.9 | 1.00 | |
| Naphthalene | ND | 11 | 3.0 | 1.00 | |
| 4-Nitroaniline | ND | 11 | 2.3 | 1.00 | |
| 3-Nitroaniline | ND | 11 | 2.4 | 1.00 | |
| 2-Nitroaniline | ND | 11 | 2.4 | 1.00 | |
| Nitrobenzene | ND | 27 | 3.2 | 1.00 | |
| 4-Nitrophenol | ND | 11 | 1.7 | 1.00 | |
| 2-Nitrophenol | ND | 11 | 2.8 | 1.00 | |
| Pentachlorophenol | ND | 11 | 4.9 | 1.00 | |
| Phenanthrene | ND | 11 | 3.1 | 1.00 | |
| Phenol | ND | 11 | 2.2 | 1.00 | |
| Pyrene | ND | 11 | 3.2 | 1.00 | |
| Pyridine | ND | 11 | 3.2 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 11 | 3.0 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 11 | 2.7 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 11 | 2.7 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 6 of 75

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 84 | 33-120 | |
| 2-Fluorophenol | 79 | 24-120 | |
| Nitrobenzene-d5 | 75 | 38-120 | |
| p-Terphenyl-d14 | 87 | 41-137 | |
| Phenol-d6 | 83 | 16-120 | |
| 2,4,6-Tribromophenol | 82 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|------------------|-----------------|-----------------------|------------------|
| HSM 231 Lead | 15-05-0402-3-H | 05/05/15 22:11 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 14:58 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 9 of 75

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 91 | 33-120 | |
| 2-Fluorophenol | 84 | 24-120 | |
| Nitrobenzene-d5 | 82 | 38-120 | |
| p-Terphenyl-d14 | 88 | 41-137 | |
| Phenol-d6 | 87 | 16-120 | |
| 2,4,6-Tribromophenol | 82 | 27-159 | |



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|------------------|-----------------|-----------------------|------------------|
| HSM 240 Lead | 15-05-0402-4-H | 05/05/15 22:49 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 15:16 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 11 of 75

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 98 | 33-120 | |
| 2-Fluorophenol | 91 | 24-120 | |
| Nitrobenzene-d5 | 88 | 38-120 | |
| p-Terphenyl-d14 | 95 | 41-137 | |
| Phenol-d6 | 96 | 16-120 | |
| 2,4,6-Tribromophenol | 91 | 27-159 | |



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|------------------|-----------------|---------------------------|------------------|
| HSM 250 Lead | 15-05-0402-5-H | 05/05/15 22:27 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 15:34 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 95 | 33-120 | |
| 2-Fluorophenol | 87 | 24-120 | |
| Nitrobenzene-d5 | 85 | 38-120 | |
| p-Terphenyl-d14 | 93 | 41-137 | |
| Phenol-d6 | 89 | 16-120 | |
| 2,4,6-Tribromophenol | 87 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|------------------|-----------------|---------------------------|------------------|
| HSM 260 Lead | 15-05-0402-6-H | 05/05/15 22:46 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 15:52 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 86 | 33-120 | |
| 2-Fluorophenol | 78 | 24-120 | |
| Nitrobenzene-d5 | 76 | 38-120 | |
| p-Terphenyl-d14 | 82 | 41-137 | |
| Phenol-d6 | 82 | 16-120 | |
| 2,4,6-Tribromophenol | 76 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|------------------|-----------------|---------------------------|------------------|
| HSM 270 Lead | 15-05-0402-7-H | 05/05/15 23:06 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 16:10 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 92 | 33-120 | |
| 2-Fluorophenol | 84 | 24-120 | |
| Nitrobenzene-d5 | 81 | 38-120 | |
| p-Terphenyl-d14 | 85 | 41-137 | |
| Phenol-d6 | 89 | 16-120 | |
| 2,4,6-Tribromophenol | 79 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|------------------|-----------------|---------------------------|------------------|
| HSM 210 Peak | 15-05-0402-8-H | 05/05/15 23:24 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 16:28 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 81 | 33-120 | |
| 2-Fluorophenol | 74 | 24-120 | |
| Nitrobenzene-d5 | 71 | 38-120 | |
| p-Terphenyl-d14 | 77 | 41-137 | |
| Phenol-d6 | 78 | 16-120 | |
| 2,4,6-Tribromophenol | 68 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|---------------------------|----------------|------------------|-----------------|---------------------------|------------------|
| HSM 230 Peak | 15-05-0402-9-H | 05/05/15 23:35 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 17:22 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.7 | 1.00 | |
| Benzidine | ND | 51 | 6.6 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.6 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.3 | 1.00 | |
| Benzoic Acid | ND | 51 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.6 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.3 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.8 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.9 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.6 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.9 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.5 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 51 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 51 | 14 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 7.0 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.2 | 1.00 | |
| Isophorone | ND | 10 | 2.6 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.9 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.1 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.7 | 1.00 | |
| Phenanthrene | ND | 10 | 3.0 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 79 | 33-120 | |
| 2-Fluorophenol | 74 | 24-120 | |
| Nitrobenzene-d5 | 70 | 38-120 | |
| p-Terphenyl-d14 | 77 | 41-137 | |
| Phenol-d6 | 77 | 16-120 | |
| 2,4,6-Tribromophenol | 70 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|------------------|-----------------|-----------------------|------------------|
| HSM 231 Peak | 15-05-0402-10-H | 05/05/15 22:11 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 16:46 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.8 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 9.8 | 2.8 | 1.00 | |
| Aniline | ND | 9.8 | 1.5 | 1.00 | |
| Anthracene | ND | 9.8 | 3.0 | 1.00 | |
| Azobenzene | ND | 9.8 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.4 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.8 | 4.6 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.8 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.8 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.8 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.8 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.8 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.8 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.8 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.8 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.8 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.8 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.8 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.8 | 2.6 | 1.00 | |
| Chrysene | ND | 9.8 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.8 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.8 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.8 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.8 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.8 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.8 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.8 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.8 | 3.0 | 1.00 | |
| Fluorene | ND | 9.8 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.8 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.8 | 1.00 | |
| Hexachloroethane | ND | 9.8 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.8 | 2.1 | 1.00 | |
| Isophorone | ND | 9.8 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.8 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.8 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.8 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.8 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 9.8 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 9.8 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.8 | 4.6 | 1.00 | |
| Phenanthrene | ND | 9.8 | 2.9 | 1.00 | |
| Phenol | ND | 9.8 | 2.0 | 1.00 | |
| Pyrene | ND | 9.8 | 2.9 | 1.00 | |
| Pyridine | ND | 9.8 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 87 | 33-120 | |
| 2-Fluorophenol | 80 | 24-120 | |
| Nitrobenzene-d5 | 77 | 38-120 | |
| p-Terphenyl-d14 | 84 | 41-137 | |
| Phenol-d6 | 82 | 16-120 | |
| 2,4,6-Tribromophenol | 75 | 27-159 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|------------------|-----------------|-----------------------|------------------|
| HSM 240 Peak | 15-05-0402-11-H | 05/06/15 00:05 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 17:04 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 87 | 33-120 | |
| 2-Fluorophenol | 81 | 24-120 | |
| Nitrobenzene-d5 | 78 | 38-120 | |
| p-Terphenyl-d14 | 83 | 41-137 | |
| Phenol-d6 | 83 | 16-120 | |
| 2,4,6-Tribromophenol | 75 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|------------------|-----------------|---------------------------|------------------|
| HSM 250 Peak | 15-05-0402-12-H | 05/05/15 23:30 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 18:01 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 84 | 33-120 | |
| 2-Fluorophenol | 76 | 24-120 | |
| Nitrobenzene-d5 | 75 | 38-120 | |
| p-Terphenyl-d14 | 80 | 41-137 | |
| Phenol-d6 | 80 | 16-120 | |
| 2,4,6-Tribromophenol | 72 | 27-159 | |



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Peak | 15-05-0402-13-H | 05/05/15 23:49 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 18:19 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 39 of 75

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 82 | 33-120 | |
| 2-Fluorophenol | 59 | 24-120 | |
| Nitrobenzene-d5 | 80 | 38-120 | |
| p-Terphenyl-d14 | 87 | 41-137 | |
| Phenol-d6 | 40 | 16-120 | |
| 2,4,6-Tribromophenol | 77 | 27-159 | |



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|------------------|-----------------|-----------------------|------------------|
| HSM 270 Peak | 15-05-0402-14-H | 05/06/15 00:06 | Aqueous | GC/MS CCC | 05/09/15 | 05/13/15 13:28 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 73 | 33-120 | |
| 2-Fluorophenol | 54 | 24-120 | |
| Nitrobenzene-d5 | 76 | 38-120 | |
| p-Terphenyl-d14 | 81 | 41-137 | |
| Phenol-d6 | 36 | 16-120 | |
| 2,4,6-Tribromophenol | 78 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-05-0402-16-H | 05/06/15 01:30 | Aqueous | GC/MS CCC | 05/09/15 | 05/13/15 13:46 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 76 | 33-120 | |
| 2-Fluorophenol | 57 | 24-120 | |
| Nitrobenzene-d5 | 79 | 38-120 | |
| p-Terphenyl-d14 | 83 | 41-137 | |
| Phenol-d6 | 37 | 16-120 | |
| 2,4,6-Tribromophenol | 76 | 27-159 | |



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Trail | 15-05-0402-17-H | 05/06/15 02:15 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 18:37 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 79 | 33-120 | |
| 2-Fluorophenol | 59 | 24-120 | |
| Nitrobenzene-d5 | 77 | 38-120 | |
| p-Terphenyl-d14 | 84 | 41-137 | |
| Phenol-d6 | 41 | 16-120 | |
| 2,4,6-Tribromophenol | 73 | 27-159 | |



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|------------------|-----------------|-----------------------|------------------|
| HSM 231 Trail | 15-05-0402-18-H | 05/06/15 02:40 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 17:39 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 87 | 33-120 | |
| 2-Fluorophenol | 63 | 24-120 | |
| Nitrobenzene-d5 | 85 | 38-120 | |
| p-Terphenyl-d14 | 92 | 41-137 | |
| Phenol-d6 | 44 | 16-120 | |
| 2,4,6-Tribromophenol | 84 | 27-159 | |



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 52 of 75

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-05-0402-19-H | 05/06/15 03:05 | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 18:55 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 81 | 33-120 | |
| 2-Fluorophenol | 59 | 24-120 | |
| Nitrobenzene-d5 | 80 | 38-120 | |
| p-Terphenyl-d14 | 85 | 41-137 | |
| Phenol-d6 | 40 | 16-120 | |
| 2,4,6-Tribromophenol | 78 | 27-159 | |



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-05-0402-20-H | 05/06/15 03:58 | Aqueous | GC/MS CCC | 05/09/15 | 05/13/15 14:04 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 78 | 33-120 | |
| 2-Fluorophenol | 49 | 24-120 | |
| Nitrobenzene-d5 | 78 | 38-120 | |
| p-Terphenyl-d14 | 81 | 41-137 | |
| Phenol-d6 | 30 | 16-120 | |
| 2,4,6-Tribromophenol | 75 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Trail | 15-05-0402-21-H | 05/06/15 04:20 | Aqueous | GC/MS CCC | 05/09/15 | 05/13/15 14:22 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 89 | 33-120 | |
| 2-Fluorophenol | 62 | 24-120 | |
| Nitrobenzene-d5 | 90 | 38-120 | |
| p-Terphenyl-d14 | 94 | 41-137 | |
| Phenol-d6 | 38 | 16-120 | |
| 2,4,6-Tribromophenol | 87 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|------------------|-----------------|---------------------------|------------------|
| HSM 270 Trail | 15-05-0402-22-H | 05/06/15 04:40 | Aqueous | GC/MS CCC | 05/09/15 | 05/13/15 14:39 | 150511L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 77 | 33-120 | |
| 2-Fluorophenol | 52 | 24-120 | |
| Nitrobenzene-d5 | 77 | 38-120 | |
| p-Terphenyl-d14 | 79 | 41-137 | |
| Phenol-d6 | 33 | 16-120 | |
| 2,4,6-Tribromophenol | 75 | 27-159 | |



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|---------------------------|----------------|------------------|-----------------|---------------------------|------------------|
| FDHSM 210 Trail | 15-05-0402-23-H | 05/06/15 01:30 | Aqueous | GC/MS CCC | 05/09/15 | 05/13/15 14:58 | 150511L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 86 | 33-120 | |
| 2-Fluorophenol | 56 | 24-120 | |
| Nitrobenzene-d5 | 83 | 38-120 | |
| p-Terphenyl-d14 | 88 | 41-137 | |
| Phenol-d6 | 35 | 16-120 | |
| 2,4,6-Tribromophenol | 81 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|------------------|-----------------|-----------------------|------------------|
| FDHSM 231 Trail | 15-05-0402-25-H | 05/06/15 02:40 | Aqueous | GC/MS CCC | 05/09/15 | 05/13/15 15:15 | 150511L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 81 | 33-120 | |
| 2-Fluorophenol | 52 | 24-120 | |
| Nitrobenzene-d5 | 80 | 38-120 | |
| p-Terphenyl-d14 | 86 | 41-137 | |
| Phenol-d6 | 32 | 16-120 | |
| 2,4,6-Tribromophenol | 77 | 27-159 | |



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-4039 | N/A | Aqueous | GC/MS CCC | 05/09/15 | 05/12/15 13:13 | 150509L06 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 89 | 33-120 | |
| 2-Fluorophenol | 80 | 24-120 | |
| Nitrobenzene-d5 | 80 | 38-120 | |
| p-Terphenyl-d14 | 86 | 41-137 | |
| Phenol-d6 | 85 | 16-120 | |
| 2,4,6-Tribromophenol | 79 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 73 of 75

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-4040 | N/A | Aqueous | GC/MS TT | 05/11/15 | 05/12/15 12:57 | 150511L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 83 | 33-120 | |
| 2-Fluorophenol | 73 | 24-120 | |
| Nitrobenzene-d5 | 71 | 38-120 | |
| p-Terphenyl-d14 | 75 | 41-137 | |
| Phenol-d6 | 71 | 16-120 | |
| 2,4,6-Tribromophenol | 87 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 1 of 60

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Lead | 15-05-0402-1-A | 05/05/15 22:10 | Aqueous | GC/MS BB | 05/09/15 | 05/09/15 14:27 | 150509L027 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 92 | 80-120 | |
| Dibromofluoromethane | 104 | 78-126 | |
| 1,2-Dichloroethane-d4 | 112 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|----------------|-----------------|-----------------------|-------------------|
| HSM 230 Lead | 15-05-0402-2-B | 05/05/15 22:20 | Aqueous | GC/MS O | 05/13/15 | 05/13/15 19:28 | 150513L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 97 | 80-120 | |
| Dibromofluoromethane | 97 | 78-126 | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|----------------|-----------------|-----------------------|-------------------|
| HSM 231 Lead | 15-05-0402-3-B | 05/05/15 22:11 | Aqueous | GC/MS O | 05/13/15 | 05/13/15 23:01 | 150513L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 99 | 78-126 | |
| 1,2-Dichloroethane-d4 | 105 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|----------------|-----------------|-----------------------|-------------------|
| HSM 240 Lead | 15-05-0402-4-B | 05/05/15 22:49 | Aqueous | GC/MS O | 05/13/15 | 05/13/15 23:31 | 150513L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 100 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Lead | 15-05-0402-5-B | 05/05/15 22:27 | Aqueous | GC/MS O | 05/13/15 | 05/14/15 00:02 | 150513L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 100 | 80-120 | |
| Dibromofluoromethane | 96 | 78-126 | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Lead | 15-05-0402-6-B | 05/05/15 22:46 | Aqueous | GC/MS O | 05/13/15 | 05/14/15 00:32 | 150513L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 97 | 78-126 | |
| 1,2-Dichloroethane-d4 | 105 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|----------------|-----------------|-----------------------|-------------------|
| HSM 270 Lead | 15-05-0402-7-B | 05/05/15 23:06 | Aqueous | GC/MS O | 05/13/15 | 05/14/15 01:03 | 150513L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 98 | 78-126 | |
| 1,2-Dichloroethane-d4 | 106 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 15 of 60

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|----------------|-----------------|-----------------------|-------------------|
| HSM 210 Peak | 15-05-0402-8-B | 05/05/15 23:24 | Aqueous | GC/MS O | 05/13/15 | 05/14/15 01:33 | 150513L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 96 | 78-126 | |
| 1,2-Dichloroethane-d4 | 106 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Peak | 15-05-0402-9-B | 05/05/15 23:35 | Aqueous | GC/MS O | 05/13/15 | 05/14/15 02:03 | 150513L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 99 | 80-120 | |
| Dibromofluoromethane | 99 | 78-126 | |
| 1,2-Dichloroethane-d4 | 107 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|----------------|-----------------|-----------------------|-------------------|
| HSM 231 Peak | 15-05-0402-10-B | 05/05/15 22:11 | Aqueous | GC/MS O | 05/13/15 | 05/14/15 02:34 | 150513L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 97 | 80-120 | |
| Dibromofluoromethane | 96 | 78-126 | |
| 1,2-Dichloroethane-d4 | 106 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|----------------|-----------------|-----------------------|-------------------|
| HSM 240 Peak | 15-05-0402-11-B | 05/06/15 00:05 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 01:56 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 96 | 80-120 | |
| Dibromofluoromethane | 91 | 78-126 | |
| 1,2-Dichloroethane-d4 | 100 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|----------------|-----------------|---------------------------|-------------------|
| HSM 250 Peak | 15-05-0402-12-B | 05/05/15 23:30 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 03:58 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 96 | 80-120 | |
| Dibromofluoromethane | 91 | 78-126 | |
| 1,2-Dichloroethane-d4 | 99 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Peak | 15-05-0402-13-B | 05/05/15 23:49 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 04:28 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 96 | 80-120 | |
| Dibromofluoromethane | 90 | 78-126 | |
| 1,2-Dichloroethane-d4 | 99 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|----------------|-----------------|-----------------------|-------------------|
| HSM 270 Peak | 15-05-0402-14-B | 05/06/15 00:06 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 04:58 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 96 | 80-120 | |
| Dibromofluoromethane | 91 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| TB06 | 15-05-0402-15-A | 05/06/15 10:11 | Aqueous | GC/MS RR | 05/18/15 | 05/18/15 19:28 | 150518L017 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 104 | 78-126 | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-05-0402-16-B | 05/06/15 01:30 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 05:29 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | 0.18 | 1.0 | 0.16 | 1.00 | J |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 96 | 80-120 | |
| Dibromofluoromethane | 90 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 33 of 60

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Trail | 15-05-0402-17-B | 05/06/15 02:15 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 05:59 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 91 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 35 of 60

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 231 Trail | 15-05-0402-18-B | 05/06/15 02:40 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 06:29 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 36 of 60

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 96 | 80-120 | |
| Dibromofluoromethane | 88 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 37 of 60

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-05-0402-19-B | 05/06/15 03:05 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 06:59 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 95 | 80-120 | |
| Dibromofluoromethane | 90 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-05-0402-20-B | 05/06/15 03:58 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 07:29 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 89 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Trail | 15-05-0402-21-B | 05/06/15 04:20 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 08:00 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 95 | 80-120 | |
| Dibromofluoromethane | 91 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Trail | 15-05-0402-22-A | 05/06/15 04:40 | Aqueous | GC/MS O | 05/12/15 | 05/12/15 22:51 | 150512L041 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 100 | 80-120 | |
| Dibromofluoromethane | 101 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 210 Trail | 15-05-0402-23-A | 05/06/15 01:30 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 08:30 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 90 | 78-126 | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 230 Trail | 15-05-0402-24-A | 05/06/15 02:15 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 09:00 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 94 | 78-126 | |
| 1,2-Dichloroethane-d4 | 106 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 49 of 60

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 231 Trail | 15-05-0402-25-A | 05/06/15 02:40 | Aqueous | GC/MS O | 05/15/15 | 05/16/15 09:30 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 50 of 60

| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 95 | 80-120 | |
| Dibromofluoromethane | 91 | 78-126 | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 51 of 60

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-17134 | N/A | Aqueous | GC/MS BB | 05/09/15 | 05/09/15 12:13 | 150509L027 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 103 | 78-126 | |
| 1,2-Dichloroethane-d4 | 110 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-17133 | N/A | Aqueous | GC/MS O | 05/12/15 | 05/12/15 22:22 | 150512L041 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 100 | 80-120 | |
| Dibromofluoromethane | 99 | 78-126 | |
| 1,2-Dichloroethane-d4 | 99 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-17137 | N/A | Aqueous | GC/MS O | 05/13/15 | 05/13/15 16:47 | 150513L015 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 95 | 78-126 | |
| 1,2-Dichloroethane-d4 | 99 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 57 of 60

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-17160 | N/A | Aqueous | GC/MS O | 05/15/15 | 05/16/15 01:26 | 150515L039 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 99 | 80-120 | |
| Dibromofluoromethane | 95 | 78-126 | |
| 1,2-Dichloroethane-d4 | 100 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-17170 | N/A | Aqueous | GC/MS RR | 05/18/15 | 05/18/15 14:01 | 150518L017 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 97 | 80-120 | |
| Dibromofluoromethane | 104 | 78-126 | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

05/07/15

Work Order:

15-05-0402

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HSM 210 Lead | 15-05-0402-1 | 05/05/15 22:10 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 257 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 257 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 595 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 3.3 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.48 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 2.7 | 0.50 | 0.24 | 1.00 | | mg/L | 05/29/15 | 05/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 4.6 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

| HSM 230 Lead | 15-05-0402-2 | 05/05/15 22:20 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.031 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 199 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 199 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 215 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 45 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.46 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.1 | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 6.6 | 0.50 | 0.24 | 1.00 | | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 8.8 | 2.5 | 1.2 | 5.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

05/07/15

Work Order:

15-05-0402

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HSM 231 Lead | 15-05-0402-3 | 05/05/15 22:11 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.038 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 253 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 253 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 350 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 8.2 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.48 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.40 | 0.50 | 0.28 | 1.00 | J | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.4 | 0.50 | 0.24 | 1.00 | | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.2 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

| | | | |
|---------------------|---------------------|-----------------------|----------------|
| HSM 240 Lead | 15-05-0402-4 | 05/05/15 22:49 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.11 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 174 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 174 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 210 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 377 | 1.00 | 0.829 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.59 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.1 | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.8 | 0.50 | 0.24 | 1.00 | | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 7.2 | 2.5 | 1.2 | 5.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

05/07/15

Work Order:

15-05-0402

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HSM 250 Lead | 15-05-0402-5 | 05/05/15 22:27 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.035 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 238 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 238 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 320 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 8.9 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.46 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 1.0 | 0.50 | 0.24 | 1.00 | | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.0 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

| | | | |
|---------------------|---------------------|-----------------------|----------------|
| HSM 260 Lead | 15-05-0402-6 | 05/05/15 22:46 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.026 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 239 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 239 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 375 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 12 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.63 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 1.6 | 0.50 | 0.24 | 1.00 | | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.0 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

05/07/15

Work Order:

15-05-0402

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HSM 270 Lead | 15-05-0402-7 | 05/05/15 23:06 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.033 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 238 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 238 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 385 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 20 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.75 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 2.1 | 0.50 | 0.24 | 1.00 | | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 1.9 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

| | | | |
|---------------------|---------------------|-----------------------|----------------|
| HSM 210 Peak | 15-05-0402-8 | 05/05/15 23:24 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 242 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 242 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 360 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.4 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.58 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.84 | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 2.6 | 0.50 | 0.24 | 1.00 | | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 3.5 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HSM 230 Peak | 15-05-0402-9 | 05/05/15 23:35 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.062 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 190 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 190 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 330 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 13 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.44 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.70 | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 4.1 | 0.50 | 0.24 | 1.00 | | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 5.4 | 2.5 | 1.2 | 5.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

| | | | |
|---------------------|----------------------|-----------------------|----------------|
| HSM 231 Peak | 15-05-0402-10 | 05/05/15 22:11 | Aqueous |
|---------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 254 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 254 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 370 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.3 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.55 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.84 | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 0.30 | 0.50 | 0.24 | 1.00 | J | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.2 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

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San Antonio, TX 78249-1618

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|----------------------|-----------------------|----------------|
| HSM 240 Peak | 15-05-0402-11 | 05/06/15 00:05 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 249 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 249 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Solids, Total Dissolved (24) | 365 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.56 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.7 | 0.50 | 0.24 | 1.00 | | mg/L | 05/29/15 | 05/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 3.7 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

| HSM 250 Peak | 15-05-0402-12 | 05/05/15 23:30 | Aqueous |
|---------------------|----------------------|-----------------------|----------------|
|---------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.028 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 236 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 236 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Solids, Total Dissolved (24) | 360 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 4.3 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.54 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 4.8 | 0.50 | 0.24 | 1.00 | | mg/L | 05/29/15 | 05/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 3.9 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

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Analytical Report

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|----------------------|-----------------------|----------------|
| HSM 260 Peak | 15-05-0402-13 | 05/05/15 23:49 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.032 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 227 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 227 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Solids, Total Dissolved (24) | 350 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 13 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.68 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.70 | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.0 | 0.50 | 0.24 | 1.00 | | mg/L | 05/29/15 | 05/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.6 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

| | | | |
|---------------------|----------------------|-----------------------|----------------|
| HSM 270 Peak | 15-05-0402-14 | 05/06/15 00:06 | Aqueous |
|---------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.042 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 225 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 225 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Solids, Total Dissolved (24) | 385 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 12 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.56 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 0.70 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|----------------------|-----------------------|----------------|
| HSM 210 Trail | 15-05-0402-16 | 05/06/15 01:30 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.054 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 240 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 240 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Solids, Total Dissolved (24) | 405 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 2.3 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.48 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 0.46 | 0.50 | 0.24 | 1.00 | J | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.4 | 2.5 | 1.2 | 5.00 | J | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

| HSM 230 Trail | 15-05-0402-17 | 05/06/15 02:15 | Aqueous |
|----------------------|----------------------|-----------------------|----------------|
|----------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.045 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 199 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 199 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 325 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 2.4 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.35 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 3.2 | 0.50 | 0.24 | 1.00 | | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.6 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

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|----------------------|----------------------|-----------------------|----------------|
| HSM 231 Trail | 15-05-0402-18 | 05/06/15 02:40 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 252 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 252 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 385 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.2 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.48 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.84 | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 4.0 | 0.50 | 0.24 | 1.00 | | mg/L | 05/29/15 | 05/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 5.0 | 2.5 | 1.2 | 5.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

| | | | |
|----------------------|----------------------|-----------------------|----------------|
| HSM 240 Trail | 15-05-0402-19 | 05/06/15 03:05 | Aqueous |
|----------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 251 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 251 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 360 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.2 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.50 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 3.1 | 0.50 | 0.24 | 1.00 | | mg/L | 05/29/15 | 05/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 7.0 | 2.5 | 1.2 | 5.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

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Analytical Report

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|----------------------|----------------------|-----------------------|----------------|
| HSM 250 Trail | 15-05-0402-20 | 05/06/15 03:58 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 244 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 244 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 365 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.8 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.53 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 05/22/15 | 05/22/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 3.8 | 0.50 | 0.24 | 1.00 | | mg/L | 05/29/15 | 05/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 2.4 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

| HSM 260 Trail | 15-05-0402-21 | 05/06/15 04:20 | Aqueous |
|----------------------|----------------------|-----------------------|----------------|
|----------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 242 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 242 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 295 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 4.3 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.61 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 05/22/15 | 05/22/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 4.5 | 0.50 | 0.24 | 1.00 | | mg/L | 05/29/15 | 05/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 6.5 | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

05/07/15

Work Order:

15-05-0402

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|----------------------|-----------------------|----------------|
| HSM 270 Trail | 15-05-0402-22 | 05/06/15 04:40 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 241 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 241 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 305 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 3.6 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.70 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 05/22/15 | 05/22/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 0.30 | 0.50 | 0.24 | 1.00 | J | mg/L | 05/26/15 | 05/26/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

| FDHSM 210 Trail | 15-05-0402-23 | 05/06/15 01:30 | Aqueous |
|-----------------|---------------|----------------|---------|
|-----------------|---------------|----------------|---------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 256 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 256 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 340 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 3.8 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.47 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 05/22/15 | 05/22/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 3.0 | 0.50 | 0.24 | 1.00 | | mg/L | 05/26/15 | 05/26/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 3.0 | 2.5 | 1.2 | 5.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

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15-05-0402

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|------------------------|----------------------|-----------------------|----------------|
| FDHSM 230 Trail | 15-05-0402-24 | 05/06/15 02:15 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.029 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 198 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 198 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 200 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | 3.3 | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.48 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 05/22/15 | 05/22/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 4.0 | 0.50 | 0.24 | 1.00 | | mg/L | 05/29/15 | 05/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 6.6 | 2.5 | 1.2 | 5.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

| FDHSM 231 Trail | 15-05-0402-25 | 05/06/15 02:40 | Aqueous |
|-----------------|---------------|----------------|---------|
|-----------------|---------------|----------------|---------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.042 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 255 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 255 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Solids, Total Dissolved (24) | 255 | 1.00 | 0.870 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| pH (24) | 7.52 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 05/07/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 05/22/15 | 05/22/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 2.5 | 0.50 | 0.24 | 1.00 | | mg/L | 05/29/15 | 05/30/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 3.0 | 2.5 | 1.2 | 5.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

05/07/15

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15-05-0402

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|-------------------|---------------------|---------|
| Method Blank | | N/A | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|-------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 05/20/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/19/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 05/18/15 | SM 2320B |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.87 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.87 | 1.00 | | mg/L | 05/12/15 | 05/12/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 05/08/15 | 05/08/15 | SM 2540 D |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 05/21/15 | 05/21/15 | SM 4500 N Org B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 05/22/15 | 05/22/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 05/29/15 | 05/30/15 | SM 5310 B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 05/22/15 | 05/23/15 | SM 5310 B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 05/26/15 | 05/26/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 05/13/15 | 05/14/15 | SM 5310 B |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-05-0429-3 | Sample | Aqueous | IC 15 | N/A | 05/07/15 13:55 | 150507S01 |
| 15-05-0429-3 | Matrix Spike | Aqueous | IC 15 | N/A | 05/07/15 14:12 | 150507S01 |
| 15-05-0429-3 | Matrix Spike Duplicate | Aqueous | IC 15 | N/A | 05/07/15 14:29 | 150507S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.1893 | 250.0 | 213.5 | 85 | 214.3 | 86 | 80-120 | 0 | 0-20 | |
| Chloride | 36.27 | 5000 | 4509 | 89 | 4502 | 89 | 80-120 | 0 | 0-20 | |
| Bromide | ND | 500.0 | 452.6 | 91 | 452.4 | 90 | 80-120 | 0 | 0-20 | |
| Nitrate (as N) | ND | 500.0 | 446.1 | 89 | 445.2 | 89 | 80-120 | 0 | 0-20 | |
| Sulfate | 227.2 | 5000 | 4575 | 87 | 4558 | 87 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-05-0447-1 | Sample | Aqueous | IC 15 | N/A | 05/08/15 01:00 | 150507S02 |
| 15-05-0447-1 | Matrix Spike | Aqueous | IC 15 | N/A | 05/08/15 01:34 | 150507S02 |
| 15-05-0447-1 | Matrix Spike Duplicate | Aqueous | IC 15 | N/A | 05/08/15 01:53 | 150507S02 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.3060 | 250.0 | 237.8 | 95 | 236.4 | 94 | 80-120 | 1 | 0-20 | |
| Chloride | 138.3 | 5000 | 5122 | 100 | 5135 | 100 | 80-120 | 0 | 0-20 | |
| Bromide | 0.6960 | 500.0 | 498.4 | 100 | 499.1 | 100 | 80-120 | 0 | 0-20 | |
| Nitrate (as N) | ND | 500.0 | 493.4 | 99 | 494.4 | 99 | 80-120 | 0 | 0-20 | |
| Sulfate | 273.5 | 5000 | 5161 | 98 | 5162 | 98 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM 210 Lead | Sample | Aqueous | ACA 1 | N/A | 05/20/15 19:29 | 150520S01 |
| HSM 210 Lead | Matrix Spike | Aqueous | ACA 1 | N/A | 05/20/15 19:29 | 150520S01 |
| HSM 210 Lead | Matrix Spike Duplicate | Aqueous | ACA 1 | N/A | 05/20/15 19:29 | 150520S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | ND | 0.2000 | 0.1598 | 80 | 0.1607 | 80 | 90-110 | 1 | 0-25 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| FDHSM 210 Trail | Sample | Aqueous | ACA 1 | N/A | 05/20/15 19:29 | 150520S02 |
| FDHSM 210 Trail | Matrix Spike | Aqueous | ACA 1 | N/A | 05/20/15 19:29 | 150520S02 |
| FDHSM 210 Trail | Matrix Spike Duplicate | Aqueous | ACA 1 | N/A | 05/20/15 19:29 | 150520S02 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | ND | 0.2000 | 0.1963 | 98 | 0.1981 | 99 | 90-110 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
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Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HSM 210 Lead | Sample | Aqueous | TOC 8 | 05/22/15 | 05/23/15 06:08 | F0522TOCS1 | | | | |
| HSM 210 Lead | Matrix Spike | Aqueous | TOC 8 | 05/22/15 | 05/23/15 06:08 | F0522TOCS1 | | | | |
| HSM 210 Lead | Matrix Spike Duplicate | Aqueous | TOC 8 | 05/22/15 | 05/23/15 06:08 | F0522TOCS1 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 1.550 | 10.00 | 10.30 | 88 | 10.30 | 88 | 31-145 | 0 | 0-23 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|---------------------|-----------------|------------|---------------|-------------------|
| HSM 270 Trail | Sample | Aqueous | TOC 8 | 05/26/15 | 05/26/15 22:23 | F0526TOCS1 | | | | |
| HSM 270 Trail | Matrix Spike | Aqueous | TOC 8 | 05/26/15 | 05/26/15 22:23 | F0526TOCS1 | | | | |
| HSM 270 Trail | Matrix Spike Duplicate | Aqueous | TOC 8 | 05/26/15 | 05/26/15 22:23 | F0526TOCS1 | | | | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Carbon, Total Organic | ND | 10.00 | 9.860 | 99 | 9.620 | 96 | 31-145 | 2 | 0-23 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|---------------------|-----------------|------------|---------------|-------------------|
| HSM 260 Trail | Sample | Aqueous | TOC 8 | 05/29/15 | 05/30/15 06:47 | F0529TOCS1 | | | | |
| HSM 260 Trail | Matrix Spike | Aqueous | TOC 8 | 05/29/15 | 05/30/15 06:47 | F0529TOCS1 | | | | |
| HSM 260 Trail | Matrix Spike Duplicate | Aqueous | TOC 8 | 05/29/15 | 05/30/15 06:47 | F0529TOCS1 | | | | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Carbon, Total Organic | 4.520 | 10.00 | 12.00 | 75 | 11.70 | 72 | 31-145 | 3 | 0-23 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|---------------------|-----------------|------------|---------------|-------------------|
| HSM 210 Lead | Sample | Aqueous | TOC 8 | 05/13/15 | 05/14/15 14:53 | F0513DOCS1 | | | | |
| HSM 210 Lead | Matrix Spike | Aqueous | TOC 8 | 05/13/15 | 05/14/15 14:53 | F0513DOCS1 | | | | |
| HSM 210 Lead | Matrix Spike Duplicate | Aqueous | TOC 8 | 05/13/15 | 05/14/15 14:53 | F0513DOCS1 | | | | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Carbon, Dissolved Organic | 4.550 | 10.00 | 15.40 | 108 | 12.70 | 82 | 31-145 | 19 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HSM 270 Trail | Sample | Aqueous | TOC 8 | 05/13/15 | 05/14/15 19:49 | F0513DOCS2 | | | | |
| HSM 270 Trail | Matrix Spike | Aqueous | TOC 8 | 05/13/15 | 05/14/15 19:49 | F0513DOCS2 | | | | |
| HSM 270 Trail | Matrix Spike Duplicate | Aqueous | TOC 8 | 05/13/15 | 05/14/15 19:49 | F0513DOCS2 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Dissolved Organic | ND | 50.00 | 47.85 | 96 | 49.40 | 99 | 31-145 | 3 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | | Instrument | Date Prepared | Date Analyzed | | MS/MSD Batch Number | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|------------------|-----------------|------------|---------------------|-------------------|
| HSM 230 Lead | Sample | | Aqueous | | ICP 7300 | 05/07/15 | 05/16/15 21:18 | | 150507SA5 | |
| HSM 230 Lead | Matrix Spike | | Aqueous | | ICP 7300 | 05/07/15 | 05/16/15 21:10 | | 150507SA5 | |
| HSM 230 Lead | Matrix Spike Duplicate | | Aqueous | | ICP 7300 | 05/07/15 | 05/16/15 21:12 | | 150507SA5 | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Calcium | 40.10 | 0.5000 | 40.38 | 4X | 39.93 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 4.765 | 0.5000 | 5.412 | 4X | 5.374 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 3.242 | 5.000 | 8.754 | 110 | 8.797 | 111 | 83-131 | 0 | 0-7 | |
| Sodium | 6.322 | 5.000 | 11.32 | 100 | 11.32 | 100 | 73-127 | 0 | 0-9 | |
| Strontium | 0.1609 | 0.5000 | 0.7410 | 116 | 0.7278 | 113 | 81-123 | 2 | 0-6 | |
| Silicon | 2.013 | 0.5000 | 2.755 | 4X | 2.729 | 4X | 24-180 | 4X | 0-15 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|---------------------------|------------------------|-------------|----------|------------|---------------|----------------|---------------------|-----|--------|------------|
| FDHSM 230 Trail | Sample | | Aqueous | ICP 7300 | 05/07/15 | 05/08/15 19:05 | 150507SA6 | | | |
| FDHSM 230 Trail | Matrix Spike | | Aqueous | ICP 7300 | 05/07/15 | 05/08/15 19:07 | 150507SA6 | | | |
| FDHSM 230 Trail | Matrix Spike Duplicate | | Aqueous | ICP 7300 | 05/07/15 | 05/08/15 19:08 | 150507SA6 | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Calcium | 86.76 | 0.5000 | 95.63 | 4X | 95.81 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 12.49 | 0.5000 | 12.53 | 4X | 12.82 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 2.048 | 5.000 | 7.081 | 101 | 7.262 | 104 | 83-131 | 3 | 0-7 | |
| Sodium | 12.91 | 5.000 | 18.80 | 118 | 19.16 | 125 | 73-127 | 2 | 0-9 | |
| Strontium | 0.4114 | 0.5000 | 1.123 | 142 | 1.146 | 147 | 81-123 | 2 | 0-6 | 3 |
| Silicon | 4.108 | 0.5000 | 4.995 | 4X | 4.993 | 4X | 24-180 | 4X | 0-15 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | |
|---------------------------|------------------------|-------------|----------|----------|------------|---------------|----------------|---------------------|--------|------------|
| HSM 210 Lead | Sample | | Aqueous | | ICP/MS 04 | 05/07/15 | 05/11/15 21:21 | 150507SA4 | | |
| HSM 210 Lead | Matrix Spike | | Aqueous | | ICP/MS 04 | 05/07/15 | 05/11/15 21:01 | 150507SA4 | | |
| HSM 210 Lead | Matrix Spike Duplicate | | Aqueous | | ICP/MS 04 | 05/07/15 | 05/11/15 21:05 | 150507SA4 | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.08492 | 85 | 0.09805 | 98 | 85-133 | 14 | 0-11 | 4 |
| Arsenic | 0.001117 | 0.1000 | 0.08380 | 83 | 0.09547 | 94 | 73-127 | 13 | 0-11 | 4 |
| Barium | 0.03955 | 0.1000 | 0.1432 | 104 | 0.1439 | 104 | 74-128 | 0 | 0-10 | |
| Beryllium | ND | 0.1000 | 0.08334 | 83 | 0.09752 | 98 | 56-122 | 16 | 0-11 | 4 |
| Cadmium | ND | 0.1000 | 0.08112 | 81 | 0.09400 | 94 | 84-114 | 15 | 0-8 | 3,4 |
| Chromium | ND | 0.1000 | 0.08843 | 88 | 0.1017 | 102 | 73-133 | 14 | 0-11 | 4 |
| Copper | ND | 0.1000 | 0.08288 | 83 | 0.09533 | 95 | 72-108 | 14 | 0-10 | 4 |
| Lead | ND | 0.1000 | 0.08491 | 85 | 0.09705 | 97 | 79-121 | 13 | 0-10 | 4 |
| Nickel | 0.002058 | 0.1000 | 0.07987 | 78 | 0.09155 | 89 | 68-122 | 14 | 0-10 | 4 |
| Selenium | ND | 0.1000 | 0.07557 | 76 | 0.08794 | 88 | 59-125 | 15 | 0-12 | 4 |
| Silver | ND | 0.05000 | 0.05111 | 102 | 0.05106 | 102 | 68-128 | 0 | 0-14 | |
| Thallium | ND | 0.1000 | 0.08391 | 84 | 0.09678 | 97 | 73-121 | 14 | 0-11 | 4 |
| Zinc | 0.01170 | 0.1000 | 0.1451 | 133 | 0.09586 | 84 | 43-145 | 41 | 0-39 | 4 |
| Aluminum | ND | 0.1000 | 0.1150 | 115 | 0.1112 | 111 | 47-161 | 3 | 0-24 | |
| Iron | 0.05564 | 5.100 | 5.569 | 108 | 5.532 | 107 | 27-201 | 1 | 0-24 | |
| Manganese | 0.01498 | 0.1000 | 0.09788 | 83 | 0.1084 | 93 | 72-126 | 10 | 0-42 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|---------------------------|------------------------|-------------|----------|------------|---------------|----------------|---------------------|-----|--------|------------|
| FDHSM 210 Trail | Sample | | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 00:48 | 150507SA6 | | | |
| FDHSM 210 Trail | Matrix Spike | | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 18:07 | 150507SA6 | | | |
| FDHSM 210 Trail | Matrix Spike Duplicate | | Aqueous | ICP/MS 04 | 05/07/15 | 05/12/15 18:11 | 150507SA6 | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.08691 | 87 | 0.08237 | 82 | 85-133 | 5 | 0-11 | 3 |
| Arsenic | 0.001241 | 0.1000 | 0.07790 | 77 | 0.07358 | 72 | 73-127 | 6 | 0-11 | 3 |
| Barium | 0.03968 | 0.1000 | 0.1330 | 93 | 0.1396 | 100 | 74-128 | 5 | 0-10 | |
| Beryllium | ND | 0.1000 | 0.09143 | 91 | 0.08651 | 87 | 56-122 | 6 | 0-11 | |
| Cadmium | ND | 0.1000 | 0.08438 | 84 | 0.08011 | 80 | 84-114 | 5 | 0-8 | 3 |
| Chromium | ND | 0.1000 | 0.09848 | 98 | 0.09234 | 92 | 73-133 | 6 | 0-11 | |
| Copper | ND | 0.1000 | 0.08187 | 82 | 0.07770 | 78 | 72-108 | 5 | 0-10 | |
| Lead | ND | 0.1000 | 0.08843 | 88 | 0.08335 | 83 | 79-121 | 6 | 0-10 | |
| Nickel | 0.002068 | 0.1000 | 0.08273 | 81 | 0.07959 | 78 | 68-122 | 4 | 0-10 | |
| Selenium | ND | 0.1000 | 0.07382 | 74 | 0.06980 | 70 | 59-125 | 6 | 0-12 | |
| Silver | ND | 0.05000 | 0.04911 | 98 | 0.05238 | 105 | 68-128 | 6 | 0-14 | |
| Thallium | ND | 0.1000 | 0.08732 | 87 | 0.08291 | 83 | 73-121 | 5 | 0-11 | |
| Zinc | 0.01100 | 0.1000 | 0.1004 | 89 | 0.1033 | 92 | 43-145 | 3 | 0-39 | |
| Aluminum | ND | 0.1000 | 0.1175 | 118 | 0.1270 | 127 | 47-161 | 8 | 0-24 | |
| Iron | 0.07078 | 5.100 | 5.429 | 105 | 5.660 | 110 | 27-201 | 4 | 0-24 | |
| Manganese | 0.02035 | 0.1000 | 0.1065 | 86 | 0.1042 | 84 | 72-126 | 2 | 0-42 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM 231 Lead | Sample | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:22 | 150511S04 |
| HSM 231 Lead | Matrix Spike | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:04 | 150511S04 |
| HSM 231 Lead | Matrix Spike Duplicate | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:07 | 150511S04 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | ND | 0.01000 | 0.009775 | 98 | 0.01015 | 101 | 57-141 | 4 | 0-10 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| FDHSM 231 Trail | Sample | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 19:19 | 150511S05 | | | | |
| FDHSM 231 Trail | Matrix Spike | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:09 | 150511S05 | | | | |
| FDHSM 231 Trail | Matrix Spike Duplicate | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:11 | 150511S05 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Mercury | ND | 0.01000 | 0.009875 | 99 | 0.009661 | 97 | 57-141 | 2 | 0-10 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HSM 231 Lead | Sample | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 18:33 | 150512S21 | | | | |
| HSM 231 Lead | Matrix Spike | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 12:44 | 150512S21 | | | | |
| HSM 231 Lead | Matrix Spike Duplicate | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 12:52 | 150512S21 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Caffeine | 368.8 | 100.0 | 472.7 | 104 | 451.5 | 83 | 70-130 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HSM 260 Trail | Sample | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 15:59 | 150512S22 | | | | |
| HSM 260 Trail | Matrix Spike | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 13:00 | 150512S22 | | | | |
| HSM 260 Trail | Matrix Spike Duplicate | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 13:10 | 150512S22 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Caffeine | 109.8 | 100.0 | 220.3 | 111 | 217.0 | 107 | 70-130 | 2 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HSM 270 Trail | Sample | Aqueous | GC/MS O | 05/12/15 | 05/12/15 22:51 | 150512S028 | | | | |
| HSM 270 Trail | Matrix Spike | Aqueous | GC/MS O | 05/12/15 | 05/12/15 23:19 | 150512S028 | | | | |
| HSM 270 Trail | Matrix Spike Duplicate | Aqueous | GC/MS O | 05/12/15 | 05/12/15 23:48 | 150512S028 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 50.76 | 102 | 51.42 | 103 | 74-122 | 1 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 53.96 | 108 | 53.80 | 108 | 60-144 | 0 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 50.28 | 101 | 51.48 | 103 | 73-120 | 2 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 54.84 | 110 | 56.27 | 113 | 80-122 | 3 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 48.10 | 96 | 48.92 | 98 | 70-120 | 2 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 51.37 | 103 | 52.16 | 104 | 64-142 | 2 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 52.74 | 105 | 47.94 | 96 | 52-136 | 10 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 51.92 | 104 | 53.23 | 106 | 77-125 | 2 | 0-24 | |
| Toluene | ND | 50.00 | 50.94 | 102 | 51.50 | 103 | 72-126 | 1 | 0-23 | |
| Trichloroethene | ND | 50.00 | 51.34 | 103 | 51.73 | 103 | 74-128 | 1 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 44.53 | 89 | 45.74 | 91 | 67-133 | 3 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 98.21 | 98 | 101.9 | 102 | 63-129 | 4 | 0-25 | |
| o-Xylene | ND | 50.00 | 48.02 | 96 | 49.41 | 99 | 62-128 | 3 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 48.56 | 97 | 45.55 | 91 | 68-134 | 6 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HSM 230 Lead | Sample | Aqueous | GC/MS O | 05/13/15 | 05/13/15 19:28 | 150513S010 | | | | |
| HSM 230 Lead | Matrix Spike | Aqueous | GC/MS O | 05/13/15 | 05/13/15 20:59 | 150513S010 | | | | |
| HSM 230 Lead | Matrix Spike Duplicate | Aqueous | GC/MS O | 05/13/15 | 05/13/15 21:30 | 150513S010 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 49.21 | 98 | 47.55 | 95 | 74-122 | 3 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 46.45 | 93 | 45.18 | 90 | 60-144 | 3 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 48.30 | 97 | 46.78 | 94 | 73-120 | 3 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 49.66 | 99 | 49.74 | 99 | 80-122 | 0 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 45.24 | 90 | 45.32 | 91 | 70-120 | 0 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 50.29 | 101 | 48.37 | 97 | 64-142 | 4 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 48.70 | 97 | 48.57 | 97 | 52-136 | 0 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 49.53 | 99 | 47.73 | 95 | 77-125 | 4 | 0-24 | |
| Toluene | ND | 50.00 | 49.38 | 99 | 47.94 | 96 | 72-126 | 3 | 0-23 | |
| Trichloroethene | ND | 50.00 | 48.52 | 97 | 47.09 | 94 | 74-128 | 3 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 48.54 | 97 | 47.98 | 96 | 67-133 | 1 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 96.10 | 96 | 93.30 | 93 | 63-129 | 3 | 0-25 | |
| o-Xylene | ND | 50.00 | 46.77 | 94 | 45.20 | 90 | 62-128 | 3 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 45.92 | 92 | 43.06 | 86 | 68-134 | 6 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HSM 240 Peak | Sample | Aqueous | GC/MS O | 05/15/15 | 05/16/15 01:56 | 150515S020 | | | | |
| HSM 240 Peak | Matrix Spike | Aqueous | GC/MS O | 05/15/15 | 05/16/15 02:27 | 150515S020 | | | | |
| HSM 240 Peak | Matrix Spike Duplicate | Aqueous | GC/MS O | 05/15/15 | 05/16/15 02:57 | 150515S020 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 47.47 | 95 | 49.71 | 99 | 74-122 | 5 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 38.29 | 77 | 43.77 | 88 | 60-144 | 13 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 47.62 | 95 | 49.36 | 99 | 73-120 | 4 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 50.63 | 101 | 52.62 | 105 | 80-122 | 4 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 46.15 | 92 | 47.55 | 95 | 70-120 | 3 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 48.39 | 97 | 51.34 | 103 | 64-142 | 6 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 38.27 | 77 | 47.48 | 95 | 52-136 | 21 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 48.80 | 98 | 49.85 | 100 | 77-125 | 2 | 0-24 | |
| Toluene | ND | 50.00 | 47.48 | 95 | 49.95 | 100 | 72-126 | 5 | 0-23 | |
| Trichloroethene | ND | 50.00 | 47.61 | 95 | 50.23 | 100 | 74-128 | 5 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 37.72 | 75 | 40.64 | 81 | 67-133 | 7 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 94.43 | 94 | 97.36 | 97 | 63-129 | 3 | 0-25 | |
| o-Xylene | ND | 50.00 | 46.09 | 92 | 47.31 | 95 | 62-128 | 3 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 33.89 | 68 | 43.46 | 87 | 68-134 | 25 | 0-21 | 4 |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HSM 210 Lead | Sample | Aqueous | GC/MS BB | 05/09/15 | 05/09/15 14:27 | 150509S015 | | | | |
| HSM 210 Lead | Matrix Spike | Aqueous | GC/MS BB | 05/09/15 | 05/09/15 15:53 | 150509S015 | | | | |
| HSM 210 Lead | Matrix Spike Duplicate | Aqueous | GC/MS BB | 05/09/15 | 05/09/15 16:21 | 150509S015 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 40.01 | 80 | 40.03 | 80 | 74-122 | 0 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 39.36 | 79 | 39.30 | 79 | 60-144 | 0 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 43.11 | 86 | 43.28 | 87 | 73-120 | 0 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 44.30 | 89 | 45.68 | 91 | 80-122 | 3 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 46.20 | 92 | 46.62 | 93 | 70-120 | 1 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 44.51 | 89 | 46.06 | 92 | 64-142 | 3 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 38.61 | 77 | 39.69 | 79 | 52-136 | 3 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 42.74 | 85 | 42.75 | 86 | 77-125 | 0 | 0-24 | |
| Toluene | ND | 50.00 | 40.92 | 82 | 41.74 | 83 | 72-126 | 2 | 0-23 | |
| Trichloroethene | ND | 50.00 | 39.65 | 79 | 39.53 | 79 | 74-128 | 0 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 48.42 | 97 | 49.52 | 99 | 67-133 | 2 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 88.69 | 89 | 88.97 | 89 | 63-129 | 0 | 0-25 | |
| o-Xylene | ND | 50.00 | 46.86 | 94 | 47.12 | 94 | 62-128 | 1 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 45.07 | 90 | 46.50 | 93 | 68-134 | 3 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-05-0849-3 | Sample | Aqueous | GC/MS RR | 05/18/15 | 05/18/15 14:28 | 150518S007 | | | | |
| 15-05-0849-3 | Matrix Spike | Aqueous | GC/MS RR | 05/18/15 | 05/18/15 12:12 | 150518S007 | | | | |
| 15-05-0849-3 | Matrix Spike Duplicate | Aqueous | GC/MS RR | 05/18/15 | 05/18/15 12:39 | 150518S007 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 51.38 | 103 | 57.63 | 115 | 74-122 | 11 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 70.81 | 142 | 77.43 | 155 | 60-144 | 9 | 0-21 | 3 |
| Chlorobenzene | ND | 50.00 | 52.96 | 106 | 59.71 | 119 | 73-120 | 12 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 50.21 | 100 | 57.78 | 116 | 80-122 | 14 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 53.12 | 106 | 59.04 | 118 | 70-120 | 11 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 49.87 | 100 | 57.32 | 115 | 64-142 | 14 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 63.51 | 127 | 69.55 | 139 | 52-136 | 9 | 0-21 | 3 |
| Ethylbenzene | ND | 50.00 | 52.99 | 106 | 58.60 | 117 | 77-125 | 10 | 0-24 | |
| Toluene | ND | 50.00 | 51.46 | 103 | 58.25 | 116 | 72-126 | 12 | 0-23 | |
| Trichloroethene | 1.321 | 50.00 | 50.72 | 99 | 55.61 | 109 | 74-128 | 9 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 48.73 | 97 | 52.08 | 104 | 67-133 | 7 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 107.7 | 108 | 120.4 | 120 | 63-129 | 11 | 0-25 | |
| o-Xylene | ND | 50.00 | 53.87 | 108 | 60.42 | 121 | 62-128 | 11 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 52.41 | 105 | 61.14 | 122 | 68-134 | 15 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|--------------|-------------|------------|----------------|----------------|-----------------------|
| HSM 210 Lead | Sample | Aqueous | ICP/MS 04 | 05/07/15 00:00 | 05/11/15 21:21 | 150507SA4 |
| HSM 210 Lead | PDS | Aqueous | ICP/MS 04 | 05/07/15 00:00 | 05/11/15 21:09 | 150507SA4 |
| Parameter | Sample Conc. | Spike Added | PDS Conc. | PDS %Rec. | %Rec. CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.1035 | 103 | 75-125 | |
| Arsenic | 0.001117 | 0.1000 | 0.1029 | 102 | 75-125 | |
| Barium | 0.03955 | 0.1000 | 0.1414 | 102 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.1056 | 106 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.09975 | 100 | 75-125 | |
| Chromium | ND | 0.1000 | 0.1093 | 109 | 75-125 | |
| Copper | ND | 0.1000 | 0.1029 | 103 | 75-125 | |
| Lead | ND | 0.1000 | 0.1043 | 104 | 75-125 | |
| Nickel | 0.002058 | 0.1000 | 0.09701 | 95 | 75-125 | |
| Selenium | ND | 0.1000 | 0.09147 | 91 | 75-125 | |
| Silver | ND | 0.05000 | 0.04678 | 94 | 75-125 | |
| Thallium | ND | 0.1000 | 0.1024 | 102 | 75-125 | |
| Zinc | 0.01170 | 0.1000 | 0.1069 | 95 | 75-125 | |
| Aluminum | ND | 0.1000 | 0.1158 | 116 | 75-125 | |
| Iron | 0.05564 | 5.100 | 5.275 | 102 | 75-125 | |
| Manganese | 0.01498 | 0.1000 | 0.1143 | 99 | 75-125 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|---------------------|--------------------|------------------|-----------------------|-----------------------|-----------------------|
| FDHSM 210 Trail | Sample | Aqueous | ICP/MS 04 | 05/07/15 00:00 | 05/12/15 00:48 | 150507SA6 |
| FDHSM 210 Trail | PDS | Aqueous | ICP/MS 04 | 05/07/15 00:00 | 05/12/15 00:36 | 150507SA6 |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>PDS Conc.</u> | <u>PDS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Antimony | ND | 0.1000 | 0.1052 | 105 | 75-125 | |
| Arsenic | 0.001241 | 0.1000 | 0.1029 | 102 | 75-125 | |
| Barium | 0.03968 | 0.1000 | 0.1449 | 105 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.1190 | 119 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.1033 | 103 | 75-125 | |
| Chromium | ND | 0.1000 | 0.1113 | 111 | 75-125 | |
| Copper | ND | 0.1000 | 0.1045 | 105 | 75-125 | |
| Lead | ND | 0.1000 | 0.1062 | 106 | 75-125 | |
| Nickel | 0.002068 | 0.1000 | 0.09876 | 97 | 75-125 | |
| Selenium | ND | 0.1000 | 0.08735 | 87 | 75-125 | |
| Silver | ND | 0.05000 | 0.04823 | 96 | 75-125 | |
| Thallium | ND | 0.1000 | 0.1039 | 104 | 75-125 | |
| Zinc | 0.01100 | 0.1000 | 0.1063 | 95 | 75-125 | |
| Aluminum | ND | 0.1000 | 0.1223 | 122 | 75-125 | |
| Iron | 0.07078 | 5.100 | 5.518 | 107 | 75-125 | |
| Manganese | 0.02035 | 0.1000 | 0.1221 | 102 | 75-125 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---|------------------|---------------------|------------------|---------------|----------------|------------------------|
| HSM 240 Peak | Sample | Aqueous | PH1/BUR03 | N/A | 05/18/15 20:05 | F0518ALKD4 |
| HSM 240 Peak | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 05/18/15 20:05 | F0518ALKD4 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Alkalinity, Total (as CaCO ₃) | | 249.0 | 249.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---|------------------|---------------------|------------------|---------------|----------------|------------------------|
| HSM 210 Lead | Sample | Aqueous | PH1/BUR03 | N/A | 05/19/15 17:17 | F0519ALKD1 |
| HSM 210 Lead | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 05/19/15 17:17 | F0519ALKD1 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Alkalinity, Total (as CaCO ₃) | | 257.0 | 256.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|-------------------------------------|------------------|---------------------|------------------|---------------|----------------|------------------------|
| HSM 240 Peak | Sample | Aqueous | PH1/BUR03 | N/A | 05/18/15 20:05 | F0518HCOD4 |
| HSM 240 Peak | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 05/18/15 20:05 | F0518HCOD4 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Bicarbonate (as CaCO ₃) | | 249.0 | 249.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|-------------------------------------|------------------|---------------------|------------------|---------------|----------------|------------------------|
| HSM 210 Lead | Sample | Aqueous | PH1/BUR03 | N/A | 05/19/15 17:17 | F0519HCOD1 |
| HSM 210 Lead | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 05/19/15 17:17 | F0519HCOD1 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Bicarbonate (as CaCO ₃) | | 257.0 | 256.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| HSM 240 Peak | Sample | Aqueous | PH1/BUR03 | N/A | 05/18/15 20:05 | F0518CO3D4 |
| HSM 240 Peak | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 05/18/15 20:05 | F0518CO3D4 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| HSM 210 Lead | Sample | Aqueous | PH1/BUR03 | N/A | 05/19/15 17:17 | F0519CO3D1 |
| HSM 210 Lead | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 05/19/15 17:17 | F0519CO3D1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-05-0399-3 | Sample | Aqueous | SC 2 | 05/12/15 00:00 | 05/12/15 17:00 | F0512TDSD1 |
| 15-05-0399-3 | Sample Duplicate | Aqueous | SC 2 | 05/12/15 00:00 | 05/12/15 17:00 | F0512TDSD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Dissolved | 12760 | 12280 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|----------------|-------------|-----------------------|-----------------------|------------------------|
| HSM 210 Lead | Sample | Aqueous | SC 2 | 05/12/15 00:00 | 05/12/15 18:00 | F0512TDSD3 |
| HSM 210 Lead | Sample Duplicate | Aqueous | SC 2 | 05/12/15 00:00 | 05/12/15 18:00 | F0512TDSD3 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Dissolved | 595.0 | 611.0 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------------------|------------------|----------------|----------------|------------------------|
| HSM 240 Lead | Sample | Aqueous | N/A | 05/08/15 00:00 | 05/08/15 18:00 | F0508TSSD3 |
| HSM 240 Lead | Sample Duplicate | Aqueous | N/A | 05/08/15 00:00 | 05/08/15 18:00 | F0508TSSD3 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Solids, Total Suspended | | 376.6 | 377.0 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-05-0399-3 | Sample | Aqueous | PH 1 | N/A | 05/07/15 15:07 | F0507PHD1 |
| 15-05-0399-3 | Sample Duplicate | Aqueous | PH 1 | N/A | 05/07/15 15:07 | F0507PHD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|------------------|---------------------|------------------|------------|---------------|-------------------|
| pH | 7.060 | 7.100 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| HSM 270 Peak | Sample | Aqueous | PH 1 | N/A | 05/07/15 18:06 | F0507PHD3 |
| HSM 270 Peak | Sample Duplicate | Aqueous | PH 1 | N/A | 05/07/15 18:06 | F0507PHD3 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|------------------|---------------------|------------------|------------|---------------|-------------------|
| pH | 7.560 | 7.580 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 4500 N Org B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|----------------|--------------|-----------------------|-----------------------|------------------------|
| HSM 240 Trail | Sample | Aqueous | BUR05 | 05/21/15 00:00 | 05/21/15 16:00 | F0521TKND1 |
| HSM 240 Trail | Sample Duplicate | Aqueous | BUR05 | 05/21/15 00:00 | 05/21/15 16:00 | F0521TKND1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Total Kjeldahl Nitrogen | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 4500 N Org B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|----------------|--------------|-----------------------|-----------------------|------------------------|
| FDHSM 231 Trail | Sample | Aqueous | BUR05 | 05/22/15 00:00 | 05/22/15 16:40 | F0522TKND2 |
| FDHSM 231 Trail | Sample Duplicate | Aqueous | BUR05 | 05/22/15 00:00 | 05/22/15 16:40 | F0522TKND2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Total Kjeldahl Nitrogen | 0.5600 | 0.5600 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 099-12-906-5739 | LCS | Aqueous | IC 15 | N/A | 05/07/15 11:22 | 150507L01 |
| 099-12-906-5739 | LCSD | Aqueous | IC 15 | N/A | 05/07/15 11:39 | 150507L01 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| Fluoride | 2.500 | 2.432 | 97 | 2.443 | 98 | 90-110 | 0 | 0-15 | |
| Chloride | 50.00 | 49.78 | 100 | 49.92 | 100 | 90-110 | 0 | 0-15 | |
| Bromide | 5.000 | 4.982 | 100 | 5.001 | 100 | 90-110 | 0 | 0-15 | |
| Nitrate (as N) | 5.000 | 4.947 | 99 | 4.952 | 99 | 90-110 | 0 | 0-15 | |
| Sulfate | 50.00 | 49.25 | 99 | 49.44 | 99 | 90-110 | 0 | 0-15 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

SWCA Environmental Consultants
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San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|------------------|
| 099-12-906-5725 | LCS | Aqueous | IC 15 | N/A | 05/07/15 21:21 | 150507L02 |

| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
|----------------|-------------|-----------------|-----------|----------|------------|
| Fluoride | 2.500 | 2.473 | 99 | 90-110 | |
| Chloride | 50.00 | 50.06 | 100 | 90-110 | |
| Bromide | 5.000 | 5.013 | 100 | 90-110 | |
| Nitrate (as N) | 5.000 | 4.966 | 99 | 90-110 | |
| Sulfate | 50.00 | 49.56 | 99 | 90-110 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
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Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-739-104 | LCS | Aqueous | ACA 1 | N/A | 05/20/15 19:29 | 150520L01 | | | |
| 099-12-739-104 | LCSD | Aqueous | ACA 1 | N/A | 05/20/15 19:29 | 150520L01 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Phosphorus, Total | 0.2000 | 0.2042 | 102 | 0.2061 | 103 | 90-110 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
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Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------------|------------|---------------|-------------------|
| 099-12-739-105 | LCS | Aqueous | ACA 1 | N/A | 05/20/15 19:29 | 150520L02 | | | |
| 099-12-739-105 | LCSD | Aqueous | ACA 1 | N/A | 05/20/15 19:29 | 150520L02 | | | |
| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Phosphorus, Total | 0.2000 | 0.2012 | 101 | 0.2067 | 103 | 90-110 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-859-708 | LCS | Aqueous | PH1/BUR03 | N/A | 05/18/15 20:05 | F0518ALKB4 | | | |
| 099-15-859-708 | LCSD | Aqueous | PH1/BUR03 | N/A | 05/18/15 20:05 | F0518ALKB4 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 100.0 | 98.00 | 98 | 98.00 | 98 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-859-703 | LCS | Aqueous | PH1/BUR03 | N/A | 05/19/15 17:17 | F0519ALKB1 | | | |
| 099-15-859-703 | LCSD | Aqueous | PH1/BUR03 | N/A | 05/19/15 17:17 | F0519ALKB1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 100.0 | 98.00 | 98 | 98.00 | 98 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4569 | LCS | Aqueous | SC 2 | 05/12/15 | 05/12/15 17:00 | F0512TDSL1 | | | |
| 099-12-180-4569 | LCSD | Aqueous | SC 2 | 05/12/15 | 05/12/15 17:00 | F0512TDSL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 95.00 | 95 | 90.00 | 90 | 80-120 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4574 | LCS | Aqueous | SC 2 | 05/12/15 | 05/12/15 18:00 | F0512TDSL3 | | | |
| 099-12-180-4574 | LCSD | Aqueous | SC 2 | 05/12/15 | 05/12/15 18:00 | F0512TDSL3 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 115.0 | 115 | 110.0 | 110 | 80-120 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------------|------------|---------------|-------------------|
| 099-09-010-7166 | LCS | Aqueous | N/A | 05/08/15 | 05/08/15 17:00 | F0508TSSL1 | | | |
| 099-09-010-7166 | LCSD | Aqueous | N/A | 05/08/15 | 05/08/15 17:00 | F0508TSSL1 | | | |
| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Solids, Total Suspended | 100.0 | 93.00 | 93 | 92.00 | 92 | 80-120 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-09-010-7169 | LCS | Aqueous | N/A | 05/08/15 | 05/08/15 18:00 | F0508TSSL3 | | | |
| 099-09-010-7169 | LCSD | Aqueous | N/A | 05/08/15 | 05/08/15 18:00 | F0508TSSL3 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Suspended | 100.0 | 88.00 | 88 | 89.00 | 89 | 80-120 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-097-5645 | LCS | Aqueous | TOC 8 | 05/22/15 | 05/23/15 06:08 | F0522TOCL1 | | | |
| 099-05-097-5645 | LCSD | Aqueous | TOC 8 | 05/22/15 | 05/23/15 06:08 | F0522TOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 10.00 | 9.470 | 95 | 9.930 | 99 | 80-120 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-097-5647 | LCS | Aqueous | TOC 8 | 05/26/15 | 05/26/15 22:23 | F0526TOCL1 | | | |
| 099-05-097-5647 | LCSD | Aqueous | TOC 8 | 05/26/15 | 05/26/15 22:23 | F0526TOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 10.00 | 9.140 | 91 | 9.530 | 95 | 80-120 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

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San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-097-5656 | LCS | Aqueous | TOC 8 | 05/29/15 | 05/30/15 06:47 | F0529TOCL1 | | | |
| 099-05-097-5656 | LCSD | Aqueous | TOC 8 | 05/29/15 | 05/30/15 06:47 | F0529TOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 10.00 | 9.410 | 94 | 9.720 | 97 | 80-120 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

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San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-115-1432 | LCS | Aqueous | TOC 8 | 05/13/15 | 05/14/15 14:53 | F0513DOCL1 | | | |
| 099-05-115-1432 | LCSD | Aqueous | TOC 8 | 05/13/15 | 05/14/15 14:53 | F0513DOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Dissolved Organic | 10.00 | 11.00 | 110 | 11.00 | 110 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

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San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-115-1433 | LCS | Aqueous | TOC 8 | 05/13/15 | 05/14/15 19:49 | F0513DOCL2 | | | |
| 099-05-115-1433 | LCSD | Aqueous | TOC 8 | 05/13/15 | 05/14/15 19:49 | F0513DOCL2 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Dissolved Organic | 10.00 | 10.30 | 103 | 10.60 | 106 | 80-120 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

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Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-15-683-1264 | LCS | Aqueous | ICP 7300 | 05/07/15 | 05/21/15 11:29 | 150507LA5 |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Calcium | | 0.5000 | 0.4872 | 97 | 80-120 | |
| Magnesium | | 0.5000 | 0.4921 | 98 | 80-120 | |
| Potassium | | 5.000 | 5.107 | 102 | 80-120 | |
| Sodium | | 5.000 | 5.213 | 104 | 80-120 | |
| Strontium | | 0.5000 | 0.5093 | 102 | 80-120 | |
| Silicon | | 0.5000 | 0.4878 | 98 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

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San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-15-683-1265 | LCS | Aqueous | ICP 7300 | 05/07/15 | 05/16/15 21:04 | 150507LA6A |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Calcium | | 0.5000 | 0.4472 | 89 | 80-120 | |
| Magnesium | | 0.5000 | 0.4943 | 99 | 80-120 | |
| Potassium | | 5.000 | 4.896 | 98 | 80-120 | |
| Sodium | | 5.000 | 4.621 | 92 | 80-120 | |
| Strontium | | 0.5000 | 0.4853 | 97 | 80-120 | |
| Silicon | | 0.5000 | 0.4440 | 89 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits

Quality Control - LCS

SWCA Environmental Consultants
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San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-820 | LCS | Aqueous | ICP/MS 04 | 05/07/15 | 05/09/15 03:00 | 150507LA4F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.09868 | 99 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.09926 | 99 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.09742 | 97 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.09743 | 97 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.09861 | 99 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.1015 | 101 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.09934 | 99 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.09646 | 96 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.09565 | 96 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.09817 | 98 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.04706 | 94 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.09544 | 95 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.09991 | 100 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.1016 | 102 | 80-120 | 73-127 | |
| Iron | 5.100 | 5.129 | 101 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.09955 | 100 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-821 | LCS | Aqueous | ICP/MS 04 | 05/07/15 | 05/09/15 03:04 | 150507LA6F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.09791 | 98 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.09904 | 99 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.09653 | 97 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.09677 | 97 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.09788 | 98 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.1027 | 103 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.09893 | 99 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.09679 | 97 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.09447 | 94 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.09717 | 97 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.04727 | 95 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.09449 | 94 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.09940 | 99 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.1004 | 100 | 80-120 | 73-127 | |
| Iron | 5.100 | 5.128 | 101 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.09839 | 98 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|------------------|
| 099-15-763-554 | LCS | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:00 | 150511L04 |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.01000 | 0.009965 | 100 | 85-121 | |

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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|------------------|
| 099-15-763-555 | LCS | Aqueous | Mercury 04 | 05/11/15 | 05/11/15 18:02 | 150511L05 |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.01000 | 0.009819 | 98 | 85-121 | |

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Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 099-16-376-13 | LCS | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 12:27 | 150512L21 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Caffeine | | 100.0 | 96.60 | 97 | 80-120 | |

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Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 1694
Method: EPA 1694 (M) Caffeine

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 099-16-376-14 | LCS | Aqueous | LC/TQ 2 | 05/12/15 | 05/13/15 16:51 | 150512L22 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Caffeine | | 100.0 | 97.59 | 98 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-529-803 | LCS | Aqueous | | GC 44 | 05/08/15 | 05/11/15 14:43 | 150508L05 | | | |
| 099-12-529-803 | LCSD | Aqueous | | GC 44 | 05/08/15 | 05/11/15 14:57 | 150508L05 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Alpha-BHC | 0.5000 | 0.4357 | 87 | 0.4463 | 89 | 50-135 | 36-149 | 2 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.4556 | 91 | 0.4650 | 93 | 50-135 | 36-149 | 2 | 0-25 | |
| Beta-BHC | 0.5000 | 0.3816 | 76 | 0.4048 | 81 | 50-135 | 36-149 | 6 | 0-25 | |
| Heptachlor | 0.5000 | 0.4470 | 89 | 0.4230 | 85 | 50-135 | 36-149 | 6 | 0-25 | |
| Delta-BHC | 0.5000 | 0.4066 | 81 | 0.4293 | 86 | 50-135 | 36-149 | 5 | 0-25 | |
| Aldrin | 0.5000 | 0.4055 | 81 | 0.3680 | 74 | 50-135 | 36-149 | 10 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.4239 | 85 | 0.4265 | 85 | 50-135 | 36-149 | 1 | 0-25 | |
| Endosulfan I | 0.5000 | 0.4150 | 83 | 0.4206 | 84 | 50-135 | 36-149 | 1 | 0-25 | |
| Dieldrin | 0.5000 | 0.4344 | 87 | 0.4364 | 87 | 50-135 | 36-149 | 0 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.4265 | 85 | 0.4297 | 86 | 50-135 | 36-149 | 1 | 0-25 | |
| Endrin | 0.5000 | 0.3544 | 71 | 0.3670 | 73 | 50-135 | 36-149 | 3 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.4570 | 91 | 0.4520 | 90 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.4255 | 85 | 0.4192 | 84 | 50-135 | 36-149 | 1 | 0-25 | |
| Endosulfan II | 0.5000 | 0.4235 | 85 | 0.4190 | 84 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.3793 | 76 | 0.3829 | 77 | 50-135 | 36-149 | 1 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.3934 | 79 | 0.3893 | 78 | 50-135 | 36-149 | 1 | 0-25 | |
| Methoxychlor | 0.5000 | 0.3855 | 77 | 0.3814 | 76 | 50-135 | 36-149 | 1 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-529-804 | LCS | Aqueous | | GC 44 | 05/09/15 | 05/11/15 19:30 | 150509L11 | | | |
| 099-12-529-804 | LCSD | Aqueous | | GC 44 | 05/09/15 | 05/11/15 19:45 | 150509L11 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Alpha-BHC | 0.5000 | 0.4492 | 90 | 0.4483 | 90 | 50-135 | 36-149 | 0 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.4664 | 93 | 0.4624 | 92 | 50-135 | 36-149 | 1 | 0-25 | |
| Beta-BHC | 0.5000 | 0.3938 | 79 | 0.3853 | 77 | 50-135 | 36-149 | 2 | 0-25 | |
| Heptachlor | 0.5000 | 0.4370 | 87 | 0.4351 | 87 | 50-135 | 36-149 | 0 | 0-25 | |
| Delta-BHC | 0.5000 | 0.4353 | 87 | 0.4263 | 85 | 50-135 | 36-149 | 2 | 0-25 | |
| Aldrin | 0.5000 | 0.4147 | 83 | 0.4141 | 83 | 50-135 | 36-149 | 0 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.4142 | 83 | 0.4139 | 83 | 50-135 | 36-149 | 0 | 0-25 | |
| Endosulfan I | 0.5000 | 0.4094 | 82 | 0.4075 | 82 | 50-135 | 36-149 | 0 | 0-25 | |
| Dieldrin | 0.5000 | 0.4194 | 84 | 0.4166 | 83 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.4131 | 83 | 0.4072 | 81 | 50-135 | 36-149 | 1 | 0-25 | |
| Endrin | 0.5000 | 0.3959 | 79 | 0.3814 | 76 | 50-135 | 36-149 | 4 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.4193 | 84 | 0.4188 | 84 | 50-135 | 36-149 | 0 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.4100 | 82 | 0.4087 | 82 | 50-135 | 36-149 | 0 | 0-25 | |
| Endosulfan II | 0.5000 | 0.4003 | 80 | 0.3966 | 79 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.3780 | 76 | 0.3567 | 71 | 50-135 | 36-149 | 6 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.3916 | 78 | 0.3862 | 77 | 50-135 | 36-149 | 1 | 0-25 | |
| Methoxychlor | 0.5000 | 0.3799 | 76 | 0.3540 | 71 | 50-135 | 36-149 | 7 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-533-1035 | LCS | Aqueous | GC 31 | 05/08/15 | 05/11/15 18:00 | 150508L15 | | | |
| 099-12-533-1035 | LCSD | Aqueous | GC 31 | 05/08/15 | 05/11/15 18:19 | 150508L15 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Aroclor-1016 | 2.000 | 1.857 | 93 | 2.217 | 111 | 50-135 | 18 | 0-25 | |
| Aroclor-1260 | 2.000 | 1.973 | 99 | 2.050 | 102 | 50-135 | 4 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-533-1034 | LCS | Aqueous | GC 31 | 05/09/15 | 05/11/15 16:43 | 150509L12 | | | |
| 099-12-533-1034 | LCSD | Aqueous | GC 31 | 05/09/15 | 05/11/15 17:02 | 150509L12 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Aroclor-1016 | 2.000 | 1.592 | 80 | 1.744 | 87 | 50-135 | 9 | 0-25 | |
| Aroclor-1260 | 2.000 | 1.740 | 87 | 1.797 | 90 | 50-135 | 3 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|--------|-----|--------|------------|
| 099-15-963-93 | LCS | Aqueous | GC 26 | 05/09/15 | 05/12/15 11:57 | 150509L04 | | | | |
| 099-15-963-93 | LCSD | Aqueous | GC 26 | 05/09/15 | 05/12/15 12:41 | 150509L04 | | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.03887 | 97 | 0.03402 | 85 | 30-130 | 13-147 | 13 | 0-30 | |
| Bolstar | 0.04000 | 0.04686 | 117 | 0.04654 | 116 | 30-130 | 13-147 | 1 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.04314 | 108 | 0.04323 | 108 | 30-130 | 13-147 | 0 | 0-30 | |
| Coumaphos | 0.04000 | 0.04329 | 108 | 0.03855 | 96 | 30-130 | 13-147 | 12 | 0-30 | |
| Diazinon | 0.04000 | 0.04485 | 112 | 0.04375 | 109 | 30-130 | 13-147 | 2 | 0-30 | |
| Disulfoton | 0.04000 | 0.05020 | 126 | 0.04471 | 112 | 30-130 | 13-147 | 12 | 0-30 | |
| Ethoprop | 0.04000 | 0.05043 | 126 | 0.04957 | 124 | 30-130 | 13-147 | 2 | 0-30 | |
| Fensulfothion | 0.04000 | 0.04603 | 115 | 0.04442 | 111 | 30-130 | 13-147 | 4 | 0-30 | |
| Fenthion | 0.04000 | 0.04298 | 107 | 0.04221 | 106 | 30-130 | 13-147 | 2 | 0-30 | |
| Merphos | 0.04000 | 0.04717 | 118 | 0.04682 | 117 | 30-130 | 13-147 | 1 | 0-30 | |
| Methyl Parathion | 0.04000 | 0.04441 | 111 | 0.04312 | 108 | 30-130 | 13-147 | 3 | 0-30 | |
| Phorate | 0.04000 | 0.04542 | 114 | 0.04621 | 116 | 30-130 | 13-147 | 2 | 0-30 | |
| Ronnel | 0.04000 | 0.04502 | 113 | 0.04438 | 111 | 30-130 | 13-147 | 1 | 0-30 | |
| Stirophos | 0.04000 | 0.03129 | 78 | 0.03018 | 75 | 30-130 | 13-147 | 4 | 0-30 | |
| Tokuthion | 0.04000 | 0.04523 | 113 | 0.04528 | 113 | 30-130 | 13-147 | 0 | 0-30 | |
| Trichloronate | 0.04000 | 0.04820 | 120 | 0.04776 | 119 | 30-130 | 13-147 | 1 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-963-94 | LCS | Aqueous | | GC 26 | 05/09/15 | 05/13/15 06:20 | 150509L05 | | | |
| 099-15-963-94 | LCSD | Aqueous | | GC 26 | 05/09/15 | 05/13/15 07:04 | 150509L05 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.03277 | 82 | 0.03283 | 82 | 30-130 | 13-147 | 0 | 0-30 | |
| Bolstar | 0.04000 | 0.04077 | 102 | 0.03964 | 99 | 30-130 | 13-147 | 3 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.03894 | 97 | 0.03751 | 94 | 30-130 | 13-147 | 4 | 0-30 | |
| Coumaphos | 0.04000 | 0.03809 | 95 | 0.03715 | 93 | 30-130 | 13-147 | 2 | 0-30 | |
| Diazinon | 0.04000 | 0.04098 | 102 | 0.04033 | 101 | 30-130 | 13-147 | 2 | 0-30 | |
| Disulfoton | 0.04000 | 0.04931 | 123 | 0.03837 | 96 | 30-130 | 13-147 | 25 | 0-30 | |
| Ethoprop | 0.04000 | 0.04532 | 113 | 0.04402 | 110 | 30-130 | 13-147 | 3 | 0-30 | |
| Fensulfothion | 0.04000 | 0.03810 | 95 | 0.03758 | 94 | 30-130 | 13-147 | 1 | 0-30 | |
| Fenthion | 0.04000 | 0.03740 | 94 | 0.03604 | 90 | 30-130 | 13-147 | 4 | 0-30 | |
| Merphos | 0.04000 | 0.04159 | 104 | 0.03825 | 96 | 30-130 | 13-147 | 8 | 0-30 | |
| Methyl Parathion | 0.04000 | 0.03571 | 89 | 0.03549 | 89 | 30-130 | 13-147 | 1 | 0-30 | |
| Phorate | 0.04000 | 0.05132 | 128 | 0.04989 | 125 | 30-130 | 13-147 | 3 | 0-30 | |
| Ronnel | 0.04000 | 0.03973 | 99 | 0.03891 | 97 | 30-130 | 13-147 | 2 | 0-30 | |
| Stirophos | 0.04000 | 0.02626 | 66 | 0.02538 | 63 | 30-130 | 13-147 | 3 | 0-30 | |
| Tokuthion | 0.04000 | 0.03960 | 99 | 0.03869 | 97 | 30-130 | 13-147 | 2 | 0-30 | |
| Trichloronate | 0.04000 | 0.04257 | 106 | 0.04161 | 104 | 30-130 | 13-147 | 2 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 095-01-034-644 | LCS | Aqueous | GC 40 | 05/08/15 | 05/15/15 00:44 | 150508L12 |
| 095-01-034-644 | LCSD | Aqueous | GC 40 | 05/08/15 | 05/15/15 01:07 | 150508L12 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| 2,4-D | 20.00 | 19.36 | 97 | 17.82 | 89 | 30-130 | 8 | 0-30 | |
| 2,4,5-T | 2.000 | 1.880 | 94 | 1.750 | 88 | 30-130 | 7 | 0-30 | |
| 2,4-DB | 20.00 | 19.66 | 98 | 18.12 | 91 | 30-130 | 8 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 095-01-034-645 | LCS | Aqueous | GC 40 | 05/08/15 | 05/15/15 01:53 | 150508L13 |
| 095-01-034-645 | LCSD | Aqueous | GC 40 | 05/08/15 | 05/15/15 02:40 | 150508L13 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| 2,4-D | 20.00 | 17.68 | 88 | 17.37 | 87 | 30-130 | 2 | 0-30 | |
| 2,4,5-T | 2.000 | 1.715 | 86 | 1.710 | 86 | 30-130 | 0 | 0-30 | |
| 2,4-DB | 20.00 | 17.88 | 89 | 17.63 | 88 | 30-130 | 1 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-4040 | LCS | Aqueous | | GC/MS TT | 05/11/15 | 05/12/15 13:16 | 150511L04 | | | |
| 095-01-003-4040 | LCSD | Aqueous | | GC/MS TT | 05/11/15 | 05/12/15 13:35 | 150511L04 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 200.0 | 189.7 | 95 | 196.8 | 98 | 61-120 | 51-130 | 4 | 0-20 | |
| Acenaphthylene | 200.0 | 179.9 | 90 | 188.1 | 94 | 55-120 | 44-131 | 4 | 0-20 | |
| Butyl Benzyl Phthalate | 200.0 | 164.6 | 82 | 169.1 | 85 | 56-122 | 45-133 | 3 | 0-20 | |
| 4-Chloro-3-Methylphenol | 200.0 | 146.8 | 73 | 147.9 | 74 | 52-120 | 41-131 | 1 | 0-20 | |
| 2-Chlorophenol | 200.0 | 169.0 | 84 | 165.1 | 83 | 47-120 | 35-132 | 2 | 0-20 | |
| 1,4-Dichlorobenzene | 200.0 | 168.9 | 84 | 165.2 | 83 | 36-120 | 22-134 | 2 | 0-20 | |
| Dimethyl Phthalate | 200.0 | 178.6 | 89 | 185.4 | 93 | 60-120 | 50-130 | 4 | 0-20 | |
| 2,4-Dinitrotoluene | 200.0 | 183.7 | 92 | 184.1 | 92 | 61-121 | 51-131 | 0 | 0-20 | |
| Fluorene | 200.0 | 187.1 | 94 | 193.5 | 97 | 67-120 | 58-129 | 3 | 0-20 | |
| N-Nitroso-di-n-propylamine | 200.0 | 149.8 | 75 | 148.9 | 74 | 39-123 | 25-137 | 1 | 0-20 | |
| Naphthalene | 200.0 | 166.4 | 83 | 169.8 | 85 | 54-120 | 43-131 | 2 | 0-20 | |
| 4-Nitrophenol | 200.0 | 95.62 | 48 | 96.86 | 48 | 14-120 | 0-138 | 1 | 0-20 | |
| Pentachlorophenol | 200.0 | 117.8 | 59 | 119.6 | 60 | 31-127 | 15-143 | 2 | 0-20 | |
| Phenol | 200.0 | 151.8 | 76 | 149.8 | 75 | 17-120 | 0-137 | 1 | 0-20 | |
| Pyrene | 200.0 | 156.5 | 78 | 161.3 | 81 | 58-124 | 47-135 | 3 | 0-20 | |
| 1,2,4-Trichlorobenzene | 200.0 | 163.7 | 82 | 165.5 | 83 | 49-120 | 37-132 | 1 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-4039 | LCS | Aqueous | | GC/MS CCC | 05/09/15 | 05/12/15 13:30 | 150509L06 | | | |
| 095-01-003-4039 | LCSD | Aqueous | | GC/MS CCC | 05/09/15 | 05/12/15 13:48 | 150509L06 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 200.0 | 179.8 | 90 | 198.6 | 99 | 61-120 | 51-130 | 10 | 0-20 | |
| Acenaphthylene | 200.0 | 177.5 | 89 | 196.1 | 98 | 55-120 | 44-131 | 10 | 0-20 | |
| Butyl Benzyl Phthalate | 200.0 | 183.2 | 92 | 201.7 | 101 | 56-122 | 45-133 | 10 | 0-20 | |
| 4-Chloro-3-Methylphenol | 200.0 | 150.4 | 75 | 169.0 | 84 | 52-120 | 41-131 | 12 | 0-20 | |
| 2-Chlorophenol | 200.0 | 157.6 | 79 | 174.7 | 87 | 47-120 | 35-132 | 10 | 0-20 | |
| 1,4-Dichlorobenzene | 200.0 | 150.9 | 75 | 162.5 | 81 | 36-120 | 22-134 | 7 | 0-20 | |
| Dimethyl Phthalate | 200.0 | 174.7 | 87 | 193.2 | 97 | 60-120 | 50-130 | 10 | 0-20 | |
| 2,4-Dinitrotoluene | 200.0 | 168.7 | 84 | 183.7 | 92 | 61-121 | 51-131 | 9 | 0-20 | |
| Fluorene | 200.0 | 181.8 | 91 | 199.6 | 100 | 67-120 | 58-129 | 9 | 0-20 | |
| N-Nitroso-di-n-propylamine | 200.0 | 168.6 | 84 | 189.4 | 95 | 39-123 | 25-137 | 12 | 0-20 | |
| Naphthalene | 200.0 | 153.3 | 77 | 167.9 | 84 | 54-120 | 43-131 | 9 | 0-20 | |
| 4-Nitrophenol | 200.0 | 160.1 | 80 | 178.9 | 89 | 14-120 | 0-138 | 11 | 0-20 | |
| Pentachlorophenol | 200.0 | 138.2 | 69 | 147.8 | 74 | 31-127 | 15-143 | 7 | 0-20 | |
| Phenol | 200.0 | 162.0 | 81 | 180.3 | 90 | 17-120 | 0-137 | 11 | 0-20 | |
| Pyrene | 200.0 | 160.9 | 80 | 177.1 | 89 | 58-124 | 47-135 | 10 | 0-20 | |
| 1,2,4-Trichlorobenzene | 200.0 | 146.5 | 73 | 158.3 | 79 | 49-120 | 37-132 | 8 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

Page 34 of 38

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-17133 | LCS | Aqueous | GC/MS O | 05/12/15 | 05/12/15 21:13 | 150512L041 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 49.24 | 98 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 50.99 | 102 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 47.41 | 95 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 54.52 | 109 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 47.80 | 96 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 50.55 | 101 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 49.63 | 99 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 49.24 | 98 | 80-123 | 73-130 | |
| Toluene | 50.00 | 48.78 | 98 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 49.21 | 98 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 45.91 | 92 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 95.47 | 95 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 46.63 | 93 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 48.85 | 98 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

Page 35 of 38

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-17137 | LCS | Aqueous | GC/MS O | 05/13/15 | 05/13/15 15:39 | 150513L015 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 48.54 | 97 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 49.47 | 99 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 47.17 | 94 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 51.81 | 104 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 46.82 | 94 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 47.56 | 95 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 47.46 | 95 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 49.07 | 98 | 80-123 | 73-130 | |
| Toluene | 50.00 | 48.58 | 97 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 49.17 | 98 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 44.37 | 89 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 96.67 | 97 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 46.64 | 93 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 49.00 | 98 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-17160 | LCS | Aqueous | GC/MS O | 05/15/15 | 05/16/15 00:25 | 150515L039 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 50.31 | 101 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 41.18 | 82 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 49.44 | 99 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 53.24 | 106 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 48.41 | 97 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 51.80 | 104 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 45.56 | 91 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 51.28 | 103 | 80-123 | 73-130 | |
| Toluene | 50.00 | 50.03 | 100 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 50.84 | 102 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 39.22 | 78 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 100.1 | 100 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 48.46 | 97 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 43.03 | 86 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-17134 | LCS | Aqueous | GC/MS BB | 05/09/15 | 05/09/15 11:06 | 150509L027 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 40.21 | 80 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 40.44 | 81 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 44.57 | 89 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 45.42 | 91 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 47.40 | 95 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 46.06 | 92 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 40.53 | 81 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 44.51 | 89 | 80-123 | 73-130 | |
| Toluene | 50.00 | 41.72 | 83 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 40.51 | 81 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 49.57 | 99 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 93.19 | 93 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 49.04 | 98 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 44.82 | 90 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 05/07/15
Work Order: 15-05-0402
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-17170 | LCS | Aqueous | GC/MS RR | 05/18/15 | 05/18/15 11:29 | 150518L017 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 47.55 | 95 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 61.59 | 123 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 50.18 | 100 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 49.22 | 98 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 50.44 | 101 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 48.05 | 96 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 55.67 | 111 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 48.75 | 97 | 80-123 | 73-130 | |
| Toluene | 50.00 | 47.72 | 95 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 45.07 | 90 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 42.54 | 85 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 99.93 | 100 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 50.79 | 102 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 50.22 | 100 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Sample Analysis Summary Report

Work Order: 15-05-0402

Page 1 of 1

| <u>Method</u> | <u>Extraction</u> | <u>Chemist ID</u> | <u>Instrument</u> | <u>Analytical Location</u> |
|-----------------------|-------------------|-------------------|-------------------|----------------------------|
| EPA 1694 (M) Caffeine | EPA 1694 | 262 | LC/TQ 2 | 1 |
| EPA 300.0 | N/A | 834 | IC 15 | 1 |
| EPA 365.1 | N/A | 735 | ACA 1 | 1 |
| EPA 6010B | EPA 3005A Filt. | 935 | ICP 7300 | 1 |
| EPA 6020 | EPA 3005A Filt. | 598 | ICP/MS 04 | 1 |
| EPA 7470A | EPA 7470A Filt. | 598 | Mercury 04 | 1 |
| EPA 8081A | EPA 3510C | 669 | GC 44 | 1 |
| EPA 8082 | EPA 3510C | 669 | GC 31 | 1 |
| EPA 8141A | EPA 3510C | 949 | GC 26 | 1 |
| EPA 8151A | EPA 8151A | 944 | GC 40 | 1 |
| EPA 8260B | EPA 5030C | 131 | GC/MS RR | 2 |
| EPA 8260B | EPA 5030C | 975 | GC/MS O | 2 |
| EPA 8260B | EPA 5030C | 975 | GC/MS BB | 2 |
| EPA 8270C | EPA 3510C | 923 | GC/MS CCC | 1 |
| SM 2320B | N/A | 688 | PH1/BUR03 | 1 |
| SM 2540 C | N/A | 689 | SC 2 | 1 |
| SM 2540 D | N/A | 689 | N/A | 1 |
| SM 4500 H+ B | N/A | 688 | PH 1 | 1 |
| SM 4500 N Org B | N/A | 685 | BUR05 | 1 |
| SM 5310 B | N/A | 735 | TOC 8 | 1 |

[Return to Contents](#)

Location 1: 7440 Lincoln Way, Garden Grove, CA 92841

Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841

Glossary of Terms and Qualifiers

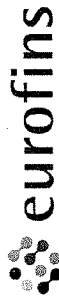
Work Order: 15-05-0402

Page 1 of 1

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|--|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| CI | See case narrative. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.



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For courier service / sample drop off information, contact us26_sales@eurofinsus.com or call us.

CHAIN OF CUSTODY RECORD

DATE: 5-6-2015

PAGE: 2 OF 3

WORK # / LAB USE ONLY

15-05-0402

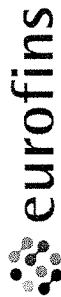
| | | | | | |
|---|--|---|--|---------------------------------|--|
| LABORATORY CLIENT: SWCA Environmental Consultants | | CLIENT PROJECT NAME / NUMBER: EAA 27122 | | P.O. NO.: 27122.02.08 | |
| ADDRESS: 6200 UTSA Blvd. Suite 102 | | PROJECT CONTACT: Philip Pearce | | SAMPLER(S) (PRINT): PHIL PEARCE | |
| CITY: San Antonio | | STATE: TX | | ZIP: 78249-1618 | |
| TEL: 210.877.2847 | | E-MAIL: ppearce@swca.com | | | |

REQUESTED ANALYSES

| | | | | | | | | | |
|---|--------------|--------|-------|----------|--------|--------------|-------------|-----------|----------------|
| TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD"): | | | | | | | | | |
| <input type="checkbox"/> SAME DAY <input type="checkbox"/> 24 HR <input type="checkbox"/> 48 HR <input type="checkbox"/> 72 HR <input checked="" type="checkbox"/> 10 Days (standard) | | | | | | | | | |
| <input type="checkbox"/> COELT EDF | | | | | | | | | |
| SPECIAL INSTRUCTIONS: | | | | | | | | | |
| Samples for EPA 6010B, EPA 6020, and EPA 7470 Metals are field filtered | | | | | | | | | |
| EPA 6020 Metals: Zn, Ti, Se, Sb, Pb, Ni, Cu, Cr, Cd, Be, Ba, As, Al, Ag, Fe, Mn | | | | | | | | | |
| Please also analyze each sample for pH | | | | | | | | | |
| LAB USE ONLY | SAMPLE ID | DATE | TIME | SAMPLING | MATRIX | NO. OF CONT. | Unpreserved | Preserved | Field Filtered |
| 8 | HSM 210 Peak | 5-5-15 | 23:24 | | NPW | 9 | 2 | 7 | 1 |
| 9 | HSM 230 Peak | 5-5-15 | 23:35 | | NPW | 9 | 2 | 7 | 1 |
| 10 | HSM 231 Peak | 5-5-15 | 22:11 | | NPW | 9 | 2 | 7 | 1 |
| 11 | HSM 240 Peak | 5-6-15 | 00:05 | | NPW | 9 | 2 | 7 | 1 |
| 12 | HSM 250 Peak | 5-5-15 | 23:30 | | NPW | 9 | 2 | 7 | 1 |
| 13 | HSM 260 Peak | 5-5-15 | 23:49 | | NPW | 9 | 2 | 7 | 1 |
| 14 | HSM 270 Peak | 5-6-15 | 00:06 | | NPW | 9 | 2 | 7 | 1 |
| 15 | pp#008 TB06 | 5-6-14 | 10:11 | | NPW | 1 | | | |

| | | | |
|------------------------------|--|--------------------------------------|--|
| Relinquished by: (Signature) | | Received by: (Signature/Affiliation) | |
| Relinquished by: (Signature) | | Received by: (Signature/Affiliation) | |
| Relinquished by: (Signature) | | Received by: (Signature/Affiliation) | |

Date: 05/07/15 Time: 1030



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CHAIN OF CUSTODY RECORD

WO # / LAB USE ONLY:
15-05-0402

DATE: 5-6-2015

PAGE: 3 OF 3

| | | | |
|---|--------------------------|---|---------------------------------|
| LABORATORY CLIENT: SWCA Environmental Consultants | | CLIENT PROJECT NAME / NUMBER: EAA 27122 | |
| ADDRESS: 6200 UTSA Blvd, Suite 102 | | P.O. NO.: 27122.02.08 | |
| CITY: San Antonio | STATE: TX | ZIP: 78249-1618 | SAMPLER(S) (PRINT): PHIL PEARCE |
| TEL: 210.877.2847 | E-MAIL: PPearce@swca.com | PROJECT CONTACT: Philip Pearce | |

| | | | | | |
|---|--|---|--|---|--|
| SPECIAL INSTRUCTIONS: Samples for EPA 6010B, EPA 6020, and EPA 7470 Metals are field filtered EPA 6020 Metals: Zn, Ti, Se, Sb, Pb, Ni, Cu, Cr, Cd, Be, Ba, As, Al, Ag, Fe, Mn Please also analyze each sample for pH | | | | LOG CODE: Unpreserved Preserved Field Filtered | |
| GLOBAL ID: | | TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD"): <input type="checkbox"/> SAME DAY <input type="checkbox"/> 24 HR <input type="checkbox"/> 48 HR <input checked="" type="checkbox"/> 72 HR <input type="checkbox"/> 10 Days (standard) | | LAB USE ONLY | |
| COELT EDF | | NO. OF CONT. | | NO. OF CONT. | |
| SAMPLE ID | | DATE | | TIME | |
| MATRIX | | NO. | | CONT. | |
| HSM 210 Trail | | 5-6-15 | | 01:30 | |
| HSM 230 Trail | | 5-6-15 | | 02:15 | |
| HSM 231 Trail | | 5-6-15 | | 02:40 | |
| HSM 240 Trail | | 5-6-15 | | 03:05 | |
| HSM 250 Trail | | 5-6-15 | | 03:58 | |
| HSM 260 Trail | | 5-6-15 | | 04:20 | |
| HSM 270 Trail | | 5-6-15 | | 04:40 | |
| FDHSM 210 Trail | | 5-6-15 | | 01:30 | |
| FDHSM 230 Trail | | 5-6-15 | | 02:15 | |
| FDHSM231 Trail | | 5-6-15 | | 02:40 | |

| | | | |
|--|---|--|---|
| Requested Analyses | | Please check box or fill in blank as needed. | |
| EPA 6010B (Si, Ca, Mg, K, Na, S) | X | EPA 8081A Organochlorine Pesticides | X |
| EPA 7470A Mercury | X | EPA 1694 (M) Caffeine | X |
| EPA 8082 PCB Aroclors | X | EPA 8141A Organophosphorus Pest | X |
| EPA 8151A Chlorinated Herbicides | X | EPA 8270C Semi-Volatile Organics | X |
| EPA 8260B Volatile Organics | X | EPA 365.1 Total Phosphorus | X |
| SM 2320B Alkalinity: Total, Bicarb, Carb | X | SM 5310B TOC | X |
| SM 5310B DOC | X | SM 5310B Solids, Total Dissolved | X |
| SM 4500 N Org B TKN | X | EPA 160.2 TSS | X |

| | |
|------------------------------|--------------------------------------|
| Relinquished by: (Signature) | Received by: (Signature/Affiliation) |
| Relinquished by: (Signature) | Received by: (Signature/Affiliation) |
| Relinquished by: (Signature) | Received by: (Signature/Affiliation) |

Date: 05/07/15 Time: 1030



From: (210) 877-2847
Philip C. Pearce, P.G.
SWCA
6200 UTSA Blvd
Ste 102
San Antonio, TX 78249

Origin ID: SATA



Ship Date: 06MAY15
ActWgt: 50.0 LB
CAD: 8347991/INET3610

Dims: 24 X 13 X 14 IN

Delivery Address Bar Code



SHIP TO: (714) 895-5494
Laboratory
Eurofins Calscience
7440 LINCOLN WAY

BILL SENDER

GARDEN GROVE, CA 92841

Ref # 27122.02.08
Invoice #
PO #
Dept #

1 of 24

THU - 07 MAY AA

STANDARD OVERNIGHT

TRK# 7735 3671 6556

0201

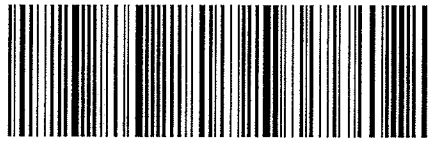
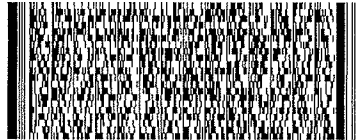
MASTER

A7 APVA

DSR

92841

CA-US

SNA

537J1ZSE2EE4B

24 Piece shipment

Tracking number

773536716556

(master)

773536716957

773536716990

773536717092

773536717221

773536717482

773536717574

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773536718250

773536718320

773536718353

773536718401

773536718456

773536718673

773536718982

773536719040

773536719084

773536719338

773536719371

773536719522

773536719566

SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.3°C); Temperature (w/o CF): 3.5 °C (w/ CF): 3.2 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☐ Yes ☒ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☒ No ☐ N/A

Sample container(s) intact and in good condition ☐ Yes ☒ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: 150413B)

Aqueous: ☐ VOA ☒ VOA ⁽³⁾ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ⁽²⁾ ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 subs ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 2 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 2.3 °C (w/ CF): 2.0 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Present but Not Intact☐ Not Present☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact☐ Present but Not Intact☒ Not Present☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples

Yes ☒ No ☐ N/A ☐

COC document(s) received complete

☐ Yes ☒ No ☐ N/A ☐☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished time

Sampler's name indicated on COC

☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC

☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition

☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested

☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested

☒ Yes ☐ No ☐ N/A

Samples received within holding time

☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container

☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace

☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation

☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAn ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 cubi ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄,

Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 3 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 3.1 °C (w/ CF): 2.8 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☐ Yes ☒ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAn ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 sub ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 4 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 4.0 °C (w/ CF): 3.7 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☐ Yes ☒ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAn ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 cubi ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 5 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 3.3 °C (w/ CF): 3.0 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☐ Yes ☒ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAn ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 cubi ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 6 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 2.7 °C (w/ CF): 2.4 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input checked="" type="checkbox"/> No relinquished date <input checked="" type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAn ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 Subi ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 7 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 2.4 °C (w/ CF): 2.1 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input checked="" type="checkbox"/> No relinquished date <input checked="" type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 cubi ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 8 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 2.4 °C (w/ CF): 2.1 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☐ Yes ☒ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAn ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 cubi ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 9 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 3.1 °C (w/ CF): 2.8 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input checked="" type="checkbox"/> No relinquished date <input checked="" type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 sub² ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300s = H₂SO₄, u = ultra-pure, z₂na = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 10 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 4.4 °C (w/ CF): 4.3 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☐ Yes ☒ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA⁽³⁾ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBz_{na} ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 Subi ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, z_{na} = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 11 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 3.7 °C (w/ CF): 3.4 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input checked="" type="checkbox"/> No relinquished date <input checked="" type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA⁽³⁾ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ⁽²⁾ ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 subⁱ ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 12 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 4.5 °C (w/ CF): 4.2 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Present but Not Intact☐ Not Present☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact☐ Present but Not Intact☒ Not Present☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples

Yes No N/A

COC document(s) received complete

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished time

Sampler's name indicated on COC

Sample container label(s) consistent with COC

Sample container(s) intact and in good condition

Proper containers for analyses requested

Sufficient volume/mass for analyses requested

Samples received within holding time

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen

Proper preservation chemical(s) noted on COC and/or sample container

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 cubi ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300s = H₂SO₄, u = ultra-pure, z₂na = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 13 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 4.0 °C (w/ CF): 3.7 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|--------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished time

| | | | |
|---------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---------------------------------------|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Aqueous samples for certain analyses received within 15-minute holding time

| | | | |
|--|--------------------------|-------------------------------------|--------------------------|
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
|--|--------------------------|-------------------------------------|--------------------------|

| | | | |
|--|-------------------------------------|--------------------------|--------------------------|
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|-------------------------------------|--------------------------|--------------------------|

Unpreserved aqueous sample(s) received for certain analyses

☒ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

| | | | |
|---|-------------------------------------|--------------------------|--------------------------|
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---|-------------------------------------|--------------------------|--------------------------|

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

| | | | |
|---|--------------------------|--------------------------|-------------------------------------|
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|---|--------------------------|--------------------------|-------------------------------------|

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 cubi ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, z₂na = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 14 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 2.9 °C (w/ CF): 2.6 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☐ Yes ☒ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAn ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 cubi ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 15 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 2.9 °C (w/ CF): 2.6 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Present but Not Intact☐ Not Present☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact☐ Present but Not Intact☒ Not Present☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☐ Yes ☒ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBz_{na} ☐ 250AGB ☐ 250CGB ☒ 250CGBs ² ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 Cubi ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® (____) ☐ TerraCores® (____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300s = H₂SO₄, u = ultra-pure, z_{na} = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 16 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 3.5 °C (w/ CF): 3.2 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input checked="" type="checkbox"/> No relinquished date <input checked="" type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAn ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 cubi ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 17 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 4.8 °C (w/ CF): 4.5 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☐ Yes ☒ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA⁽³⁾ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs⁽²⁾ ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 cubi ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 18 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.3°C); Temperature (w/o CF): 4.4 °C (w/ CF): 4.1 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☐ Yes ☒ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAn ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 cu bi ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 19 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 2.8 °C (w/ CF): 2.5 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Present but Not Intact☐ Not Present☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact☐ Present but Not Intact☒ Not Present☐ N/A

Checked by: 300

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input checked="" type="checkbox"/> No relinquished date <input checked="" type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBz_{na} ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 sub¹ ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® (____) ☐ TerraCores® (____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300s = H₂SO₄, u = ultra-pure, z_{na} = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 20 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.3°C); Temperature (w/o CF): 4.0 °C (w/ CF): 3.7 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☐ Yes ☒ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA⁽³⁾ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 cubi ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 2 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 2.7 °C (w/ CF): 2.4 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input checked="" type="checkbox"/> No relinquished date <input checked="" type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB
☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs
☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 Subi ☐ _____ ☐ _____ ☐ _____
Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____
Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, z₂na = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 22 OF 24CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 3.2 °C (w/ CF): 2.9 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ FilterChecked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Present but Not Intact☐ Not Present☐ N/AChecked by: 300Sample(s) ☐ Present and Intact☐ Present but Not Intact☒ Not Present☐ N/AChecked by: 300

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input checked="" type="checkbox"/> No relinquished date <input checked="" type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAn ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 Cubi ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOHReviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 23 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 3.6 °C (w/ CF): 3.3 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 300

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☐ Yes ☒ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☒ No relinquished date ☒ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA⁽³⁾ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBz^{na} ☐ 250AGB ☐ 250CGB ☒ 250CGBs ⁽²⁾ ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 ^{sub} ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, z^{na} = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE RECEIPT CHECKLIST

COOLER 24 OF 24

CLIENT: SWCA Env'l Consultants

DATE: 05 / 07 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.3°C); Temperature (w/o CF): 4.6 °C (w/ CF): 4.3 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 300

CUSTODY SEAL:

Cooler ☒ Present and Intact

☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 300

Sample(s) ☐ Present and Intact

☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 300

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input checked="" type="checkbox"/> No relinquished date <input checked="" type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAn ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznn ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5Cubi ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 300

s = H₂SO₄, u = ultra-pure, znn = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 972

SAMPLE ANOMALY REPORT

DATE: 05 / 07 / 2015

SAMPLES, CONTAINERS, AND LABELS:

- ☐ Sample(s) NOT RECEIVED but listed on COC
- ☐ Sample(s) received but NOT LISTED on COC
- ☐ Holding time expired (list client or ECI sample ID and analysis)
- ☐ Insufficient sample amount for requested analysis (list analysis)
- ☐ Improper container(s) used (list analysis)
- ☐ Improper preservative used (list analysis)
- ☐ No preservative noted on COC or label (list analysis and notify lab)
- ☐ Sample container(s) not labeled
- ☐ Client sample label(s) illegible (list container type and analysis)
- ☒ Client sample label(s) do not match COC (comment)
 - ☐ Project information
 - ☐ Client sample ID
 - ☐ Sampling date and/or time
 - ☒ Number of container(s)
 - ☒ Requested analysis
- ☐ Sample container(s) compromised (comment)
 - ☐ Broken
 - ☐ Water present in sample container
- ☐ Air sample container(s) compromised (comment)
 - ☐ Flat
 - ☐ Very low in volume
 - ☐ Leaking (not transferred; duplicate bag submitted)
 - ☐ Leaking (transferred into ECI Tedlar™ bags*)
 - ☐ Leaking (transferred into client's Tedlar™ bags*)

* Transferred at client's request.

MISCELLANEOUS: (Describe)

Limited Volume

HEADSPACE:

(Containers with bubble > 6 mm or ¼ inch for volatile organic or dissolved gas analysis)

| ECI Sample ID | ECI Container ID | Total Number** | ECI Sample ID | ECI Container ID | Total Number** |
|---------------|------------------|----------------|---------------|------------------|----------------|
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

Comments

(-24) FDHSM 23 D Trail received only 6 of 9 containers
Not received: 2 of 3 vials (8260)
1 Liter Amber Glass (8270)

1 Liter Amber Glass w/ H₂SO₄ (TKA) received not full (~400ml)

(-15) TB06 received vial for 8260 (requested 8270).

Comments

Comments: _____

** Record the total number of containers (i.e., vials or bottles) for the affected sample.

Reported by: 972

Reviewed by: 300

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS210 Lead
Sample Number: AA97930
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 11:46
Receipt Date/Time: 10/23/15 15:18

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97930-A | E. coli | √ | SM 9223B-2004 | 77000 | MPN/100 mL | | 1 | 44747 | 10/23/15 | 16:49 | PAL/CKK |
| AA97930-A | E. Coli Holding Time - IDEXX Colilert | | NA | 5.05 | hours | | 0.00 | 44746 | 10/23/15 | 16:49 | PAL/CKK |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44747

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

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Sample Location: HCS240 Lead
Sample Number: AA97931
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 12:00
Receipt Date/Time: 10/23/15 15:18

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97931-A | E. coli | √ | SM 9223B-2004 | 20000 | MPN/100 mL | | 1 | 44747 | 10/23/15 | 16:49 | PAL/CKK |
| AA97931-A | E. Coli Holding Time - IDEXX Colilert | | NA | 4.82 | hours | | 0.00 | 44746 | 10/23/15 | 16:49 | PAL/CKK |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44747

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

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6200 UTSA Blvd Ste 102
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Sample Location: HCS250 Lead
Sample Number: AA97932
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 12:03
Receipt Date/Time: 10/23/15 15:18

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97932-A | E. coli | √ | SM 9223B-2004 | 25000 | MPN/100 mL | | 1 | 44747 | 10/23/15 | 16:49 | PAL/CKK |
| AA97932-A | E. Coli Holding Time - IDEXX Colilert | | NA | 4.77 | hours | | 0.00 | 44746 | 10/23/15 | 16:49 | PAL/CKK |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44747

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

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Sample Location: HCS260 Lead
Sample Number: AA97933
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 12:24
Receipt Date/Time: 10/23/15 15:18

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|---------------------------------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97933-A | E. coli | √ | SM 9223B-2004 | 17000 | MPN/100 mL | | 1 | 44747 | 10/23/15 | 16:49 | PAL/CKK |
| AA97933-A | E. Coli Holding Time - IDEXX Colilert | | NA | 4.42 | hours | | 0.00 | 44746 | 10/23/15 | 16:49 | PAL/CKK |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44747

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

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Sample Location: HCS270 Lead
Sample Number: AA97934
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 12:20
Receipt Date/Time: 10/23/15 15:18

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

No sample and/or analysis comment(s)

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97934-A | E. coli | √ | SM 9223B-2004 | 24000 | MPN/100 mL | | 1 | 44747 | 10/23/15 | 16:49 | PAL/CKK |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44747

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

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Sample Location: HSM210 LEAD
Sample Number: AA97943
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 15:25
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97943-A | E. coli | √ | SM 9223B-2004 | 41 | MPN/100 mL | H | 1 | 44773 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44773

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

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Sample Location: HSM230 LEAD
Sample Number: AA97944
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 15:41
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed. Limited air space.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97944-A | E. coli | √ | SM 9223B-2004 | 87000 | MPN/100 mL | H | 1 | 44773 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44773

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

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Sample Location: HSM231 LEAD
Sample Number: AA97945
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 16:02
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed. Minimal sample volume.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97945-A | E. coli | √ | SM 9223B-2004 | 11000 | MPN/100 mL | H | 1 | 44773 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44773

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

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Sample Location: HSM240 LEAD
Sample Number: AA97946
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 15:30
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97946-A | E. coli | √ | SM 9223B-2004 | 28000 | MPN/100 mL | H | 1 | 44773 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44773

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

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Sample Location: HSM250 LEAD
Sample Number: AA97947
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 16:21
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97947-A | E. coli | √ | SM 9223B-2004 | 9200 | MPN/100 mL | H | 1 | 44773 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44773

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM260 LEAD
Sample Number: AA97948
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 15:58
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97948-A | E. coli | √ | SM 9223B-2004 | 7300 | MPN/100 mL | H | 1 | 44773 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44773

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

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Sample Location: HCS240 PEAK
Sample Number: AA97949
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 14:36
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97949-A | E. coli | √ | SM 9223B-2004 | 2000 | MPN/100 mL | H | 1 | 44773 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44773

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

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Sample Location: HCS250 PEAK
Sample Number: AA97950
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 14:59
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97950-A | E. coli | √ | SM 9223B-2004 | 17000 | MPN/100 mL | H | 1 | 44773 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44773

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



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Page 1 of 2

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Sample Location: HCS260 PEAK
Sample Number: AA97951
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 15:01
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed. Minmial sample volume.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97951-A | E. coli | √ | SM 9223B-2004 | 20000 | MPN/100 mL | H | 1 | 44773 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44773

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

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Sample Location: HCS270 PEAK
Sample Number: AA97952
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 15:16
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97952-A | E. coli | √ | SM 9223B-2004 | 20000 | MPN/100 mL | H | 1 | 44775 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44775

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

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Sample Location: HSM260 PEAK
Sample Number: AA97953
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 17:25
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97953-A | E. coli | √ | SM 9223B-2004 | 24000 | MPN/100 mL | H | 1 | 44775 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44775

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



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Page 1 of 2

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Sample Location: HSM270 PEAK
Sample Number: AA97954
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 17:45
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97954-A | E. coli | √ | SM 9223B-2004 | 8700 | MPN/100 mL | H | 1 | 44775 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44775

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

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Sample Location: HSM210 TRAIL
Sample Number: AA97955
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 20:35
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97955-A | E. coli | √ | SM 9223B-2004 | 450 | MPN/100 mL | H | 1 | 44775 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44775

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

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Sample Location: FDHSM210 TRAIL
Sample Number: AA97956
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 20:35
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97956-A | E. coli | √ | SM 9223B-2004 | 180 | MPN/100 mL | H | 1 | 44775 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44775

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



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Page 1 of 2

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Sample Location: HSM230 TRAIL
Sample Number: AA97957
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 21:01
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97957-A | E. coli | √ | SM 9223B-2004 | 73000 | MPN/100 mL | H | 1 | 44775 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44775

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



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Sample Location: FDHSM230 TRAIL
Sample Number: AA97958
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 21:01
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed. Minimal sample volume.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97958-A | E. coli | √ | SM 9223B-2004 | 65000 | MPN/100 mL | H | 1 | 44775 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
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QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44775

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HCS270 TRAIL
Sample Number: AA97959
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 17:34
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97959-A | E. coli | √ | SM 9223B-2004 | 20000 | MPN/100 mL | H | 1 | 44775 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44775

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

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Sample Location: FDHCS270 TRAIL
Sample Number: AA97960
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 17:34
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97960-A | E. coli | √ | SM 9223B-2004 | 17000 | MPN/100 mL | H | 1 | 44775 | 10/26/15 | 15:42 | HH/ST |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44775

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

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Sample Location: HSM231 TRAIL
Sample Number: AA97961
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 21:21
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97961-A | E. coli | √ | SM 9223B-2004 | 1700 | MPN/100 mL | H | 1 | 44777 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44777

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
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Sample Location: FDHSM231 TRAIL
Sample Number: AA97962
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 21:21
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97962-A | E. coli | √ | SM 9223B-2004 | 1700 | MPN/100 mL | H | 1 | 44777 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44777

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

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Sample Location: HSM240 TRAIL
Sample Number: AA97963
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 20:51
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97963-A | E. coli | √ | SM 9223B-2004 | 750 | MPN/100 mL | H | 1 | 44777 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44777

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

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Sample Location: HSM250 TRAIL
Sample Number: AA97964
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 20:06
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97964-A | E. coli | √ | SM 9223B-2004 | 1900 | MPN/100 mL | H | 1 | 44777 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44777

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

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Sample Location: HSM260 TRAIL
Sample Number: AA97965
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 21:35
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97965-A | E. coli | √ | SM 9223B-2004 | 4600 | MPN/100 mL | H | 1 | 44777 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44777

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
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Sample Location: HSM270 TRAIL
Sample Number: AA97966
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 21:55
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97966-A | E. coli | √ | SM 9223B-2004 | 4100 | MPN/100 mL | H | 1 | 44777 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44777

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

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Sample Location: HCS210 TRAIL
Sample Number: AA97967
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 16:14
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97967-A | E. coli | √ | SM 9223B-2004 | 69000 | MPN/100 mL | H | 1 | 44777 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44777

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

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Sample Location: HCS240 TRAIL
Sample Number: AA97968
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 16:29
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97968-A | E. coli | √ | SM 9223B-2004 | 2900 | MPN/100 mL | H | 1 | 44777 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44777

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

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Sample Location: HCS250 TRAIL
Sample Number: AA97969
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 17:09
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97969-A | E. coli | √ | SM 9223B-2004 | 16000 | MPN/100 mL | H | 1 | 44777 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44777

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HCS260 TRAIL
Sample Number: AA97970
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 16:50
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97970-A | E. coli | √ | SM 9223B-2004 | 20000 | MPN/100 mL | H | 1 | 44779 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44779

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: FDHCS260 TRAIL
Sample Number: AA97971
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 16:50
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97971-A | E. coli | √ | SM 9223B-2004 | 24000 | MPN/100 mL | H | 1 | 44779 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44779

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HSM270 LEAD
Sample Number: AA97972
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 16:15
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97972-A | E. coli | √ | SM 9223B-2004 | 14000 | MPN/100 mL | H | 1 | 44779 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44779

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HSM210 PEAK
Sample Number: AA97973
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 17:02
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97973-A | E. coli | √ | SM 9223B-2004 | 41 | MPN/100 mL | H | 1 | 44779 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44779

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HSM230 PEAK
Sample Number: AA97974
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 17:16
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97974-A | E. coli | √ | SM 9223B-2004 | 11000 | MPN/100 mL | H | 1 | 44779 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44779

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HSM231 PEAK
Sample Number: AA97975
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 17:06
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97975-A | E. coli | √ | SM 9223B-2004 | 2200 | MPN/100 mL | H | 1 | 44779 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44779

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HSM240 PEAK
Sample Number: AA97976
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 17:34
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative .

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97976-A | E. coli | √ | SM 9223B-2004 | 1000 | MPN/100 mL | H | 1 | 44779 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44779

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory

ANALYTICAL REPORT



November 03, 2015

Page 1 of 2

Client:

Phil Pearce
6200 UTSA Blvd Ste 102
San Antonio, TX 78249

Fax #: NA

This analytical report is intended exclusively for the individual or entity to which it is addressed. Recipient is not authorized to print or copy this report, except in full without written approval of the laboratory. If you have received this report in error, please notify the San Antonio River Authority.

Sample Location: HSM250 PEAK
Sample Number: AA97977
Sample Matrix: Non Potable Water

Collection Date/Time: 10/23/15 17:05
Receipt Date/Time: 10/26/15 08:32

CASE NARRATIVE

This report provides results related only to the referenced sample ID numbers. All samples were received in acceptable condition unless otherwise noted. For questions regarding this report, please contact Greg Mateo, Laboratory Supervisor, at (210) 302-3290.

Analysis identified with a "√" complies with NELAP requirements unless otherwise specified in the case narrative.

Sample Comments: Exceeded hold time, per customer proceed.

ANALYTICAL RESULTS

| | <u>Analyses</u> | <u>NELAC</u> | <u>Analysis Method</u> | <u>Result</u> | <u>Units</u> | <u>Qualifier</u> | <u>Reporting</u> | <u>QC Batch</u> | <u>Analysis</u> | | <u>Analyst</u> |
|-----------|-----------------|--------------|------------------------|---------------|--------------|------------------|------------------|-----------------|-----------------|-------------|----------------|
| | | | | | | | <u>Limit</u> | <u>Number</u> | <u>Date</u> | <u>Time</u> | |
| AA97977-A | E. coli | √ | SM 9223B-2004 | 6500 | MPN/100 mL | H | 1 | 44779 | 10/26/15 | 15:55 | HH |

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable

Environmental Sciences Department Laboratory
ANALYTICAL REPORT



November 03, 2015

Page 2 of 2

QC ANALYTICAL RESULTS

QC Batch Name: E_COLI_QUANTITRAY-44779

Acceptance Criteria

QC Analyte Name

Initial Blank for E. coli

Result

Absent

Units

Qualifier

Lower

Target

Absent

Upper



Patricia M. Carvajal
Quality Assurance Supervisor

11/3/2015

Date

A - Outside upper acceptance criteria
D - Outside lower acceptance criteria
T - Microbiological Controls were unacceptable

H - Hold Time for preparation or analysis exceeded
J - Analyte detected outside quantitation limit

* - See Case Narrative
--- - Not Applicable



WORK ORDER NUMBER: 15-10-1856

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: SWCA Environmental Consultants

Client Project Name: EAA 27122

Attention: Philip Pearce
6200 UTSA Blvd.
Suite 102
San Antonio, TX 78249-1618

A handwritten signature in black ink, appearing to read "Don Burley".

Approved for release on 11/23/2015 by:
Don Burley
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

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Work Order Number: 15-10-1856

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Work Order Narrative

Work Order: 15-10-1856

Page 1 of 1

Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 10/24/15. They were assigned to Work Order 15-10-1856.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.



Calscience

Sample Summary

| | |
|--|------------------------------------|
| Client: SWCA Environmental Consultants | Work Order: 15-10-1856 |
| 6200 UTSA Blvd., Suite 102 | Project Name: EAA 27122 |
| San Antonio, TX 78249-1618 | PO Number: |
| | Date/Time Received: 10/24/15 11:40 |
| | Number of Containers: 75 |
| Attn: Philip Pearce | |

| Sample Identification | Lab Number | Collection Date and Time | Number of Containers | Matrix |
|-----------------------|--------------|--------------------------|----------------------|---------|
| HCS 210 Lead | 15-10-1856-1 | 10/23/15 11:46 | 9 | Aqueous |
| HCS 240 Lead | 15-10-1856-2 | 10/23/15 12:00 | 9 | Aqueous |
| HCS 250 Lead | 15-10-1856-3 | 10/23/15 12:03 | 9 | Aqueous |
| HCS 260 Lead | 15-10-1856-4 | 10/23/15 12:24 | 9 | Aqueous |
| HCS 270 Lead | 15-10-1856-5 | 10/23/15 12:20 | 11 | Aqueous |
| TB14 | 15-10-1856-6 | 10/23/15 19:30 | 1 | Aqueous |
| HCS 210 Peak | 15-10-1856-7 | 10/23/15 14:19 | 9 | Aqueous |
| HCS 240 Peak | 15-10-1856-8 | 10/23/15 14:36 | 9 | Aqueous |
| HCS 250 Peak | 15-10-1856-9 | 10/23/15 14:59 | 9 | Aqueous |

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Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-10-1856-1-I | 10/23/15 11:46 | Aqueous | IC 15 | N/A | 10/24/15 18:28 | 151024L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.030 | 0.10 | 0.027 | 1.00 | J |
| Chloride | 1.8 | 1.0 | 0.52 | 1.00 | |
| Bromide | ND | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 0.44 | 0.10 | 0.053 | 1.00 | |
| Sulfate | 2.7 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|--------------|----------------|-------------------|---------|-------|-----|-------------------|-----------|
| HCS 240 Lead | 15-10-1856-2-I | 10/23/15 12:00 | Aqueous | IC 15 | N/A | 10/24/15 18:46 | 151024L01 |
|--------------|----------------|-------------------|---------|-------|-----|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.16 | 0.10 | 0.027 | 1.00 | |
| Chloride | 11 | 1.0 | 0.52 | 1.00 | |
| Bromide | ND | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | |
| Sulfate | 17 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|--------------|----------------|-------------------|---------|-------|-----|-------------------|-----------|
| HCS 250 Lead | 15-10-1856-3-I | 10/23/15 12:03 | Aqueous | IC 15 | N/A | 10/24/15 19:05 | 151024L01 |
|--------------|----------------|-------------------|---------|-------|-----|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.15 | 0.10 | 0.027 | 1.00 | |
| Chloride | 13 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.086 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.3 | 0.10 | 0.053 | 1.00 | |
| Sulfate | 19 | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-10-1856-4-I | 10/23/15 12:24 | Aqueous | IC 15 | N/A | 10/24/15 20:04 | 151024L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.19 | 0.10 | 0.027 | 1.00 | |
| Chloride | 17 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.11 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.7 | 0.10 | 0.053 | 1.00 | |
| Sulfate | 30 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|-------|-----|----------------|-----------|
| HCS 270 Lead | 15-10-1856-5-I | 10/23/15 12:20 | Aqueous | IC 15 | N/A | 10/24/15 20:22 | 151024L01 |
|--------------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.16 | 0.10 | 0.027 | 1.00 | |
| Chloride | 13 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.086 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.4 | 0.10 | 0.053 | 1.00 | |
| Sulfate | 22 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|-------|-----|----------------|-----------|
| HCS 210 Peak | 15-10-1856-7-I | 10/23/15 14:19 | Aqueous | IC 15 | N/A | 10/24/15 20:41 | 151024L01 |
|--------------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.081 | 0.10 | 0.027 | 1.00 | J |
| Chloride | 4.8 | 1.0 | 0.52 | 1.00 | |
| Bromide | ND | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 0.61 | 0.10 | 0.053 | 1.00 | |
| Sulfate | 7.5 | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-10-1856-8-I | 10/23/15 14:36 | Aqueous | IC 15 | N/A | 10/24/15 20:59 | 151024L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.18 | 0.10 | 0.027 | 1.00 | |
| Chloride | 17 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.11 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.7 | 0.10 | 0.053 | 1.00 | |
| Sulfate | 26 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|--------------|----------------|-------------------|---------|-------|-----|-------------------|-----------|
| HCS 250 Peak | 15-10-1856-9-I | 10/23/15 14:59 | Aqueous | IC 15 | N/A | 10/24/15 21:39 | 151024L01 |
|--------------|----------------|-------------------|---------|-------|-----|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.19 | 0.10 | 0.027 | 1.00 | |
| Chloride | 16 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.095 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.6 | 0.10 | 0.053 | 1.00 | |
| Sulfate | 24 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|--------------|-----------------|-----|---------|-------|-----|-------------------|-----------|
| Method Blank | 099-12-906-6192 | N/A | Aqueous | IC 15 | N/A | 10/24/15 15:40 | 151024L01 |
|--------------|-----------------|-----|---------|-------|-----|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | ND | 0.10 | 0.027 | 1.00 | |
| Chloride | ND | 1.0 | 0.52 | 1.00 | |
| Bromide | ND | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | ND | 0.10 | 0.053 | 1.00 | |
| Sulfate | ND | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-10-1856-1-D | 10/23/15 11:46 | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 20:44 | 151026LA4 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 14.1 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 1.28 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 3.29 | 0.500 | 0.103 | 1.00 | |
| Sodium | 1.32 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.0449 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 1.07 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|----------|----------|----------------|-----------|
| HCS 240 Lead | 15-10-1856-2-D | 10/23/15 12:00 | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 20:50 | 151026LA4 |
|--------------|----------------|----------------|---------|----------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 54.5 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 11.7 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.77 | 0.500 | 0.103 | 1.00 | |
| Sodium | 8.60 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.438 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.28 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|----------|----------|----------------|-----------|
| HCS 250 Lead | 15-10-1856-3-D | 10/23/15 12:03 | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 20:53 | 151026LA4 |
|--------------|----------------|----------------|---------|----------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 62.5 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 13.1 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.30 | 0.500 | 0.103 | 1.00 | |
| Sodium | 9.56 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.497 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.66 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-10-1856-4-D | 10/23/15 12:24 | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 20:55 | 151026LA4 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 77.6 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.0 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.86 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.657 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.92 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|----------|----------|----------------|-----------|
| HCS 270 Lead | 15-10-1856-5-D | 10/23/15 12:20 | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 20:57 | 151026LA4 |
|--------------|----------------|----------------|---------|----------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 61.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 13.2 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.50 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.512 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.82 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|----------|----------|----------------|-----------|
| HCS 210 Peak | 15-10-1856-7-D | 10/23/15 14:19 | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 21:00 | 151026LA4 |
|--------------|----------------|----------------|---------|----------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 26.7 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 4.61 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 3.18 | 0.500 | 0.103 | 1.00 | |
| Sodium | 3.50 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.174 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 2.17 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HCS 240 Peak | 15-10-1856-8-D | 10/23/15 14:36 | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 21:02 | 151026LA4 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 80.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 16.9 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.65 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.9 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.653 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 6.04 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HCS 250 Peak | 15-10-1856-9-D | 10/23/15 14:59 | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 21:04 | 151026LA4 |
|---------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 73.1 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 15.3 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.07 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.596 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.51 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|------------|----------------|-----------------|-----------------|-----------------------|------------------|
| Method Blank | 099-15-683-1457 | N/A | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 19:06 | 151026LA4 |
|---------------------|------------------------|------------|----------------|-----------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | ND | 0.100 | 0.0118 | 1.00 | |
| Magnesium | ND | 0.100 | 0.00336 | 1.00 | |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-10-1856-1-D | 10/23/15 11:46 | Aqueous | ICP/MS 03 | 10/26/15 | 11/02/15 11:25 | 151026LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000140 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.00942 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000720 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00183 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000270 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00110 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0125 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0258 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0665 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Lead | 15-10-1856-2-D | 10/23/15 12:00 | Aqueous | ICP/MS 03 | 10/26/15 | 10/30/15 05:25 | 151026LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0381 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000449 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.000975 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | 0.000105 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00106 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000329 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0228 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00911 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0704 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00399 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-10-1856-3-D | 10/23/15 12:03 | Aqueous | ICP/MS 03 | 10/26/15 | 10/30/15 05:29 | 151026LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|-----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | 0.000391 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | 0.0435 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000420 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.000787 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | 0.0000992 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00102 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000383 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0117 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00464 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0467 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00522 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-10-1856-4-D | 10/23/15 12:24 | Aqueous | ICP/MS 03 | 10/26/15 | 10/30/15 05:32 | 151026LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | 0.00763 | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0553 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000570 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00120 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00147 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000413 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0112 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00530 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0534 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00319 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

Page 5 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-10-1856-5-D | 10/23/15 12:20 | Aqueous | ICP/MS 03 | 10/26/15 | 10/30/15 05:36 | 151026LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0433 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00159 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000103 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00218 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000327 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0111 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0101 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0586 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00264 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Peak | 15-10-1856-7-D | 10/23/15 14:19 | Aqueous | ICP/MS 03 | 10/26/15 | 10/30/15 05:39 | 151026LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0190 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000455 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00181 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000142 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.000836 | 0.00100 | 0.000132 | 1.00 | J |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0171 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0114 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0525 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00116 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-10-1856-8-D | 10/23/15 14:36 | Aqueous | ICP/MS 03 | 10/26/15 | 10/30/15 05:43 | 151026LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0549 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000261 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00150 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000496 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0110 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00453 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0519 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00105 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Peak | 15-10-1856-9-D | 10/23/15 14:59 | Aqueous | ICP/MS 03 | 10/26/15 | 10/30/15 05:46 | 151026LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | 0.000446 | 0.00100 | 0.000386 | 1.00 | J |
| Barium | 0.0506 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000656 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00144 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000394 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00915 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0105 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0511 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00257 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-950 | N/A | Aqueous | ICP/MS 03 | 10/26/15 | 10/30/15 03:39 | 151026LA1F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

Page 1 of 2

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-10-1856-1-D | 10/23/15 11:46 | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 21:04 | 151029LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 240 Lead | 15-10-1856-2-D | 10/23/15 12:00 | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 21:06 | 151029LA3F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 250 Lead | 15-10-1856-3-D | 10/23/15 12:03 | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 21:08 | 151029LA3F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 260 Lead | 15-10-1856-4-D | 10/23/15 12:24 | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 21:11 | 151029LA3F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 270 Lead | 15-10-1856-5-D | 10/23/15 12:20 | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 21:13 | 151029LA3F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 210 Peak | 15-10-1856-7-D | 10/23/15 14:19 | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 21:15 | 151029LA3F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-10-1856-8-D | 10/23/15 14:36 | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 21:17 | 151029LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 250 Peak | 15-10-1856-9-D | 10/23/15 14:59 | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 21:24 | 151029LA3F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|-----|---------|------------|----------|----------------|------------|
| Method Blank | 099-15-763-647 | N/A | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 20:22 | 151029LA3F |
|--------------|----------------|-----|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-10-1856-1-I | 10/23/15 11:46 | Aqueous | GC 44 | 10/29/15 | 10/31/15 03:55 | 151029L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 87 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 88 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

Page 2 of 9

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Lead | 15-10-1856-2-I | 10/23/15 12:00 | Aqueous | GC 44 | 10/29/15 | 10/31/15 04:09 | 151029L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 109 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 87 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-10-1856-3-I | 10/23/15 12:03 | Aqueous | GC 44 | 10/29/15 | 10/31/15 04:23 | 151029L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 86 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-10-1856-4-I | 10/23/15 12:24 | Aqueous | GC 44 | 10/29/15 | 10/31/15 04:37 | 151029L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 86 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 85 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-10-1856-5-I | 10/23/15 12:20 | Aqueous | GC 44 | 10/29/15 | 10/31/15 04:52 | 151029L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 82 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 83 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Peak | 15-10-1856-7-I | 10/23/15 14:19 | Aqueous | GC 44 | 10/29/15 | 10/31/15 05:06 | 151029L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 83 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 79 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-10-1856-8-I | 10/23/15 14:36 | Aqueous | GC 44 | 10/29/15 | 10/31/15 05:20 | 151029L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 84 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 84 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Peak | 15-10-1856-9-I | 10/23/15 14:59 | Aqueous | GC 44 | 10/29/15 | 10/31/15 05:35 | 151029L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 87 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 91 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-850 | N/A | Aqueous | GC 44 | 10/29/15 | 10/29/15 21:20 | 151029L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 88 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 92 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

Page 1 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-10-1856-1-I | 10/23/15 11:46 | Aqueous | GC 66 | 10/29/15 | 10/30/15 17:16 | 151029L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 75 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 81 | 50-135 | |

| | | | | | | | |
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
| HCS 240 Lead | 15-10-1856-2-I | 10/23/15 12:00 | Aqueous | GC 66 | 10/29/15 | 10/30/15 17:34 | 151029L03 |
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 74 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 80 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-10-1856-3-I | 10/23/15 12:03 | Aqueous | GC 66 | 10/29/15 | 10/30/15 18:58 | 151029L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 76 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 79 | 50-135 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|
| HCS 260 Lead | 15-10-1856-4-I | 10/23/15 12:24 | Aqueous | GC 66 | 10/29/15 | 10/30/15 19:16 | 151029L03 |
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 78 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 80 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-10-1856-5-I | 10/23/15 12:20 | Aqueous | GC 66 | 10/29/15 | 10/30/15 19:34 | 151029L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 73 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 76 | 50-135 | |

| HCS 210 Peak | 15-10-1856-7-I | 10/23/15 14:19 | Aqueous | GC 66 | 10/29/15 | 10/30/15 19:52 | 151029L03 |
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 79 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 81 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

Page 4 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-10-1856-8-I | 10/23/15 14:36 | Aqueous | GC 66 | 10/29/15 | 10/30/15 20:09 | 151029L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 78 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 80 | 50-135 | |

| | | | | | | | |
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
| HCS 250 Peak | 15-10-1856-9-I | 10/23/15 14:59 | Aqueous | GC 66 | 10/29/15 | 10/30/15 20:27 | 151029L03 |
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 80 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 81 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-533-1100 | N/A | Aqueous | GC 66 | 10/29/15 | 10/30/15 11:16 | 151029L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 95 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 81 | 50-135 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-10-1856-1-I | 10/23/15 11:46 | Aqueous | GC 26 | 10/27/15 | 10/29/15 18:05 | 151027L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 77 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Lead | 15-10-1856-2-I | 10/23/15 12:00 | Aqueous | GC 26 | 10/27/15 | 10/29/15 18:49 | 151027L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 66 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-10-1856-3-I | 10/23/15 12:03 | Aqueous | GC 26 | 10/27/15 | 10/29/15 19:33 | 151027L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 83 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-10-1856-4-I | 10/23/15 12:24 | Aqueous | GC 26 | 10/27/15 | 10/29/15 20:17 | 151027L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 75 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-10-1856-5-I | 10/23/15 12:20 | Aqueous | GC 26 | 10/27/15 | 10/29/15 21:01 | 151027L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 74 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Peak | 15-10-1856-7-I | 10/23/15 14:19 | Aqueous | GC 26 | 10/27/15 | 10/29/15 21:45 | 151027L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 74 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-10-1856-8-I | 10/23/15 14:36 | Aqueous | GC 26 | 10/27/15 | 10/29/15 22:29 | 151027L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 74 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

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Date Received: 10/24/15
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Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Peak | 15-10-1856-9-I | 10/23/15 14:59 | Aqueous | GC 26 | 10/27/15 | 10/29/15 23:13 | 151027L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 73 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-119 | N/A | Aqueous | GC 26 | 10/27/15 | 10/29/15 17:20 | 151027L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 79 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-10-1856-1-I | 10/23/15 11:46 | Aqueous | GC 40 | 10/26/15 | 11/03/15 21:25 | 151026L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 95 | 0-123 | |

| | | | | | | | |
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
| HCS 240 Lead | 15-10-1856-2-I | 10/23/15 12:00 | Aqueous | GC 40 | 10/26/15 | 11/03/15 21:48 | 151026L01 |
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 104 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-10-1856-3-I | 10/23/15 12:03 | Aqueous | GC 40 | 10/26/15 | 11/03/15 22:11 | 151026L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 104 | 0-123 | |

| | | | | | | | |
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
| HCS 260 Lead | 15-10-1856-4-I | 10/23/15 12:24 | Aqueous | GC 40 | 10/26/15 | 11/03/15 22:34 | 151026L01 |
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 90 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
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Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-10-1856-5-I | 10/23/15 12:20 | Aqueous | GC 40 | 10/26/15 | 11/03/15 22:57 | 151026L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 87 | 0-123 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|
| HCS 210 Peak | 15-10-1856-7-I | 10/23/15 14:19 | Aqueous | GC 40 | 10/26/15 | 11/03/15 23:20 | 151026L01 |
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 98 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
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Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-10-1856-8-I | 10/23/15 14:36 | Aqueous | GC 40 | 10/26/15 | 11/03/15 23:43 | 151026L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 81 | 0-123 | |

| HCS 250 Peak | 15-10-1856-9-I | 10/23/15 14:59 | Aqueous | GC 40 | 10/26/15 | 11/04/15 00:06 | 151026L01 |
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 83 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-034-665 | N/A | Aqueous | GC 40 | 10/26/15 | 11/03/15 18:21 | 151026L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPPE | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 55 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

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Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-10-1856-1-H | 10/23/15 11:46 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 13:30 | 151030L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 74 | 33-120 | |
| 2-Fluorophenol | 77 | 24-120 | |
| Nitrobenzene-d5 | 67 | 38-120 | |
| p-Terphenyl-d14 | 67 | 41-137 | |
| Phenol-d6 | 74 | 16-120 | |
| 2,4,6-Tribromophenol | 86 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Lead | 15-10-1856-2-H | 10/23/15 12:00 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 13:49 | 151030L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 66 | 33-120 | |
| 2-Fluorophenol | 68 | 24-120 | |
| Nitrobenzene-d5 | 60 | 38-120 | |
| p-Terphenyl-d14 | 61 | 41-137 | |
| Phenol-d6 | 66 | 16-120 | |
| 2,4,6-Tribromophenol | 74 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-10-1856-3-H | 10/23/15 12:03 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 14:08 | 151030L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 61 | 33-120 | |
| 2-Fluorophenol | 62 | 24-120 | |
| Nitrobenzene-d5 | 55 | 38-120 | |
| p-Terphenyl-d14 | 58 | 41-137 | |
| Phenol-d6 | 59 | 16-120 | |
| 2,4,6-Tribromophenol | 68 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-10-1856-4-H | 10/23/15 12:24 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 14:27 | 151030L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 65 | 33-120 | |
| 2-Fluorophenol | 67 | 24-120 | |
| Nitrobenzene-d5 | 58 | 38-120 | |
| p-Terphenyl-d14 | 60 | 41-137 | |
| Phenol-d6 | 63 | 16-120 | |
| 2,4,6-Tribromophenol | 71 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-10-1856-5-G | 10/23/15 12:20 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 14:45 | 151030L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 65 | 33-120 | |
| 2-Fluorophenol | 64 | 24-120 | |
| Nitrobenzene-d5 | 56 | 38-120 | |
| p-Terphenyl-d14 | 64 | 41-137 | |
| Phenol-d6 | 61 | 16-120 | |
| 2,4,6-Tribromophenol | 75 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Peak | 15-10-1856-7-H | 10/23/15 14:19 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 15:04 | 151030L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 70 | 33-120 | |
| 2-Fluorophenol | 71 | 24-120 | |
| Nitrobenzene-d5 | 63 | 38-120 | |
| p-Terphenyl-d14 | 63 | 41-137 | |
| Phenol-d6 | 67 | 16-120 | |
| 2,4,6-Tribromophenol | 79 | 27-159 | |



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-10-1856-8-H | 10/23/15 14:36 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 15:23 | 151030L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 66 | 33-120 | |
| 2-Fluorophenol | 70 | 24-120 | |
| Nitrobenzene-d5 | 61 | 38-120 | |
| p-Terphenyl-d14 | 64 | 41-137 | |
| Phenol-d6 | 66 | 16-120 | |
| 2,4,6-Tribromophenol | 74 | 27-159 | |



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Peak | 15-10-1856-9-H | 10/23/15 14:59 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 15:42 | 151030L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 70 | 33-120 | |
| 2-Fluorophenol | 71 | 24-120 | |
| Nitrobenzene-d5 | 62 | 38-120 | |
| p-Terphenyl-d14 | 65 | 41-137 | |
| Phenol-d6 | 68 | 16-120 | |
| 2,4,6-Tribromophenol | 78 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-4131 | N/A | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 16:38 | 151030L11 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 71 | 33-120 | |
| 2-Fluorophenol | 73 | 24-120 | |
| Nitrobenzene-d5 | 64 | 38-120 | |
| p-Terphenyl-d14 | 66 | 41-137 | |
| Phenol-d6 | 68 | 16-120 | |
| 2,4,6-Tribromophenol | 78 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Lead | 15-10-1856-1-A | 10/23/15 11:46 | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 17:08 | 151028L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 103 | 78-126 | |
| 1,2-Dichloroethane-d4 | 106 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Lead | 15-10-1856-2-A | 10/23/15 12:00 | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 17:35 | 151028L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 105 | 78-126 | |
| 1,2-Dichloroethane-d4 | 105 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Lead | 15-10-1856-3-A | 10/23/15 12:03 | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 18:02 | 151028L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 105 | 78-126 | |
| 1,2-Dichloroethane-d4 | 106 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Lead | 15-10-1856-4-A | 10/23/15 12:24 | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 18:29 | 151028L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 95 | 80-120 | |
| Dibromofluoromethane | 105 | 78-126 | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | |
| Toluene-d8 | 96 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Lead | 15-10-1856-5-A | 10/23/15 12:20 | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 13:59 | 151028L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 10 of 20

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 95 | 80-120 | |
| Dibromofluoromethane | 103 | 78-126 | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | |
| Toluene-d8 | 98 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| TB14 | 15-10-1856-6-A | 10/23/15 19:30 | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 13:32 | 151028L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|----------|----------------|------------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| 1,4-Bromofluorobenzene | 95 | 80-120 | | | |
| Dibromofluoromethane | 101 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | | | |
| Toluene-d8 | 97 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Peak | 15-10-1856-7-A | 10/23/15 14:19 | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 18:56 | 151028L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | 12 | 20 | 10 | 1.00 | J |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 93 | 80-120 | |
| Dibromofluoromethane | 104 | 78-126 | |
| 1,2-Dichloroethane-d4 | 106 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Peak | 15-10-1856-8-A | 10/23/15 14:36 | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 19:23 | 151028L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 105 | 78-126 | |
| 1,2-Dichloroethane-d4 | 107 | 75-135 | |
| Toluene-d8 | 96 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Peak | 15-10-1856-9-A | 10/23/15 14:59 | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 19:51 | 151028L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 104 | 78-126 | |
| 1,2-Dichloroethane-d4 | 106 | 75-135 | |
| Toluene-d8 | 95 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-18614 | N/A | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 13:06 | 151028L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 94 | 80-120 | |
| Dibromofluoromethane | 100 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

10/24/15

Work Order:

15-10-1856

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 210 Lead | 15-10-1856-1 | 10/23/15 11:46 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.16 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 43.0 | 1.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 43.0 | 1.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 158 | 1.00 | 0.870 | 1.00 | | mg/L | 10/29/15 | 10/29/15 | SM 2540 C |
| Solids, Total Suspended (24) | 206 | 1.00 | 0.829 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.27 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 2.4 | 0.50 | 0.28 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 8.8 | 0.50 | 0.24 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 8.9 | 0.50 | 0.24 | 1.00 | | mg/L | 10/24/15 | 10/24/15 | SM 5310 B |

| | | | |
|---------------------|---------------------|-----------------------|----------------|
| HCS 240 Lead | 15-10-1856-2 | 10/23/15 12:00 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.12 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 164 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 164 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 195 | 1.00 | 0.870 | 1.00 | | mg/L | 10/29/15 | 10/29/15 | SM 2540 C |
| Solids, Total Suspended (24) | 58 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.53 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.1 | 0.50 | 0.28 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.9 | 0.50 | 0.24 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 8.4 | 0.50 | 0.24 | 1.00 | | mg/L | 10/24/15 | 10/24/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

10/24/15

Work Order:

15-10-1856

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 250 Lead | 15-10-1856-3 | 10/23/15 12:03 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.091 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 195 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 195 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 300 | 1.00 | 0.870 | 1.00 | | mg/L | 10/29/15 | 10/29/15 | SM 2540 C |
| Solids, Total Suspended (24) | 162 | 1.00 | 0.829 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.56 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.1 | 0.50 | 0.28 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.4 | 0.50 | 0.24 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 6.1 | 0.50 | 0.24 | 1.00 | | mg/L | 10/24/15 | 10/24/15 | SM 5310 B |

| | | | |
|---------------------|---------------------|-----------------------|----------------|
| HCS 260 Lead | 15-10-1856-4 | 10/23/15 12:24 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.032 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 227 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 227 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 490 | 1.00 | 0.870 | 1.00 | | mg/L | 10/29/15 | 10/29/15 | SM 2540 C |
| Solids, Total Suspended (24) | 30 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.68 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.84 | 0.50 | 0.28 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 0.73 | 0.50 | 0.24 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 0.79 | 0.50 | 0.24 | 1.00 | | mg/L | 10/24/15 | 10/24/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

10/24/15

Work Order:

15-10-1856

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 270 Lead | 15-10-1856-5 | 10/23/15 12:20 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.036 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 183 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 183 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 360 | 1.00 | 0.870 | 1.00 | | mg/L | 10/29/15 | 10/29/15 | SM 2540 C |
| Solids, Total Suspended (24) | 9.9 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.71 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.70 | 0.50 | 0.28 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 10/24/15 | 10/24/15 | SM 5310 B |

| | | | |
|---------------------|---------------------|-----------------------|----------------|
| HCS 210 Peak | 15-10-1856-7 | 10/23/15 14:19 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.36 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 82.0 | 1.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 82.0 | 1.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 265 | 1.00 | 0.870 | 1.00 | | mg/L | 10/29/15 | 10/29/15 | SM 2540 C |
| Solids, Total Suspended (24) | 17 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.69 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.3 | 0.50 | 0.28 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 12 | 0.50 | 0.24 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 10 | 0.50 | 0.24 | 1.00 | | mg/L | 10/24/15 | 10/24/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

10/24/15

Work Order:

15-10-1856

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 240 Peak | 15-10-1856-8 | 10/23/15 14:36 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 166 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 166 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 360 | 1.00 | 0.870 | 1.00 | | mg/L | 10/29/15 | 10/29/15 | SM 2540 C |
| Solids, Total Suspended (24) | 6.1 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.68 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 10/24/15 | 10/24/15 | SM 5310 B |

| | | | |
|---------------------|---------------------|-----------------------|----------------|
| HCS 250 Peak | 15-10-1856-9 | 10/23/15 14:59 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.021 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 232 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 232 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 480 | 1.00 | 0.870 | 1.00 | | mg/L | 10/29/15 | 10/29/15 | SM 2540 C |
| Solids, Total Suspended (24) | 3.1 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.65 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/24/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.3 | 0.50 | 0.28 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 10/24/15 | 10/24/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

10/24/15

Work Order:

15-10-1856

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|-------------------|---------------------|----------------|
| Method Blank | | N/A | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Results</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> | <u>Units</u> | <u>Date Prepared</u> | <u>Date Analyzed</u> | <u>Method</u> |
|--|----------------|-----------|------------|-----------|-------------------|--------------|----------------------|----------------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/09/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.87 | 1.00 | | mg/L | 10/29/15 | 10/29/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 10/24/15 | 10/24/15 | SM 5310 B |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HCS 250 Peak | Sample | Aqueous | IC 15 | N/A | 10/24/15 21:39 | 151024S1 |
| HCS 250 Peak | Matrix Spike | Aqueous | IC 15 | N/A | 10/24/15 22:53 | 151024S1 |
| HCS 250 Peak | Matrix Spike Duplicate | Aqueous | IC 15 | N/A | 10/24/15 23:11 | 151024S1 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.1862 | 2.500 | 2.538 | 94 | 2.539 | 94 | 80-120 | 0 | 0-20 | |
| Chloride | 15.90 | 50.00 | 67.55 | 103 | 67.57 | 103 | 80-120 | 0 | 0-20 | |
| Bromide | ND | 5.000 | 5.036 | 101 | 5.042 | 101 | 80-120 | 0 | 0-20 | |
| Nitrate (as N) | 1.598 | 5.000 | 6.715 | 102 | 6.700 | 102 | 80-120 | 0 | 0-20 | |
| Sulfate | 23.92 | 50.00 | 77.79 | 108 | 77.64 | 107 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-10-1947-7 | Sample | Aqueous | ACA 1 | N/A | 11/09/15 17:16 | 151109S02 |
| 15-10-1947-7 | Matrix Spike | Aqueous | ACA 1 | N/A | 11/09/15 17:16 | 151109S02 |
| 15-10-1947-7 | Matrix Spike Duplicate | Aqueous | ACA 1 | N/A | 11/09/15 17:16 | 151109S02 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | ND | 0.2000 | 0.2580 | 129 | 0.2465 | 123 | 90-110 | 5 | 0-25 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HCS 240 Peak | Sample | Aqueous | TOC 8 | 11/05/15 | 11/05/15 23:04 | F1105TOCS1 |
| HCS 240 Peak | Matrix Spike | Aqueous | TOC 8 | 11/05/15 | 11/05/15 23:04 | F1105TOCS1 |
| HCS 240 Peak | Matrix Spike Duplicate | Aqueous | TOC 8 | 11/05/15 | 11/05/15 23:04 | F1105TOCS1 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Carbon, Total Organic | ND | 10.00 | 17.80 | 178 | 19.00 | 190 | 31-145 | 7 | 0-23 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-10-1826-2 | Sample | Aqueous | TOC 8 | 10/24/15 | 10/24/15 17:36 | F1024DOCS2 |
| 15-10-1826-2 | Matrix Spike | Aqueous | TOC 8 | 10/24/15 | 10/24/15 17:36 | F1024DOCS2 |
| 15-10-1826-2 | Matrix Spike Duplicate | Aqueous | TOC 8 | 10/24/15 | 10/24/15 17:36 | F1024DOCS2 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|---------------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Carbon, Dissolved Organic | 11.60 | 50.00 | 65.50 | 108 | 67.00 | 111 | 31-145 | 2 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HCS 270 Lead | Sample | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 20:57 | 151026SA4A |
| HCS 270 Lead | Matrix Spike | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 20:40 | 151026SA4A |
| HCS 270 Lead | Matrix Spike Duplicate | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 20:42 | 151026SA4A |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Calcium | 61.15 | 0.5000 | 61.77 | 4X | 62.24 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 13.25 | 0.5000 | 13.68 | 4X | 13.59 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 2.500 | 5.000 | 7.555 | 101 | 7.526 | 101 | 83-131 | 0 | 0-7 | |
| Sodium | 11.07 | 5.000 | 16.21 | 103 | 16.35 | 105 | 73-127 | 1 | 0-9 | |
| Strontium | 0.5118 | 0.5000 | 1.034 | 104 | 1.035 | 105 | 81-123 | 0 | 0-6 | |
| Silicon | 4.818 | 0.5000 | 5.262 | 4X | 5.241 | 4X | 24-180 | 4X | 0-15 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3020A Total
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| 15-10-1826-4 | Sample | Aqueous | ICP/MS 03 | 10/26/15 | 10/30/15 04:47 | 151026SA1 | | | | |
| 15-10-1826-4 | Matrix Spike | Aqueous | ICP/MS 03 | 10/26/15 | 10/30/15 03:46 | 151026SA1 | | | | |
| 15-10-1826-4 | Matrix Spike Duplicate | Aqueous | ICP/MS 03 | 10/26/15 | 10/30/15 03:50 | 151026SA1 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.1031 | 103 | 0.1007 | 101 | 85-133 | 2 | 0-11 | |
| Arsenic | 0.005215 | 0.1000 | 0.1006 | 95 | 0.09807 | 93 | 73-127 | 3 | 0-11 | |
| Barium | 0.1099 | 0.1000 | 0.2068 | 97 | 0.1907 | 81 | 74-128 | 8 | 0-10 | |
| Beryllium | ND | 0.1000 | 0.07817 | 78 | 0.07682 | 77 | 56-122 | 2 | 0-11 | |
| Cadmium | ND | 0.1000 | 0.09293 | 93 | 0.09018 | 90 | 84-114 | 3 | 0-8 | |
| Chromium | 0.01641 | 0.1000 | 0.1265 | 110 | 0.1230 | 107 | 73-133 | 3 | 0-11 | |
| Copper | 0.003928 | 0.1000 | 0.1020 | 98 | 0.09896 | 95 | 72-108 | 3 | 0-10 | |
| Lead | ND | 0.1000 | 0.1161 | 116 | 0.1134 | 113 | 79-121 | 2 | 0-10 | |
| Nickel | 0.01517 | 0.1000 | 0.1236 | 108 | 0.1112 | 96 | 68-122 | 11 | 0-10 | 4 |
| Selenium | ND | 0.1000 | 0.07781 | 78 | 0.07515 | 75 | 59-125 | 3 | 0-12 | |
| Silver | ND | 0.05000 | 0.04182 | 84 | 0.03878 | 78 | 68-128 | 8 | 0-14 | |
| Thallium | ND | 0.1000 | 0.1133 | 113 | 0.1114 | 111 | 73-121 | 2 | 0-11 | |
| Zinc | 0.005715 | 0.1000 | 0.08870 | 83 | 0.08288 | 77 | 43-145 | 7 | 0-39 | |
| Aluminum | 0.2697 | 0.1000 | 0.3112 | 42 | 0.3196 | 50 | 47-161 | 3 | 0-24 | 3 |
| Iron | 1.181 | 5.100 | 6.660 | 107 | 6.644 | 107 | 27-201 | 0 | 0-24 | |
| Manganese | 0.8608 | 0.1000 | 0.8803 | 4X | 0.8503 | 4X | 72-126 | 4X | 0-42 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-10-1850-3 | Sample | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 20:31 | 151029SA3 |
| 15-10-1850-3 | Matrix Spike | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 20:33 | 151029SA3 |
| 15-10-1850-3 | Matrix Spike Duplicate | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 20:35 | 151029SA3 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | ND | 0.01000 | 0.009809 | 98 | 0.009738 | 97 | 55-133 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HCS 270 Lead | Sample | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 13:59 | 151028S005 | | | | |
| HCS 270 Lead | Matrix Spike | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 11:18 | 151028S005 | | | | |
| HCS 270 Lead | Matrix Spike Duplicate | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 11:45 | 151028S005 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 47.46 | 95 | 47.55 | 95 | 74-122 | 0 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 46.89 | 94 | 47.36 | 95 | 60-144 | 1 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 50.73 | 101 | 50.94 | 102 | 73-120 | 0 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 47.99 | 96 | 48.97 | 98 | 80-122 | 2 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 50.69 | 101 | 51.67 | 103 | 70-120 | 2 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 45.75 | 92 | 46.85 | 94 | 64-142 | 2 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 46.08 | 92 | 47.10 | 94 | 52-136 | 2 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 52.07 | 104 | 51.88 | 104 | 77-125 | 0 | 0-24 | |
| Toluene | ND | 50.00 | 47.89 | 96 | 48.45 | 97 | 72-126 | 1 | 0-23 | |
| Trichloroethene | ND | 50.00 | 47.70 | 95 | 48.27 | 97 | 74-128 | 1 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 49.34 | 99 | 50.73 | 101 | 67-133 | 3 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 104.3 | 104 | 104.0 | 104 | 63-129 | 0 | 0-25 | |
| o-Xylene | ND | 50.00 | 51.83 | 104 | 52.01 | 104 | 62-128 | 0 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 43.00 | 86 | 44.85 | 90 | 68-134 | 4 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3020A Total
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|--------------|-------------|------------|----------------|----------------|-----------------------|
| 15-10-1826-4 | Sample | Aqueous | ICP/MS 03 | 10/26/15 00:00 | 10/30/15 04:47 | 151026SA1 |
| 15-10-1826-4 | PDS | Aqueous | ICP/MS 03 | 10/26/15 00:00 | 10/30/15 04:36 | 151026SA1 |
| Parameter | Sample Conc. | Spike Added | PDS Conc. | PDS %Rec. | %Rec. CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.09345 | 93 | 75-125 | |
| Arsenic | 0.005215 | 0.1000 | 0.09921 | 94 | 75-125 | |
| Barium | 0.1099 | 0.1000 | 0.2055 | 96 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.08161 | 82 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.08386 | 84 | 75-125 | |
| Chromium | 0.01641 | 0.1000 | 0.1158 | 99 | 75-125 | |
| Copper | 0.003928 | 0.1000 | 0.1010 | 97 | 75-125 | |
| Lead | ND | 0.1000 | 0.1075 | 107 | 75-125 | |
| Nickel | 0.01517 | 0.1000 | 0.1144 | 99 | 75-125 | |
| Selenium | ND | 0.1000 | 0.07543 | 75 | 75-125 | |
| Silver | ND | 0.05000 | 0.03921 | 78 | 75-125 | |
| Thallium | ND | 0.1000 | 0.1051 | 105 | 75-125 | |
| Zinc | 0.005715 | 0.1000 | 0.08520 | 79 | 75-125 | |
| Aluminum | 0.2697 | 0.1000 | 0.3689 | 99 | 75-125 | |
| Iron | 1.181 | 5.100 | 6.733 | 109 | 75-125 | |
| Manganese | 0.8608 | 0.1000 | 0.9460 | 4X | 75-125 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-10-2300-2 | Sample | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:09 | F1105ALKD1 |
| 15-10-2300-2 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:09 | F1105ALKD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|---|---------------------|------------------|------------|---------------|-------------------|
| Alkalinity, Total (as CaCO ₃) | 378.0 | 374.0 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-10-2300-2 | Sample | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:09 | F1105HCOD1 |
| 15-10-2300-2 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:09 | F1105HCOD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Bicarbonate (as CaCO ₃) | 378.0 | 374.0 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-10-2300-2 | Sample | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:09 | F1105CO3D1 |
| 15-10-2300-2 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:09 | F1105CO3D1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-10-1826-4 | Sample | Aqueous | SC 2 | 10/29/15 00:00 | 10/29/15 16:00 | F1029TDSD1 |
| 15-10-1826-4 | Sample Duplicate | Aqueous | SC 2 | 10/29/15 00:00 | 10/29/15 16:00 | F1029TDSD1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Dissolved | 2100 | 2155 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-10-1748-2 | Sample | Aqueous | N/A | 10/30/15 00:00 | 10/30/15 19:30 | F1030TSSD2 |
| 15-10-1748-2 | Sample Duplicate | Aqueous | N/A | 10/30/15 00:00 | 10/30/15 19:30 | F1030TSSD2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Suspended | 396.0 | 416.0 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------------------|------------------|---------------|----------------|------------------------|
| HCS 210 Lead | Sample | Aqueous | PH 1 | N/A | 10/24/15 15:40 | F1024PHD1 |
| HCS 210 Lead | Sample Duplicate | Aqueous | PH 1 | N/A | 10/24/15 15:40 | F1024PHD1 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| pH | | 7.270 | 7.290 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 4500 N Org B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-10-2174-3 | Sample | Aqueous | BUR05 | 10/30/15 00:00 | 10/30/15 16:21 | F1030TKND2 |
| 15-10-2174-3 | Sample Duplicate | Aqueous | BUR05 | 10/30/15 00:00 | 10/30/15 16:21 | F1030TKND2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Total Kjeldahl Nitrogen | 54.46 | 54.04 | 1 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|----------------|--------------|---------------|-----------------------|------------------|
| 099-12-906-6192 | LCS | Aqueous | IC 15 | N/A | 10/24/15 15:59 | 151024L01 |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Fluoride | 2.500 | 2.530 | 101 | 90-110 | |
| Chloride | 50.00 | 48.71 | 97 | 90-110 | |
| Bromide | 5.000 | 4.899 | 98 | 90-110 | |
| Nitrate (as N) | 5.000 | 4.865 | 97 | 90-110 | |
| Sulfate | 50.00 | 48.42 | 97 | 90-110 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------------|------------|---------------|-------------------|
| 099-12-739-131 | LCS | Aqueous | ACA 1 | N/A | 11/09/15 17:16 | 151109L02 | | | |
| 099-12-739-131 | LCSD | Aqueous | ACA 1 | N/A | 11/09/15 17:16 | 151109L02 | | | |
| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Phosphorus, Total | 0.2000 | 0.1915 | 96 | 0.1917 | 96 | 90-110 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-859-848 | LCS | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:09 | F1105ALKB1 | | | |
| 099-15-859-848 | LCSD | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:09 | F1105ALKB1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 100.0 | 99.00 | 99 | 99.00 | 99 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4818 | LCS | Aqueous | SC 2 | 10/29/15 | 10/29/15 16:00 | F1029TDSB1 | | | |
| 099-12-180-4818 | LCSD | Aqueous | SC 2 | 10/29/15 | 10/29/15 16:00 | F1029TDSB1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 105.0 | 105 | 110.0 | 110 | 80-120 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------------|------------|---------------|-------------------|
| 099-09-010-7378 | LCS | Aqueous | N/A | 10/30/15 | 10/30/15 19:30 | F1030TSSL2 | | | |
| 099-09-010-7378 | LCSD | Aqueous | N/A | 10/30/15 | 10/30/15 19:30 | F1030TSSL2 | | | |
| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Solids, Total Suspended | 100.0 | 97.00 | 97 | 98.00 | 98 | 80-120 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 099-05-097-5777 | LCS | Aqueous | TOC 8 | 11/05/15 | 11/05/15 23:04 | F1105TOCL1 |
| 099-05-097-5777 | LCSD | Aqueous | TOC 8 | 11/05/15 | 11/05/15 23:04 | F1105TOCL1 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| Carbon, Total Organic | 10.00 | 9.960 | 100 | 10.40 | 104 | 80-120 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------------|------------|---------------|-------------------|
| 099-05-115-1465 | LCS | Aqueous | TOC 8 | 10/24/15 | 10/24/15 21:49 | F1024DOCL2 | | | |
| 099-05-115-1465 | LCSD | Aqueous | TOC 8 | 10/24/15 | 10/24/15 21:49 | F1024DOCL2 | | | |
| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Carbon, Dissolved Organic | 10.00 | 10.80 | 108 | 11.50 | 115 | 80-120 | 6 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|--------------------|------------------------|------------------|-----------------|-------------------|
| 099-15-683-1457 | LCS | Aqueous | ICP 7300 | 10/26/15 | 10/30/15 19:09 | 151026LA4 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Calcium | | 0.5000 | 0.4763 | 95 | 80-120 | |
| Magnesium | | 0.5000 | 0.5002 | 100 | 80-120 | |
| Potassium | | 5.000 | 4.716 | 94 | 80-120 | |
| Sodium | | 5.000 | 4.826 | 97 | 80-120 | |
| Strontium | | 0.5000 | 0.4954 | 99 | 80-120 | |
| Silicon | | 0.5000 | 0.4716 | 94 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-950 | LCS | Aqueous | ICP/MS 03 | 10/26/15 | 10/30/15 03:43 | 151026LA1F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.09937 | 99 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.09725 | 97 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.09679 | 97 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.1015 | 102 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.09851 | 99 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.09831 | 98 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.1022 | 102 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.09866 | 99 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.09921 | 99 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.09641 | 96 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.04342 | 87 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.09668 | 97 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.1002 | 100 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.09933 | 99 | 80-120 | 73-127 | |
| Iron | 5.100 | 5.186 | 102 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.09909 | 99 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|------------------|
| 099-15-763-647 | LCS | Aqueous | Mercury 04 | 10/29/15 | 10/29/15 20:24 | 151029LA3F |

| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
|-----------|-------------|-----------------|-----------|----------|------------|
| Mercury | 0.01000 | 0.009787 | 98 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|--------|-----|--------|------------|
| 099-12-529-850 | LCS | Aqueous | GC 44 | 10/29/15 | 10/30/15 13:00 | 151029L02 | | | | |
| 099-12-529-850 | LCSD | Aqueous | GC 44 | 10/29/15 | 10/30/15 13:14 | 151029L02 | | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Alpha-BHC | 0.5000 | 0.4916 | 98 | 0.5149 | 103 | 50-135 | 36-149 | 5 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.5020 | 100 | 0.5372 | 107 | 50-135 | 36-149 | 7 | 0-25 | |
| Beta-BHC | 0.5000 | 0.4658 | 93 | 0.5019 | 100 | 50-135 | 36-149 | 7 | 0-25 | |
| Heptachlor | 0.5000 | 0.4412 | 88 | 0.4614 | 92 | 50-135 | 36-149 | 4 | 0-25 | |
| Delta-BHC | 0.5000 | 0.4988 | 100 | 0.5145 | 103 | 50-135 | 36-149 | 3 | 0-25 | |
| Aldrin | 0.5000 | 0.3977 | 80 | 0.4112 | 82 | 50-135 | 36-149 | 3 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.4588 | 92 | 0.4810 | 96 | 50-135 | 36-149 | 5 | 0-25 | |
| Endosulfan I | 0.5000 | 0.4614 | 92 | 0.4850 | 97 | 50-135 | 36-149 | 5 | 0-25 | |
| Dieldrin | 0.5000 | 0.4796 | 96 | 0.5024 | 100 | 50-135 | 36-149 | 5 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.4696 | 94 | 0.4895 | 98 | 50-135 | 36-149 | 4 | 0-25 | |
| Endrin | 0.5000 | 0.5846 | 117 | 0.6066 | 121 | 50-135 | 36-149 | 4 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.4422 | 88 | 0.4500 | 90 | 50-135 | 36-149 | 2 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.4761 | 95 | 0.4979 | 100 | 50-135 | 36-149 | 4 | 0-25 | |
| Endosulfan II | 0.5000 | 0.4734 | 95 | 0.4940 | 99 | 50-135 | 36-149 | 4 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.4900 | 98 | 0.5154 | 103 | 50-135 | 36-149 | 5 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.4892 | 98 | 0.4970 | 99 | 50-135 | 36-149 | 2 | 0-25 | |
| Methoxychlor | 0.5000 | 0.5550 | 111 | 0.5500 | 110 | 50-135 | 36-149 | 1 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-533-1100 | LCS | Aqueous | GC 66 | 10/29/15 | 10/30/15 10:41 | 151029L03 | | | |
| 099-12-533-1100 | LCSD | Aqueous | GC 66 | 10/29/15 | 10/30/15 10:59 | 151029L03 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Aroclor-1016 | 2.000 | 1.821 | 91 | 1.965 | 98 | 50-135 | 8 | 0-25 | |
| Aroclor-1260 | 2.000 | 2.035 | 102 | 2.218 | 111 | 50-135 | 9 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

Page 13 of 16

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|--------|-----|--------|------------|
| 099-15-963-119 | LCS | Aqueous | GC 26 | 10/27/15 | 10/29/15 15:52 | 151027L14 | | | | |
| 099-15-963-119 | LCSD | Aqueous | GC 26 | 10/27/15 | 10/29/15 16:36 | 151027L14 | | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.03598 | 90 | 0.03538 | 88 | 30-130 | 13-147 | 2 | 0-30 | |
| Bolstar | 0.04000 | 0.04100 | 102 | 0.04076 | 102 | 30-130 | 13-147 | 1 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.03870 | 97 | 0.03957 | 99 | 30-130 | 13-147 | 2 | 0-30 | |
| Coumaphos | 0.04000 | 0.03569 | 89 | 0.03607 | 90 | 30-130 | 13-147 | 1 | 0-30 | |
| Diazinon | 0.04000 | 0.04221 | 106 | 0.04391 | 110 | 30-130 | 13-147 | 4 | 0-30 | |
| Disulfoton | 0.04000 | 0.04147 | 104 | 0.04173 | 104 | 30-130 | 13-147 | 1 | 0-30 | |
| Ethoprop | 0.04000 | 0.04319 | 108 | 0.04444 | 111 | 30-130 | 13-147 | 3 | 0-30 | |
| Fensulfothion | 0.04000 | 0.04216 | 105 | 0.04249 | 106 | 30-130 | 13-147 | 1 | 0-30 | |
| Fenthion | 0.04000 | 0.04163 | 104 | 0.04243 | 106 | 30-130 | 13-147 | 2 | 0-30 | |
| Merphos | 0.04000 | 0.04436 | 111 | 0.04425 | 111 | 30-130 | 13-147 | 0 | 0-30 | |
| Methyl Parathion | 0.04000 | 0.03863 | 97 | 0.03946 | 99 | 30-130 | 13-147 | 2 | 0-30 | |
| Phorate | 0.04000 | 0.04844 | 121 | 0.04948 | 124 | 30-130 | 13-147 | 2 | 0-30 | |
| Ronnel | 0.04000 | 0.03579 | 89 | 0.03690 | 92 | 30-130 | 13-147 | 3 | 0-30 | |
| Stirophos | 0.04000 | 0.03327 | 83 | 0.03405 | 85 | 30-130 | 13-147 | 2 | 0-30 | |
| Tokuthion | 0.04000 | 0.03606 | 90 | 0.03654 | 91 | 30-130 | 13-147 | 1 | 0-30 | |
| Trichloronate | 0.04000 | 0.03931 | 98 | 0.04024 | 101 | 30-130 | 13-147 | 2 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-034-665 | LCS | Aqueous | GC 40 | 10/26/15 | 11/03/15 17:34 | 151026L01 | | | |
| 095-01-034-665 | LCSD | Aqueous | GC 40 | 10/26/15 | 11/03/15 17:58 | 151026L01 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| 2,4-D | 20.00 | 9.110 | 46 | 10.28 | 51 | 30-130 | 12 | 0-30 | |
| 2,4,5-T | 2.000 | 1.305 | 65 | 1.410 | 70 | 30-130 | 8 | 0-30 | |
| 2,4-DB | 20.00 | 9.355 | 47 | 10.66 | 53 | 30-130 | 13 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-4131 | LCS | Aqueous | | GC/MS TT | 10/30/15 | 10/31/15 16:00 | 151030L11 | | | |
| 095-01-003-4131 | LCSD | Aqueous | | GC/MS TT | 10/30/15 | 10/31/15 16:19 | 151030L11 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 200.0 | 146.9 | 73 | 151.9 | 76 | 61-120 | 51-130 | 3 | 0-20 | |
| Acenaphthylene | 200.0 | 141.1 | 71 | 147.6 | 74 | 55-120 | 44-131 | 4 | 0-20 | |
| Butyl Benzyl Phthalate | 200.0 | 137.7 | 69 | 144.3 | 72 | 56-122 | 45-133 | 5 | 0-20 | |
| 4-Chloro-3-Methylphenol | 200.0 | 140.3 | 70 | 147.6 | 74 | 52-120 | 41-131 | 5 | 0-20 | |
| 2-Chlorophenol | 200.0 | 146.5 | 73 | 154.0 | 77 | 47-120 | 35-132 | 5 | 0-20 | |
| 1,4-Dichlorobenzene | 200.0 | 140.1 | 70 | 144.5 | 72 | 36-120 | 22-134 | 3 | 0-20 | |
| Dimethyl Phthalate | 200.0 | 138.4 | 69 | 144.9 | 72 | 60-120 | 50-130 | 5 | 0-20 | |
| 2,4-Dinitrotoluene | 200.0 | 147.7 | 74 | 157.5 | 79 | 61-121 | 51-131 | 6 | 0-20 | |
| Fluorene | 200.0 | 144.4 | 72 | 153.6 | 77 | 67-120 | 58-129 | 6 | 0-20 | |
| N-Nitroso-di-n-propylamine | 200.0 | 139.5 | 70 | 149.0 | 75 | 39-123 | 25-137 | 7 | 0-20 | |
| Naphthalene | 200.0 | 140.4 | 70 | 147.0 | 73 | 54-120 | 43-131 | 5 | 0-20 | |
| 4-Nitrophenol | 200.0 | 136.1 | 68 | 140.8 | 70 | 14-120 | 0-138 | 3 | 0-20 | |
| Pentachlorophenol | 200.0 | 144.6 | 72 | 146.4 | 73 | 31-127 | 15-143 | 1 | 0-20 | |
| Phenol | 200.0 | 146.4 | 73 | 153.6 | 77 | 17-120 | 0-137 | 5 | 0-20 | |
| Pyrene | 200.0 | 140.2 | 70 | 146.5 | 73 | 58-124 | 47-135 | 4 | 0-20 | |
| 1,2,4-Trichlorobenzene | 200.0 | 139.6 | 70 | 143.7 | 72 | 49-120 | 37-132 | 3 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/24/15
Work Order: 15-10-1856
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-18614 | LCS | Aqueous | GC/MS JJ | 10/28/15 | 10/28/15 10:44 | 151028L006 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 44.01 | 88 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 44.89 | 90 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 48.40 | 97 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 45.93 | 92 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 50.53 | 101 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 44.07 | 88 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 45.85 | 92 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 49.58 | 99 | 80-123 | 73-130 | |
| Toluene | 50.00 | 44.78 | 90 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 46.04 | 92 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 45.14 | 90 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 100.0 | 100 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 49.25 | 98 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 40.72 | 81 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Sample Analysis Summary Report

Work Order: 15-10-1856

Page 1 of 1

| <u>Method</u> | <u>Extraction</u> | <u>Chemist ID</u> | <u>Instrument</u> | <u>Analytical Location</u> |
|-----------------|-------------------|-------------------|-------------------|----------------------------|
| EPA 300.0 | N/A | 1027 | IC 15 | 1 |
| EPA 365.1 | N/A | 735 | ACA 1 | 1 |
| EPA 6010B | EPA 3005A Filt. | 935 | ICP 7300 | 1 |
| EPA 6020 | EPA 3005A Filt. | 598 | ICP/MS 03 | 1 |
| EPA 7470A | EPA 7470A Filt. | 915 | Mercury 04 | 1 |
| EPA 8081A | EPA 3510C | 669 | GC 44 | 1 |
| EPA 8082 | EPA 3510C | 669 | GC 66 | 1 |
| EPA 8141A | EPA 3510C | 960 | GC 26 | 1 |
| EPA 8151A | EPA 8151A | 669 | GC 40 | 1 |
| EPA 8260B | EPA 5030C | 486 | GC/MS JJ | 2 |
| EPA 8270C | EPA 3510C | 923 | GC/MS TT | 1 |
| SM 2320B | N/A | 688 | PH1/BUR03 | 1 |
| SM 2540 C | N/A | 689 | SC 2 | 1 |
| SM 2540 D | N/A | 689 | N/A | 1 |
| SM 4500 H+ B | N/A | 650 | PH 1 | 1 |
| SM 4500 N Org B | N/A | 685 | BUR05 | 1 |
| SM 5310 B | N/A | 735 | TOC 8 | 1 |

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Location 1: 7440 Lincoln Way, Garden Grove, CA 92841

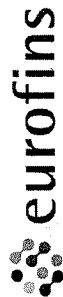
Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841

Glossary of Terms and Qualifiers

Work Order: 15-10-1856

Page 1 of 1

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|---|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| CI | See case narrative. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |
| | Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time. |
| | A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations. |



Calscience

7440 Lincoln Way, Garden Grove, CA 92641-1427 • (714) 895-5494
For courier service / sample drop off information, contact us26_sales@eurofins.com or call us.

LABORATORY CLIENT: SWCA Environmental Consultants

ADDRESS: 6200 UTSA Blvd. Suite 102

CITY: San Antonio

STATE: TX

ZIP: 78249-1618

TEL: 210.877.2847

E-MAIL: P.Pearce@swca.com

CLIENT PROJECT NAME / NUMBER: EAA 27122

PROJECT CONTACT: Philip Pearce

SAMPLER(S) (PRINT) Philip Pearce,
Jenna Cantwell, Jennifer
Moreland, + Brittany Rios

TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD"):

☐ SAME DAY ☐ 24 HR ☐ 48 HR ☒ 72 HR ☒ 10 Days (standard)

GLOBAL ID:

LOG CODE:

SPECIAL INSTRUCTIONS:

Samples for EPA 6010B, EPA 6020, and EPA 7470 Metals are field filtered
EPA 6020 Metals: Zn, Ti, Se, Sb, Pb, Ni, Cu, Cr, Cd, Be, Ba, As, Al, Ag, Fe, Mn
Please also analyze each sample for pH.

Please check box or fill in blank as needed.

REQUESTED ANALYSES

| LAB USE ONLY | SAMPLE ID | SAMPLING | | MATRIX | NO. OF CONT. | LOG CODE | | | EPA 6010B (Si, Ca, Mg, K, Na, Sr) | EPA 6020 (See Comments) | EPA 7470A Mercury | EPA 1694 (M) Caffeine | EPA 8081A Organochlorine Pesticides | EPA 8082 PCB Aroclors | EPA 8141A Organophosphorus Pest | EPA 8151A Chlorinated Herbicides | EPA 8270C Semi-Volatile Organics | EPA 8260B Volatile Organics | EPA 300 Anions (F, Cl, Br, NO ₃ , SO ₄ , CO ₃) | EPA 365.1 Total Phosphorus | SM 2320B Alkalinity: Total, Bicarb, Carb | SM 5310B TOC | SM 5310B DOC | SM 5310B Solids, Total Dissolved | SM 4500 N Org B TKN | EPA 160.2 TSS |
|--------------------|--------------|----------|-------|--------|--------------------|-------------|-----------|----------------|-----------------------------------|-------------------------|-------------------|-----------------------|-------------------------------------|-----------------------|---------------------------------|----------------------------------|----------------------------------|-----------------------------|--|----------------------------|--|--------------|--------------|----------------------------------|---------------------|---------------|
| | | DATE | TIME | | | Unpreserved | Preserved | Field Filtered | | | | | | | | | | | | | | | | | | |
| 1 | HCS 210 Lead | 10/23/15 | 11:46 | NPW | 9 | 2 | 7 | 1 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 2 | HCS 240 Lead | 10/23/15 | 12:00 | NPW | 9 | 2 | 7 | 1 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 3 | HCS 250 Lead | 10/23/15 | 12:03 | NPW | 9 | 2 | 7 | 1 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 4 | HCS 260 Lead | 10/23/15 | 12:24 | NPW | 9 | 2 | 7 | 1 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 5 | HCS 270 Lead | 10/23/15 | 12:20 | NPW | 11 | 2 | 7 | 1 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 6 | TB14 | 10/23/15 | 19:30 | NPW | 8 | 2 | 7 | 1 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 7 | HCS 210 PEAK | 10/23/15 | 14:19 | NPW | 9 | 2 | 7 | 1 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 8 | HCS 240 PEAK | 10/23/15 | 14:36 | NPW | 9 | 2 | 7 | 1 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |
| 9 | HCS 250 PEAK | 10/23/15 | 14:59 | NPW | 9 | 2 | 7 | 1 | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X | X |

Received by: (Signature/Affiliation)

Date: 10/23/15 Time: 19:35

Received by: (Signature/Affiliation)

Date: 10/24/15 Time: 17:40

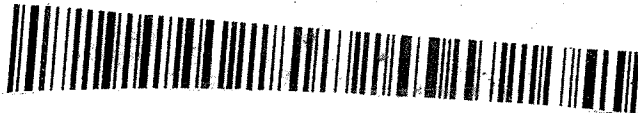
Received by: (Signature/Affiliation)

Date: Time:

1 of 8
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0215
MASTER

XO APVA

SATURDAY 12:00
PRIORITY OVERNIGHT
DS
9284
CA-US SN



MPS# 7815 7747 7992
0681
Mstr# 8079 4368 0270

XO APVA

SATURDAY 12:00
PRIORITY OVERNIGHT
DS
9284
CA-US SNA

0215

1856



3 of 8
MPS# 7815 7747 8006
0681
Mstr# 8079 4368 0270

XO APVA

SATURDAY 12:00
PRIORITY OVERNIGHT
DSF
92841
CA-US SNA

0215



4 of 8
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0681
Mstr# 8079 4368 0270

XO APVA

SATURDAY 12:00
PRIORITY OVERNIGHT
DSR
92841
CA-US SNA

0215



5 of 8
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Mstr# 8079 4368 0270

XO APVA

SATURDAY 12:00
PRIORITY OVERNIGHT
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92841
CA-US SNA

0215



6 of 8
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XO APVA

SATURDAY 12:00
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92841
CA-US SNA

0215

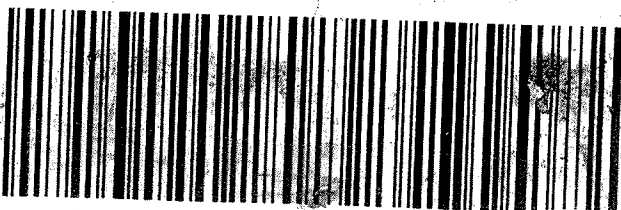


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XO APVA

SATURDAY 12:00
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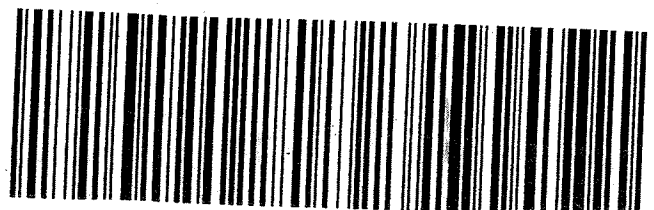


8 of 8
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0681
Mstr# 8079 4368 0270

XO APVA

SATURDAY 12:00
PRIORITY OVERNIGHT
DS
9284
CA-US SNA

0215



SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 8

CLIENT: SWCA Environmental Consultants

DATE: 10/24/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 3.5 °C (w/ CF): 3.1 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 862

CUSTODY SEAL:

Cooler ☒ Present and Intact

☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 862

Sample(s) ☐ Present and Intact

☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 778

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: 150831A)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na2} ☐ 100PJ ☐ 100PJ_{na2} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☒ 250PB_h ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s
☐ 500PB ☒ 1AGB ☐ 1AGB_{na2} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 125AGB_h ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 778

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 862

SAMPLE RECEIPT CHECKLIST

COOLER 2 OF 8

CLIENT: SWCA Environmental Consultants

DATE: 10/24/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 3.4 °C (w/ CF): 3.0 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 862

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 862

Sample(s) ☐ Present and Intact ☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 778

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|--------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBnp ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 25 g/Lube ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 778

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 862

SAMPLE RECEIPT CHECKLIST

COOLER 3 OF 8

CLIENT: SWCA Environmental Consultants

DATE: 10/24/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 3.7 °C (w/ CF): 3.3 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 862

CUSTODY SEAL:

Cooler ☒ Present and Intact

☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 862

Sample(s) ☐ Present and Intact

☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 778

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBnf ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 25galabe ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 778

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 862

SAMPLE RECEIPT CHECKLIST

COOLER 4 OF 8

CLIENT: SWCA Environmental Consultants

DATE: 10/24/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 3.5 °C (w/ CF): 3.1 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 862

CUSTODY SEAL:

Cooler ☒ Present and Intact

☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 862

Sample(s) ☐ Present and Intact

☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 228

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOAn₂ ☐ 100PJ ☐ 100PJn₂ ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PBz_{nn} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☒ 250PB_{nf} ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBn₂ ☒ 1AGBs ☐ 1PB ☐ 1PBn₂ ☒ 2.5 gal lake ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 228

s = H₂SO₄, u = ultra-pure, z_{nn}a = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 862

SAMPLE RECEIPT CHECKLIST

COOLER 5 OF 8

CLIENT: SWCA Environmental Consultants

DATE: 10/24/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.4°C); Temperature (w/o CF): 3.2 °C (w/ CF): 2.8 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 862

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 862

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 778

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☒ 250PB_{nf} ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5gal cube ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄,

Labeled/Checked by: 778

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 862

SAMPLE RECEIPT CHECKLIST

COOLER 6 OF 8

CLIENT: SWCA Environmental Consultants

DATE: 10/24/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 3.3 °C (w/ CF): 2.9 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 862

CUSTODY SEAL:

Cooler ☒ Present and Intact

☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 862

Sample(s) ☐ Present and Intact

☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 728

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☒ Yes ☐ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAn₂ ☐ 100PJ ☐ 100PJn₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBn² ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBn₂ ☒ 1AGBs ☐ 1PB ☐ 1PBn₂ ☒ 2.5 gal cube ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 728

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 862

SAMPLE RECEIPT CHECKLIST

COOLER 7 OF 8

CLIENT: SWCA Environmental Consultants

DATE: 10/24/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 3.4 °C (w/ CF): 3.0 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 862

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 862

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 728

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA³ ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☒ 250PBnf ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5 gal cube ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄,

Labeled/Checked by: 728

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 862

SAMPLE RECEIPT CHECKLIST

COOLER 8 OF 8

CLIENT: SWCA Environmental Consultants

DATE: 10/24/2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 3.3 °C (w/ CF): 2.9 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 862

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 862

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 778

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOAn₂ ☐ 100PJ ☐ 100PJn₂ ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PBz_{nn} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☒ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s
☐ 500PB ☒ 1AGB ☐ 1AGBn₂ ☒ 1AGB_s ☐ 1PB ☐ 1PBn₂ ☒ 2.5 gal chbr ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄,

Labeled/Checked by: 778

s = H₂SO₄, u = ultra-pure, z_{nn}a = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 862

SAMPLE ANOMALY REPORT

DATE: **10 / 24 / 2015**

SAMPLES, CONTAINERS, AND LABELS:

- ☐ Sample(s) NOT RECEIVED but listed on COC
- ☐ Sample(s) received but NOT LISTED on COC
- ☐ Holding time expired (list client or ECI sample ID and analysis)
- ☐ Insufficient sample amount for requested analysis (list analysis)
- ☐ Improper container(s) used (list analysis)
- ☐ Improper preservative used (list analysis)
- ☐ No preservative noted on COC or label (list analysis and notify lab)
- ☐ Sample container(s) not labeled
- ☐ Client sample label(s) illegible (list container type and analysis)
- ☐ Client sample label(s) do not match COC (comment)
 - ☐ Project information
 - ☐ Client sample ID
 - ☐ Sampling date and/or time
 - ☐ Number of container(s)
 - ☐ Requested analysis
- ☒ Sample container(s) compromised (comment)
 - ☒ Broken
 - ☐ Water present in sample container
- ☐ Air sample container(s) compromised (comment)
 - ☐ Flat
 - ☐ Very low in volume
 - ☐ Leaking (not transferred; duplicate bag submitted)
 - ☐ Leaking (transferred into ECI Tedlar™ bags*)
 - ☐ Leaking (transferred into client's Tedlar™ bags*)

* Transferred at client's request.

MISCELLANEOUS: (Describe)

HEADSPACE:

(Containers with bubble > 6 mm or ¼ inch for volatile organic or dissolved gas analysis)

| ECI Sample ID | ECI Container ID | Total Number** | ECI Sample ID | ECI Container ID | Total Number** |
|---------------|------------------|----------------|---------------|------------------|----------------|
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

Comments: _____

** Record the total number of containers (i.e., vials or bottles) for the affected sample.

Comments

57) 1 of 2 250ml glass / H₂SO₄
received broken

Comments

(Containers with bubble for other analysis)

| ECI Sample ID | ECI Container ID | Total Number** | Requested Analysis |
|---------------|------------------|----------------|--------------------|
| | | | |
| | | | |
| | | | |
| | | | |

Reported by: 228

Reviewed by: 862



Calscience

Subcontractor Analysis Report

Work Order: 15-10-1856Page 1 of 1

One or more samples in this work order have tests that were subcontracted. The subcontract report(s) follows.

For subcontracted tests, please reference the laboratory information noted below.

1. ALS - Columbia Analytical Services, Inc. - Kelso, WA CA ELAP 2286, NELAP WA100010
Method 1694 Caffeine


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Analytical Report for Service Request No:

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RE: 15-10-1856

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or o r re ere e e e e ee ed o r er ere e er **K1512169.**

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ALS Group USA, Corp. dba ALS Environmental

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Acronyms

Qualifiers

State Certifications, Accreditations, And Licenses

Case Narrative

Chain of Custody

Steroids and Endocrine Disrupting Compounds

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Acronyms

| | |
|------------|--|
| ASTM | American Society for Testing and Materials |
| A2LA | American Association for Laboratory Accreditation |
| CARB | California Air Resources Board |
| CAS Number | Chemical Abstract Service registry Number |
| CFC | Chlorofluorocarbon |
| CFU | Colony-Forming Unit |
| DEC | Department of Environmental Conservation |
| DEQ | Department of Environmental Quality |
| DHS | Department of Health Services |
| DOE | Department of Ecology |
| DOH | Department of Health |
| EPA | U. S. Environmental Protection Agency |
| ELAP | Environmental Laboratory Accreditation Program |
| GC | Gas Chromatography |
| GC/MS | Gas Chromatography/Mass Spectrometry |
| LOD | Limit of Detection |
| LOQ | Limit of Quantitation |
| LUFT | Leaking Underground Fuel Tank |
| M | Modified |
| MCL | Maximum Contaminant Level is the highest permissible concentration of a substance allowed in drinking water as established by the USEPA. |
| MDL | Method Detection Limit |
| MPN | Most Probable Number |
| MRL | Method Reporting Limit |
| NA | Not Applicable |
| NC | Not Calculated |
| NCASI | National Council of the Paper Industry for Air and Stream Improvement |
| ND | Not Detected |
| NIOSH | National Institute for Occupational Safety and Health |
| PQL | Practical Quantitation Limit |
| RCRA | Resource Conservation and Recovery Act |
| SIM | Selected Ion Monitoring |
| TPH | Total Petroleum Hydrocarbons |
| tr | Trace level is the concentration of an analyte that is less than the PQL but greater than or equal to the MDL. |

Inorganic Data Qualifiers

- * The result is an outlier. See case narrative.
- # The control limit criteria is not applicable. See case narrative.
- B The analyte was found in the associated method blank at a level that is significant relative to the sample result as defined by the DOD or NELAC standards.
- E The result is an estimate amount because the value exceeded the instrument calibration range.
- J The result is an estimated value.
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.
DOD-QSM 4.2 definition : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- i The MRL/MDL or LOQ/LOD is elevated due to a matrix interference.
- X See case narrative.
- Q See case narrative. One or more quality control criteria was outside the limits.
- H The holding time for this test is immediately following sample collection. The samples were analyzed as soon as possible after receipt by the laboratory.

Metals Data Qualifiers

- # The control limit criteria is not applicable. See case narrative.
- J The result is an estimated value.
- E The percent difference for the serial dilution was greater than 10%, indicating a possible matrix interference in the sample.
- M The duplicate injection precision was not met.
- N The Matrix Spike sample recovery is not within control limits. See case narrative.
- S The reported value was determined by the Method of Standard Additions (MSA).
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.
DOD-QSM 4.2 definition : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- W The post-digestion spike for furnace AA analysis is out of control limits, while sample absorbance is less than 50% of spike absorbance.
- i The MRL/MDL or LOQ/LOD is elevated due to a matrix interference.
- X See case narrative.
- + The correlation coefficient for the MSA is less than 0.995.
- Q See case narrative. One or more quality control criteria was outside the limits.

Organic Data Qualifiers

- * The result is an outlier. See case narrative.
- # The control limit criteria is not applicable. See case narrative.
- A A tentatively identified compound, a suspected aldol-condensation product.
- B The analyte was found in the associated method blank at a level that is significant relative to the sample result as defined by the DOD or NELAC standards.
- C The analyte was qualitatively confirmed using GC/MS techniques, pattern recognition, or by comparing to historical data.
- D The reported result is from a dilution.
- E The result is an estimated value.
- J The result is an estimated value.
- N The result is presumptive. The analyte was tentatively identified, but a confirmation analysis was not performed.
- P The GC or HPLC confirmation criteria was exceeded. The relative percent difference is greater than 40% between the two analytical results.
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.
DOD-QSM 4.2 definition : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- i The MRL/MDL or LOQ/LOD is elevated due to a chromatographic interference.
- X See case narrative.
- Q See case narrative. One or more quality control criteria was outside the limits.

Additional Petroleum Hydrocarbon Specific Qualifiers

- F The chromatographic fingerprint of the sample matches the elution pattern of the calibration standard.
- L The chromatographic fingerprint of the sample resembles a petroleum product, but the elution pattern indicates the presence of a greater amount of lighter molecular weight constituents than the calibration standard.
- H The chromatographic fingerprint of the sample resembles a petroleum product, but the elution pattern indicates the presence of a greater amount of heavier molecular weight constituents than the calibration standard.
- O The chromatographic fingerprint of the sample resembles an oil, but does not match the calibration standard.
- Y The chromatographic fingerprint of the sample resembles a petroleum product eluting in approximately the correct carbon range, but the elution pattern does not match the calibration standard.
- Z The chromatographic fingerprint does not resemble a petroleum product.

ALS Group USA Corp. dba ALS Environmental (ALS) - Kelso
State Certifications, Accreditations, and Licenses

| Agency | Web Site | Number |
|--------------------------|---|---------------|
| Alaska DEC UST | http://dec.alaska.gov/applications/eh/ehllabreports/USTLabs.aspx | UST-040 |
| Arizona DHS | http://www.azdhs.gov/lab/license/env.htm | AZ0339 |
| Arkansas - DEQ | http://www.adeq.state.ar.us/techsvs/labcert.htm | 88-0637 |
| California DHS (ELAP) | http://www.cdph.ca.gov/certlic/labs/Pages/ELAP.aspx | 2795 |
| DOD ELAP | http://www.denix.osd.mil/edqw/Accreditation/AccreditedLabs.cfm | L14-51 |
| Florida DOH | http://www.doh.state.fl.us/lab/EnvLabCert/WaterCert.htm | E87412 |
| Hawaii DOH | Not available | - |
| Idaho DHW | http://www.healthandwelfare.idaho.gov/Health/Labs/CertificationDrinkingWaterLabs/tabid/1833/Default.aspx | - |
| ISO 17025 | http://www.pjllabs.com/ | L14-50 |
| Louisiana DEQ | http://www.deq.louisiana.gov/portal/DIVISIONS/PublicParticipationandPermitSupport/LouisianaLaboratoryAccreditationProgram.aspx | 03016 |
| Maine DHS | Not available | WA01276 |
| Michigan DEQ | http://www.michigan.gov/deq/0,1607,7-135-3307_4131_4156---,00.html | 9949 |
| Minnesota DOH | http://www.health.state.mn.us/accreditation | 053-999-457 |
| Montana DPHHS | http://www.dphhs.mt.gov/publichealth/ | CERT0047 |
| Nevada DEP | http://ndep.nv.gov/bsdwlabservice.htm | WA01276 |
| New Jersey DEP | http://www.nj.gov/dep/oqa/ | WA005 |
| North Carolina DWQ | http://www.dwqlab.org/ | 605 |
| Oklahoma DEQ | http://www.deq.state.ok.us/CSDnew/labcert.htm | 9801 |
| Oregon – DEQ (NELAP) | http://public.health.oregon.gov/LaboratoryServices/EnvironmentalLaboratoryAccreditation/Pages/index.aspx | WA100010 |
| South Carolina DHEC | http://www.scdhec.gov/environment/envserv/ | 61002 |
| Texas CEQ | http://www.tceq.texas.gov/field/qa/env_lab_accreditation.html | T104704427 |
| Washington DOE | http://www.ecy.wa.gov/programs/eap/labs/lab-accreditation.html | C544 |
| Wisconsin DNR | http://dnr.wi.gov/ | 998386840 |
| Wyoming (EPA Region 8) | http://www.epa.gov/region8/water/dwhome/wyomingdi.html | - |
| Kelso Laboratory Website | www.alsglobal.com | NA |

Analyses were performed according to our laboratory's NELAP-approved quality assurance program. A complete listing of specific NELAP-certified analytes, can be found in the certification section at www.ALSGlobal.com or at the accreditation bodies web site.

Please refer to the certification and/or accreditation body's web site if samples are submitted for compliance purposes. The states highlighted above, require the analysis be listed on the state certification if used for compliance purposes and if the method/analyte is offered by that state.



Case Narrative


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www.alsglobal.com

ALS ENVIRONMENTAL

| | | | |
|-----------------------|--|-----------------------------|----------|
| Client: | Eurofins Calscience Environmental Laboratory | Service Request No.: | K1512169 |
| Project: | 15-10-1856 | Date Received: | 10/27/15 |
| Sample Matrix: | Water | | |

Case Narrative

All analyses were performed consistent with the quality assurance program of ALS Environmental. This report contains analytical results for samples designated for Tier II data deliverables. When appropriate to the method, method blank results have been reported with each analytical test. Surrogate recoveries have been reported for all applicable organic analyses. Additional quality control analyses reported herein include: Laboratory Control Sample (LCS), and Laboratory/Duplicate Laboratory Control Sample (LCS/DLCS).

Sample Receipt

Eight water samples were received for analysis at ALS Environmental on 10/27/15. The samples were received in good condition and consistent with the accompanying chain of custody form. The samples were stored in a refrigerator at 4°C upon receipt at the laboratory.

Steroids and Endocrine Disrupting Compounds by Method 1694**Elevated Detection Limits:**

Samples HCS 210 Lead and HCS 250 Lead required dilution due to the presence of elevated levels of target analyte. The reporting limits were adjusted to reflect the dilution.

No other anomalies associated with the analysis of these samples were observed.

Approved by _____



Chain of Custody


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www.alsglobal.com

To: ALS - Kelso

DATE: 10/26/15

PAGE: 1 OF 1

K1512169

[illegible]



Cooler Receipt and Preservation Form

Client / Project: EUROFINS Service Request K15 12169
Received: 10/27/15 Opened: 10/27/15 By: [Signature] Unloaded: 10/27/15 By: [Signature]

1. Samples were received via? Mail Fed Ex UPS DHL PDX Courier Hand Delivered
2. Samples were received in: (circle) Cooler Box Envelope Other NA
3. Were custody seals on coolers? NA Y N If yes, how many and where? 1 FRONT
If present, were custody seals intact? Y N If present, were they signed and dated? Y N

| Raw Cooler Temp | Corrected Cooler Temp | Raw Temp Blank | Corrected Temp Blank | Corr. Factor | Thermometer ID | Cooler/COC ID | Tracking Number | NA | Filed |
|-----------------|-----------------------|----------------|----------------------|--------------|----------------|---------------|-----------------|----|-------|
| 0.0 | 0.2 | 1.1 | 1.3 | 0.2 | 342 | | 774828956036 | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |

4. Packing material: Inserts Baggies Bubble Wrap Gel Packs Wet Ice Dry Ice Sleeves
5. Were custody papers properly filled out (ink, signed, etc.)? NA Y N
6. Did all bottles arrive in good condition (unbroken)? Indicate in the table below. NA Y N
7. Were all sample labels complete (i.e analysis, preservation, etc.)? NA Y N
8. Did all sample labels and tags agree with custody papers? Indicate major discrepancies in the table on page 2. NA Y N
9. Were appropriate bottles/containers and volumes received for the tests indicated? NA Y N
10. Were the pH-preserved bottles (*see SMO GEN SOP*) received at the appropriate pH? Indicate in the table below NA Y N
11. Were VOA vials received without headspace? Indicate in the table below. NA Y N
12. Was C12/Res negative? NA Y N

| Sample ID on Bottle | Sample ID on COC | Identified by: |
|---------------------|------------------|----------------|
| | | |
| | | |
| | | |

| Sample ID | Bottle Count | Bottle Type | Out of Temp | Head-space | Broke | pH | Reagent | Volume added | Reagent Lot Number | Initials | Time |
|-----------|--------------|-------------|-------------|------------|-------|----|---------|--------------|--------------------|----------|------|
| | | | | | | | | | | | |
| | | | | | | | | | | | |
| | | | | | | | | | | | |
| | | | | | | | | | | | |
| | | | | | | | | | | | |

Notes, Discrepancies, & Resolutions: _____



Steroids and Endocrine Disrupting Compounds


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Calscience



WORK ORDER NUMBER: 15-10-1995

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: SWCA Environmental Consultants

Client Project Name: EAA 27122

Attention: Philip Pearce
6200 UTSA Blvd.
Suite 102
San Antonio, TX 78249-1618

Approved for release on 11/23/2015 by:
Don Burley
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

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Work Order Number: 15-10-1995

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Work Order Narrative

Work Order: 15-10-1995

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Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 10/27/15. They were assigned to Work Order 15-10-1995.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.



Calscience

Sample Summary

| | | |
|--|-----------------------|----------------|
| Client: SWCA Environmental Consultants | Work Order: | 15-10-1995 |
| 6200 UTSA Blvd., Suite 102 | Project Name: | EAA 27122 |
| San Antonio, TX 78249-1618 | PO Number: | |
| | Date/Time Received: | 10/27/15 14:30 |
| | Number of Containers: | 307 |

Attn: Philip Pearce

| Sample Identification | Lab Number | Collection Date and Time | Number of Containers | Matrix |
|-----------------------|---------------|--------------------------|----------------------|---------|
| HCS 260 Peak | 15-10-1995-1 | 10/23/15 15:01 | 9 | Aqueous |
| HCS 270 Peak | 15-10-1995-2 | 10/23/15 15:16 | 11 | Aqueous |
| HCS 210 Trail | 15-10-1995-3 | 10/23/15 16:14 | 9 | Aqueous |
| HCS 240 Trail | 15-10-1995-4 | 10/23/15 16:29 | 9 | Aqueous |
| HCS 250 Trail | 15-10-1995-5 | 10/23/15 17:09 | 11 | Aqueous |
| HCS 260 Trail | 15-10-1995-6 | 10/23/15 16:50 | 11 | Aqueous |
| HCS 270 Trail | 15-10-1995-7 | 10/23/15 17:34 | 9 | Aqueous |
| FDHCS 260 Trail | 15-10-1995-8 | 10/23/15 16:50 | 9 | Aqueous |
| FDHCS 270 Trail | 15-10-1995-9 | 10/23/15 17:34 | 9 | Aqueous |
| TB15 | 15-10-1995-10 | 10/23/15 00:00 | 1 | Aqueous |
| HSM 210 Trail | 15-10-1995-11 | 10/23/15 20:35 | 9 | Aqueous |
| HSM 230 Trail | 15-10-1995-12 | 10/23/15 21:01 | 9 | Aqueous |
| HSM 231 Trail | 15-10-1995-13 | 10/23/15 21:21 | 9 | Aqueous |
| HSM 240 Trail | 15-10-1995-14 | 10/23/15 20:51 | 9 | Aqueous |
| HSM 250 Trail | 15-10-1995-15 | 10/23/15 21:06 | 9 | Aqueous |
| HSM 260 Trail | 15-10-1995-16 | 10/23/15 21:35 | 9 | Aqueous |
| HSM 270 Trail | 15-10-1995-17 | 10/23/15 21:55 | 11 | Aqueous |
| FDHSM 210 Trail | 15-10-1995-18 | 10/23/15 20:35 | 11 | Aqueous |
| FDHSM 230 Trail | 15-10-1995-19 | 10/23/15 21:01 | 9 | Aqueous |
| FDHSM 231 Trail | 15-10-1995-20 | 10/23/15 21:21 | 9 | Aqueous |
| HSM 210 Lead | 15-10-1995-21 | 10/23/15 15:25 | 11 | Aqueous |
| HSM 230 Lead | 15-10-1995-22 | 10/23/15 15:41 | 9 | Aqueous |
| HSM 231 Lead | 15-10-1995-23 | 10/23/15 16:02 | 9 | Aqueous |
| HSM 240 Lead | 15-10-1995-24 | 10/23/15 15:30 | 9 | Aqueous |
| HSM 250 Lead | 15-10-1995-25 | 10/23/15 16:21 | 11 | Aqueous |
| HSM 260 Lead | 15-10-1995-26 | 10/23/15 15:58 | 9 | Aqueous |
| HSM 270 Lead | 15-10-1995-27 | 10/23/15 16:15 | 11 | Aqueous |
| HSM 210 Peak | 15-10-1995-28 | 10/23/15 17:02 | 9 | Aqueous |
| HSM 230 Peak | 15-10-1995-29 | 10/23/15 17:16 | 9 | Aqueous |
| HSM 231 Peak | 15-10-1995-30 | 10/23/15 17:06 | 9 | Aqueous |
| HSM 240 Peak | 15-10-1995-31 | 10/23/15 17:34 | 9 | Aqueous |
| HSM 250 Peak | 15-10-1995-32 | 10/23/15 17:05 | 0 | Aqueous |
| HSM 260 Peak | 15-10-1995-33 | 10/23/15 17:25 | 9 | Aqueous |
| HSM 270 Peak | 15-10-1995-34 | 10/23/15 17:45 | 11 | Aqueous |


 Return to Contents



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-10-1995-1-I | 10/23/15 15:01 | Aqueous | IC 15 | N/A | 10/31/15 22:20 | 151031L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.19 | 0.10 | 0.027 | 1.00 | |
| Chloride | 16 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.090 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.6 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 27 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|-------|-----|----------------|-----------|
| HCS 270 Peak | 15-10-1995-2-I | 10/23/15 15:16 | Aqueous | IC 15 | N/A | 10/31/15 22:39 | 151031L02 |
|--------------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.19 | 0.10 | 0.027 | 1.00 | |
| Chloride | 17 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.091 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.6 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 27 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|-------|-----|----------------|-----------|
| HCS 210 Trail | 15-10-1995-3-I | 10/23/15 16:14 | Aqueous | IC 15 | N/A | 10/31/15 22:57 | 151031L02 |
|---------------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.10 | 0.10 | 0.027 | 1.00 | |
| Chloride | 7.2 | 1.0 | 0.52 | 1.00 | |
| Bromide | ND | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 0.68 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 11 | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-10-1995-4-I | 10/23/15 16:29 | Aqueous | IC 15 | N/A | 10/31/15 23:15 | 151031L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.20 | 0.10 | 0.027 | 1.00 | |
| Chloride | 17 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.086 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.8 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 26 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|-------|-----|----------------|-----------|
| HCS 250 Trail | 15-10-1995-5-I | 10/23/15 17:09 | Aqueous | IC 15 | N/A | 10/31/15 23:34 | 151031L02 |
|---------------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.18 | 0.10 | 0.027 | 1.00 | |
| Chloride | 16 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.073 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.6 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 24 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|-------|-----|----------------|-----------|
| HCS 260 Trail | 15-10-1995-6-I | 10/23/15 16:50 | Aqueous | IC 15 | N/A | 10/31/15 23:52 | 151031L02 |
|---------------|----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.19 | 0.10 | 0.027 | 1.00 | |
| Chloride | 17 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.099 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.7 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 30 | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-10-1995-7-I | 10/23/15 17:34 | Aqueous | IC 15 | N/A | 11/01/15 00:10 | 151031L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.20 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.088 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.6 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 28 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|-----------------|----------------|-------------------|---------|-------|-----|-------------------|-----------|
| FDHCS 260 Trail | 15-10-1995-8-I | 10/23/15 16:50 | Aqueous | IC 15 | N/A | 11/01/15 00:29 | 151031L02 |
|-----------------|----------------|-------------------|---------|-------|-----|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.20 | 0.10 | 0.027 | 1.00 | |
| Chloride | 17 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.10 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.7 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 30 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|-----------------|----------------|-------------------|---------|-------|-----|-------------------|-----------|
| FDHCS 270 Trail | 15-10-1995-9-I | 10/23/15 17:34 | Aqueous | IC 15 | N/A | 11/01/15 00:47 | 151031L02 |
|-----------------|----------------|-------------------|---------|-------|-----|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.19 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.089 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.6 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 28 | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-10-1995-11-I | 10/23/15 20:35 | Aqueous | IC 15 | N/A | 11/01/15 01:05 | 151031L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.21 | 0.10 | 0.027 | 1.00 | |
| Chloride | 27 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.17 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 0.61 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 34 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|-------|-----|----------------|-----------|
| HSM 230 Trail | 15-10-1995-12-I | 10/23/15 21:01 | Aqueous | IC 15 | N/A | 11/01/15 01:04 | 151031L02 |
|---------------|-----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.18 | 0.10 | 0.027 | 1.00 | |
| Chloride | 19 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.10 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.5 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 30 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|-------|-----|----------------|-----------|
| HSM 231 Trail | 15-10-1995-13-I | 10/23/15 21:21 | Aqueous | IC 15 | N/A | 11/01/15 01:23 | 151031L02 |
|---------------|-----------------|----------------|---------|-------|-----|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------|--------|------|-------|------|------------|
| Fluoride | 0.18 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.10 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.3 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 25 | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| HSM 240 Trail | 15-10-1995-14-I | 10/23/15 20:51 | Aqueous | IC 15 | N/A | 11/01/15 01:41 | 151031L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.18 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.096 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 25 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 250 Trail | 15-10-1995-15-I | 10/23/15 21:06 | Aqueous | IC 15 | N/A | 11/01/15 01:59 | 151031L02 |
|----------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.18 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.11 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 26 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 260 Trail | 15-10-1995-16-I | 10/23/15 21:35 | Aqueous | IC 15 | N/A | 11/01/15 02:18 | 151031L02 |
|----------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.17 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.098 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 26 | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| HSM 270 Trail | 15-10-1995-17-I | 10/23/15 21:55 | Aqueous | IC 15 | N/A | 11/01/15 02:36 | 151031L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.19 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.10 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 26 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|------------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| FDHSM 210 Trail | 15-10-1995-18-I | 10/23/15 20:35 | Aqueous | IC 15 | N/A | 11/01/15 02:55 | 151031L02 |
|------------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.20 | 0.10 | 0.027 | 1.00 | |
| Chloride | 27 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.17 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 0.62 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 34 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|------------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| FDHSM 230 Trail | 15-10-1995-19-I | 10/23/15 21:01 | Aqueous | IC 15 | N/A | 11/01/15 03:13 | 151031L02 |
|------------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.20 | 0.10 | 0.027 | 1.00 | |
| Chloride | 19 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.11 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.6 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 30 | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| FDHSM 231 Trail | 15-10-1995-20-I | 10/23/15 21:21 | Aqueous | IC 15 | N/A | 11/01/15 03:31 | 151031L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.18 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.10 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 25 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 210 Lead | 15-10-1995-21-I | 10/23/15 15:25 | Aqueous | IC 15 | N/A | 11/01/15 03:50 | 151031L02 |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.21 | 0.10 | 0.027 | 1.00 | |
| Chloride | 27 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.17 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 0.62 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 33 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 230 Lead | 15-10-1995-22-I | 10/23/15 15:41 | Aqueous | IC 15 | N/A | 10/31/15 14:32 | 151031L01 |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.12 | 0.10 | 0.027 | 1.00 | |
| Chloride | 10 | 1.0 | 0.52 | 1.00 | |
| Bromide | ND | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 0.94 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 14 | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| HSM 231 Lead | 15-10-1995-23-I | 10/23/15 16:02 | Aqueous | IC 15 | N/A | 10/31/15 14:50 | 151031L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.17 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.092 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 24 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 240 Lead | 15-10-1995-24-I | 10/23/15 15:30 | Aqueous | IC 15 | N/A | 10/31/15 15:08 | 151031L01 |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.16 | 0.10 | 0.027 | 1.00 | |
| Chloride | 16 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.084 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.1 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 23 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 250 Lead | 15-10-1995-25-I | 10/23/15 16:21 | Aqueous | IC 15 | N/A | 10/31/15 15:27 | 151031L01 |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.17 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.092 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 24 | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| HSM 260 Lead | 15-10-1995-26-I | 10/23/15 15:58 | Aqueous | IC 15 | N/A | 10/31/15 15:45 | 151031L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.16 | 0.10 | 0.027 | 1.00 | |
| Chloride | 17 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.086 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 24 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 270 Lead | 15-10-1995-27-I | 10/23/15 16:15 | Aqueous | IC 15 | N/A | 10/31/15 17:59 | 151031L01 |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.19 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.10 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.1 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 25 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 210 Peak | 15-10-1995-28-I | 10/23/15 17:02 | Aqueous | IC 15 | N/A | 10/31/15 18:18 | 151031L01 |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.21 | 0.10 | 0.027 | 1.00 | |
| Chloride | 27 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.16 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 0.60 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 33 | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| HSM 230 Peak | 15-10-1995-29-I | 10/23/15 17:16 | Aqueous | IC 15 | N/A | 10/31/15 18:36 | 151031L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.18 | 0.10 | 0.027 | 1.00 | |
| Chloride | 21 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.11 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.6 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 28 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 231 Peak | 15-10-1995-30-I | 10/23/15 17:06 | Aqueous | IC 15 | N/A | 10/31/15 18:54 | 151031L01 |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.17 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.10 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 25 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 240 Peak | 15-10-1995-31-I | 10/23/15 17:34 | Aqueous | IC 15 | N/A | 10/31/15 19:13 | 151031L01 |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.18 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.10 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 25 | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|---------------|-----------------------|------------------|
| HSM 260 Peak | 15-10-1995-33-I | 10/23/15 17:25 | Aqueous | IC 15 | N/A | 10/31/15 19:31 | 151031L01 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.17 | 0.10 | 0.027 | 1.00 | |
| Chloride | 17 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.096 | 0.10 | 0.058 | 1.00 | J |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 23 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|
| HSM 270 Peak | 15-10-1995-34-I | 10/23/15 17:45 | Aqueous | IC 15 | N/A | 10/31/15 19:49 | 151031L01 |
|---------------------|------------------------|-----------------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | 0.17 | 0.10 | 0.027 | 1.00 | |
| Chloride | 18 | 1.0 | 0.52 | 1.00 | |
| Bromide | 0.10 | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | 1.2 | 0.10 | 0.053 | 1.00 | BU |
| Sulfate | 25 | 1.0 | 0.27 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|------------|----------------|--------------|------------|-----------------------|------------------|
| Method Blank | 099-12-906-6212 | N/A | Aqueous | IC 15 | N/A | 10/31/15 12:03 | 151031L01 |
|---------------------|------------------------|------------|----------------|--------------|------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | ND | 0.10 | 0.027 | 1.00 | |
| Chloride | ND | 1.0 | 0.52 | 1.00 | |
| Bromide | ND | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | ND | 0.10 | 0.053 | 1.00 | |
| Sulfate | ND | 1.0 | 0.27 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-906-6215 | N/A | Aqueous | IC 15 | N/A | 10/31/15 21:43 | 151031L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Fluoride | ND | 0.10 | 0.027 | 1.00 | |
| Chloride | ND | 1.0 | 0.52 | 1.00 | |
| Bromide | ND | 0.10 | 0.058 | 1.00 | |
| Nitrate (as N) | ND | 0.10 | 0.053 | 1.00 | |
| Sulfate | ND | 1.0 | 0.27 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

Page 1 of 12

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-10-1995-1-F | 10/23/15 15:01 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:11 | 151028LA8F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 66.4 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 13.9 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.82 | 0.500 | 0.103 | 1.00 | |
| Sodium | 10.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.550 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.92 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS 270 Peak | 15-10-1995-2-F | 10/23/15 15:16 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:18 | 151028LA8F |
|--------------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 73.4 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 15.4 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.05 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.598 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.47 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS 210 Trail | 15-10-1995-3-F | 10/23/15 16:14 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:20 | 151028LA8F |
|---------------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 34.0 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 7.00 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.77 | 0.500 | 0.103 | 1.00 | |
| Sodium | 5.09 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.267 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 2.66 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

Page 2 of 12

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-10-1995-4-F | 10/23/15 16:29 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:22 | 151028LA8F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 86.8 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 17.9 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.86 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.7 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.697 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 6.56 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS 250 Trail | 15-10-1995-5-F | 10/23/15 17:09 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:29 | 151028LA8F |
|---------------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 71.8 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 15.3 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.42 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.587 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.66 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|----------|----------|----------------|------------|
| HCS 260 Trail | 15-10-1995-6-F | 10/23/15 16:50 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:31 | 151028LA8F |
|---------------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 79.4 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 16.6 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.87 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.7 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.651 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.90 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-10-1995-7-F | 10/23/15 17:34 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:33 | 151028LA8F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 73.8 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 15.3 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.04 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.605 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.52 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|-----------------|----------------|----------------|---------|----------|----------|----------------|------------|
| FDHCS 260 Trail | 15-10-1995-8-F | 10/23/15 16:50 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:36 | 151028LA8F |
|-----------------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 78.0 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 16.5 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.83 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.7 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.647 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.83 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|-----------------|----------------|----------------|---------|----------|----------|----------------|------------|
| FDHCS 270 Trail | 15-10-1995-9-F | 10/23/15 17:34 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:38 | 151028LA8F |
|-----------------|----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 71.5 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 14.8 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.99 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.593 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.29 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-10-1995-11-F | 10/23/15 20:35 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:40 | 151028LA8F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 78.6 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 20.0 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.01 | 0.500 | 0.103 | 1.00 | |
| Sodium | 15.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.658 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.31 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|----------|----------|----------------|------------|
| HSM 230 Trail | 15-10-1995-12-F | 10/23/15 21:01 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:43 | 151028LA8F |
|---------------|-----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 83.2 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 15.1 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.25 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.3 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.471 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.22 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|----------|----------|----------------|------------|
| HSM 231 Trail | 15-10-1995-13-F | 10/23/15 21:21 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:45 | 151028LA8F |
|---------------|-----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 87.7 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 18.6 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.51 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.529 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.74 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-10-1995-14-F | 10/23/15 20:51 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:47 | 151028LA8F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 88.1 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 18.3 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.52 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.5 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.529 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.67 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|----------|----------|----------------|------------|
| HSM 250 Trail | 15-10-1995-15-F | 10/23/15 21:06 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:50 | 151028LA8F |
|---------------|-----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 93.0 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 19.2 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.62 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.2 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.562 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.96 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|----------|----------|----------------|------------|
| HSM 260 Trail | 15-10-1995-16-F | 10/23/15 21:35 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:56 | 151028LA8F |
|---------------|-----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 86.2 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 18.4 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.72 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.6 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.523 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.70 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Trail | 15-10-1995-17-F | 10/23/15 21:55 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:58 | 151028LA8F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 89.1 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 18.5 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.66 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.537 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.77 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|-----------------|-----------------|----------------|---------|----------|----------|----------------|------------|
| FDHSM 210 Trail | 15-10-1995-18-F | 10/23/15 20:35 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:01 | 151028LA8F |
|-----------------|-----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 80.4 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 20.7 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.08 | 0.500 | 0.103 | 1.00 | |
| Sodium | 15.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.675 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.45 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|-----------------|-----------------|----------------|---------|----------|----------|----------------|------------|
| FDHSM 230 Trail | 15-10-1995-19-F | 10/23/15 21:01 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:03 | 151028LA8F |
|-----------------|-----------------|----------------|---------|----------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 80.9 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 15.1 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.22 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.2 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.465 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.18 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| FDHSM 231 Trail | 15-10-1995-20-F | 10/23/15 21:21 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:05 | 151028LA8F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 86.2 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 18.2 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.55 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.4 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.524 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.62 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 210 Lead | 15-10-1995-21-F | 10/23/15 15:25 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:08 | 151028LA8F |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 77.9 | 0.100 | 0.0118 | 1.00 | B |
| Magnesium | 19.9 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.99 | 0.500 | 0.103 | 1.00 | |
| Sodium | 15.3 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.648 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.28 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 230 Lead | 15-10-1995-22-F | 10/23/15 15:41 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 12:30 | 151028LA9F |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|--------|---------|------|------------|
| Calcium | 44.0 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 7.44 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 4.13 | 0.500 | 0.103 | 1.00 | |
| Sodium | 7.86 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.226 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 2.71 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 231 Lead | 15-10-1995-23-F | 10/23/15 16:02 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:10 | 151028LA9F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 83.4 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.6 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.85 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.3 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.504 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.44 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 240 Lead | 15-10-1995-24-F | 10/23/15 15:30 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:12 | 151028LA9F |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 75.4 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 15.8 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.72 | 0.500 | 0.103 | 1.00 | |
| Sodium | 10.3 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.454 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.08 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 250 Lead | 15-10-1995-25-F | 10/23/15 16:21 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:15 | 151028LA9F |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 87.3 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 18.3 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.86 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.530 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.68 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 260 Lead | 15-10-1995-26-F | 10/23/15 15:58 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:17 | 151028LA9F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 82.9 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.8 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.73 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.502 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.49 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 270 Lead | 15-10-1995-27-F | 10/23/15 16:15 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:23 | 151028LA9F |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 81.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.2 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.83 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.6 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.502 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.34 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 210 Peak | 15-10-1995-28-F | 10/23/15 17:02 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:26 | 151028LA9F |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 78.1 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 20.2 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.99 | 0.500 | 0.103 | 1.00 | |
| Sodium | 15.5 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.659 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.34 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 230 Peak | 15-10-1995-29-F | 10/23/15 17:16 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:28 | 151028LA9F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 83.2 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 16.3 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 2.02 | 0.500 | 0.103 | 1.00 | |
| Sodium | 12.9 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.498 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.43 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 231 Peak | 15-10-1995-30-F | 10/23/15 17:06 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:30 | 151028LA9F |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 87.3 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 18.7 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.61 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.8 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.536 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.73 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 240 Peak | 15-10-1995-31-F | 10/23/15 17:34 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:33 | 151028LA9F |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 98.7 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 20.6 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.71 | 0.500 | 0.103 | 1.00 | |
| Sodium | 13.1 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.605 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 6.32 | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 260 Peak | 15-10-1995-33-F | 10/23/15 17:25 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:35 | 151028LA9F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 81.9 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 17.0 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.80 | 0.500 | 0.103 | 1.00 | |
| Sodium | 11.0 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.494 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 5.26 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 270 Peak | 15-10-1995-34-F | 10/23/15 17:45 | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 20:37 | 151028LA9F |
|---------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 75.8 | 0.100 | 0.0118 | 1.00 | |
| Magnesium | 15.8 | 0.100 | 0.00336 | 1.00 | |
| Potassium | 1.80 | 0.500 | 0.103 | 1.00 | |
| Sodium | 10.9 | 0.500 | 0.103 | 1.00 | |
| Strontium | 0.461 | 0.0300 | 0.00277 | 1.00 | |
| Silicon | 4.86 | 0.0500 | 0.0279 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| Method Blank | 099-15-683-1458 | N/A | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:02 | 151028LA8F |
|---------------------|------------------------|------------|----------------|-----------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | 0.0244 | 0.100 | 0.0118 | 1.00 | J |
| Magnesium | ND | 0.100 | 0.00336 | 1.00 | |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-683-1442 | N/A | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 12:20 | 151028LA9F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Calcium | ND | 0.100 | 0.0118 | 1.00 | |
| Magnesium | ND | 0.100 | 0.00336 | 1.00 | |
| Potassium | ND | 0.500 | 0.103 | 1.00 | |
| Sodium | ND | 0.500 | 0.103 | 1.00 | |
| Strontium | ND | 0.0300 | 0.00277 | 1.00 | |
| Silicon | ND | 0.0500 | 0.0279 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-10-1995-1-F | 10/23/15 15:01 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 22:14 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0471 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000655 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00139 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000363 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0225 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00593 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0501 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00576 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Peak | 15-10-1995-2-F | 10/23/15 15:16 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 21:39 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000215 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0510 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000431 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.000375 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | 0.000152 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00152 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000376 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0754 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00684 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0548 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00325 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Trail | 15-10-1995-3-F | 10/23/15 16:14 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 22:18 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0290 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000449 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.00136 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000135 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.000992 | 0.00100 | 0.000132 | 1.00 | J |
| Selenium | 0.000214 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0118 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0145 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0640 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00133 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-10-1995-4-F | 10/23/15 16:29 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 22:21 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0585 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000382 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00179 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000418 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00341 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | 0.00615 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0596 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00112 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Trail | 15-10-1995-5-F | 10/23/15 17:09 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 22:25 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0499 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000814 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00155 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000392 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00549 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0540 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00168 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Trail | 15-10-1995-6-F | 10/23/15 16:50 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 22:28 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0535 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000560 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00152 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000270 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00363 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | 0.00338 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0502 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00299 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-10-1995-7-F | 10/23/15 17:34 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 23:17 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0512 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000721 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00145 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000350 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00559 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00724 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0553 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.000998 | 0.00100 | 0.000139 | 1.00 | J |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 260 Trail | 15-10-1995-8-F | 10/23/15 16:50 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 23:21 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0529 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000267 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00152 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000327 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0140 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00574 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0559 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 270 Trail | 15-10-1995-9-F | 10/23/15 17:34 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 23:24 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0490 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000496 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00139 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000318 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0101 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00581 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0502 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.000298 | 0.00100 | 0.000139 | 1.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-10-1995-11-F | 10/23/15 20:35 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 23:28 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0341 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00157 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000261 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0193 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00570 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0519 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0865 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Trail | 15-10-1995-12-F | 10/23/15 21:01 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 23:31 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000173 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0357 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00118 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00192 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000374 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0401 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00775 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0568 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00236 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 231 Trail | 15-10-1995-13-F | 10/23/15 21:21 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 23:35 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0367 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000193 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00170 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000309 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0136 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00682 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0550 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.000472 | 0.00100 | 0.000139 | 1.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-10-1995-14-F | 10/23/15 20:51 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 23:39 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000130 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0370 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000110 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00182 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000270 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0462 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00873 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0820 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-10-1995-15-F | 10/23/15 21:06 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 23:42 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0388 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000657 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00227 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000301 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0120 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00982 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0588 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Trail | 15-10-1995-16-F | 10/23/15 21:35 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 23:46 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0372 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000208 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00172 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000303 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0102 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00679 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0496 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Trail | 15-10-1995-17-F | 10/23/15 21:55 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 23:49 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0379 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000336 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00186 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000318 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0101 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.0244 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0655 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.000618 | 0.00100 | 0.000139 | 1.00 | J |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 210 Trail | 15-10-1995-18-F | 10/23/15 20:35 | Aqueous | ICP/MS 03 | 10/28/15 | 10/31/15 00:03 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0347 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000179 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00165 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00791 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00569 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0609 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0846 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 230 Trail | 15-10-1995-19-F | 10/23/15 21:01 | Aqueous | ICP/MS 03 | 10/28/15 | 10/31/15 00:07 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000143 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0369 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00151 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00188 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000270 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00928 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00881 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0529 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00354 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 231 Trail | 15-10-1995-20-F | 10/23/15 21:21 | Aqueous | ICP/MS 03 | 10/28/15 | 10/31/15 00:10 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0374 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00183 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000292 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0502 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00514 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0707 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Lead | 15-10-1995-21-F | 10/23/15 15:25 | Aqueous | ICP/MS 03 | 10/28/15 | 10/31/15 00:14 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0343 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00161 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000255 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0383 | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | 0.00560 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0473 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.0940 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Lead | 15-10-1995-22-F | 10/23/15 15:41 | Aqueous | ICP/MS 03 | 10/28/15 | 10/31/15 00:17 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000927 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0238 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.00131 | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00545 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000194 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00173 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000242 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0602 | 0.00500 | 0.000479 | 1.00 | B |
| Aluminum | 0.0223 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0947 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0126 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|------------------|-----------------|-----------------------|-------------------|
| HSM 231 Lead | 15-10-1995-23-F | 10/23/15 16:02 | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 22:11 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Antimony | 0.000167 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0362 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000675 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00177 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000294 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0290 | 0.00500 | 0.000479 | 1.00 | B |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0572 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00412 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|------------------|-----------------|-----------------------|-------------------|
| HSM 240 Lead | 15-10-1995-24-F | 10/23/15 15:30 | Aqueous | ICP/MS 03 | 10/28/15 | 10/31/15 00:21 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Antimony | 0.000363 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0354 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00226 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000336 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00191 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000332 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0138 | 0.00500 | 0.000479 | 1.00 | B |
| Aluminum | 0.0887 | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.116 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00920 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Lead | 15-10-1995-25-F | 10/23/15 16:21 | Aqueous | ICP/MS 03 | 10/28/15 | 10/31/15 00:24 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000145 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0393 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000572 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00189 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000279 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0181 | 0.00500 | 0.000479 | 1.00 | B |
| Aluminum | 0.00580 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0599 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00117 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Lead | 15-10-1995-26-F | 10/23/15 15:58 | Aqueous | ICP/MS 03 | 10/28/15 | 10/31/15 00:28 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0363 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000280 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00181 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000309 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0135 | 0.00500 | 0.000479 | 1.00 | B |
| Aluminum | 0.0102 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0565 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Lead | 15-10-1995-27-F | 10/23/15 16:15 | Aqueous | ICP/MS 03 | 10/28/15 | 10/31/15 00:31 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000111 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0370 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000360 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00180 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000275 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0149 | 0.00500 | 0.000479 | 1.00 | B |
| Aluminum | 0.00652 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0490 | 0.0500 | 0.00926 | 1.00 | J |
| Manganese | 0.00131 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Peak | 15-10-1995-28-F | 10/23/15 17:02 | Aqueous | ICP/MS 03 | 10/28/15 | 10/31/15 00:35 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0347 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00104 | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00172 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000233 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0161 | 0.00500 | 0.000479 | 1.00 | B |
| Aluminum | 0.00726 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0547 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0803 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Peak | 15-10-1995-29-F | 10/23/15 17:16 | Aqueous | ICP/MS 03 | 10/28/15 | 11/03/15 00:39 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000176 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0381 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000905 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00150 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000198 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0148 | 0.00500 | 0.000479 | 1.00 | B |
| Aluminum | 0.00555 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0560 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00716 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 231 Peak | 15-10-1995-30-F | 10/23/15 17:06 | Aqueous | ICP/MS 03 | 10/28/15 | 11/03/15 00:43 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0379 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.000644 | 0.00100 | 0.000402 | 1.00 | J |
| Copper | 0.000487 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00155 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000180 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0121 | 0.00500 | 0.000479 | 1.00 | B |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0709 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0117 | 0.00100 | 0.000139 | 1.00 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Peak | 15-10-1995-31-F | 10/23/15 17:34 | Aqueous | ICP/MS 03 | 10/28/15 | 11/03/15 00:46 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000150 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0418 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000355 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00171 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000214 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0257 | 0.00500 | 0.000479 | 1.00 | B |
| Aluminum | 0.00399 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.0681 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00266 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Peak | 15-10-1995-33-F | 10/23/15 17:25 | Aqueous | ICP/MS 03 | 10/28/15 | 11/03/15 00:50 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000465 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0369 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.00102 | 0.00100 | 0.000140 | 1.00 | |
| Lead | 0.000128 | 0.00100 | 0.0000898 | 1.00 | J |
| Nickel | 0.00153 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0322 | 0.00500 | 0.000479 | 1.00 | B |
| Aluminum | 0.0185 | 0.0500 | 0.00331 | 1.00 | J |
| Iron | 0.184 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.00516 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Peak | 15-10-1995-34-F | 10/23/15 17:45 | Aqueous | ICP/MS 03 | 10/28/15 | 11/03/15 00:53 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|----------|---------|-----------|------|------------|
| Antimony | 0.000144 | 0.00100 | 0.0000995 | 1.00 | J |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | 0.0331 | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | 0.00100 | 0.00100 | 0.000402 | 1.00 | |
| Copper | 0.000851 | 0.00100 | 0.000140 | 1.00 | J |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | 0.00150 | 0.00100 | 0.000132 | 1.00 | |
| Selenium | 0.000190 | 0.00100 | 0.000168 | 1.00 | J |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.0159 | 0.00500 | 0.000479 | 1.00 | B |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | 0.0634 | 0.0500 | 0.00926 | 1.00 | |
| Manganese | 0.0106 | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-963 | N/A | Aqueous | ICP/MS 03 | 10/28/15 | 11/06/15 16:19 | 151028LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | ND | 0.00500 | 0.000479 | 1.00 | |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-693-964 | N/A | Aqueous | ICP/MS 03 | 10/28/15 | 11/06/15 16:22 | 151028LA5F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|---------|-----------|------|------------|
| Antimony | ND | 0.00100 | 0.0000995 | 1.00 | |
| Arsenic | ND | 0.00100 | 0.000386 | 1.00 | |
| Barium | ND | 0.00100 | 0.0000986 | 1.00 | |
| Beryllium | ND | 0.00100 | 0.000290 | 1.00 | |
| Cadmium | ND | 0.00100 | 0.000128 | 1.00 | |
| Chromium | ND | 0.00100 | 0.000402 | 1.00 | |
| Copper | ND | 0.00100 | 0.000140 | 1.00 | |
| Lead | ND | 0.00100 | 0.0000898 | 1.00 | |
| Nickel | ND | 0.00100 | 0.000132 | 1.00 | |
| Selenium | ND | 0.00100 | 0.000168 | 1.00 | |
| Silver | ND | 0.00100 | 0.000111 | 1.00 | |
| Thallium | ND | 0.00100 | 0.000101 | 1.00 | |
| Zinc | 0.00185 | 0.00500 | 0.000479 | 1.00 | J |
| Aluminum | ND | 0.0500 | 0.00331 | 1.00 | |
| Iron | ND | 0.0500 | 0.00926 | 1.00 | |
| Manganese | ND | 0.00100 | 0.000139 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-10-1995-1-F | 10/23/15 15:01 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:25 | 151105LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 270 Peak | 15-10-1995-2-F | 10/23/15 15:16 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:27 | 151105LA3F |
|--------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 210 Trail | 15-10-1995-3-F | 10/23/15 16:14 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:18 | 151105LA3F |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 240 Trail | 15-10-1995-4-F | 10/23/15 16:29 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:34 | 151105LA3F |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 250 Trail | 15-10-1995-5-F | 10/23/15 17:09 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:36 | 151105LA3F |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|
| HCS 260 Trail | 15-10-1995-6-F | 10/23/15 16:50 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:39 | 151105LA3F |
|---------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-10-1995-7-F | 10/23/15 17:34 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:41 | 151105LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|-----------------|----------------|----------------|---------|------------|----------|----------------|------------|
| FDHCS 260 Trail | 15-10-1995-8-F | 10/23/15 16:50 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:43 | 151105LA3F |
|-----------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|-----------------|----------------|----------------|---------|------------|----------|----------------|------------|
| FDHCS 270 Trail | 15-10-1995-9-F | 10/23/15 17:34 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:45 | 151105LA3F |
|-----------------|----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|
| HSM 210 Trail | 15-10-1995-11-F | 10/23/15 20:35 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:48 | 151105LA3F |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|
| HSM 230 Trail | 15-10-1995-12-F | 10/23/15 21:01 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:50 | 151105LA3F |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|
| HSM 231 Trail | 15-10-1995-13-F | 10/23/15 21:21 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:52 | 151105LA3F |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-10-1995-14-F | 10/23/15 20:51 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:54 | 151105LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|
| HSM 250 Trail | 15-10-1995-15-F | 10/23/15 21:06 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:01 | 151105LA3F |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|
| HSM 260 Trail | 15-10-1995-16-F | 10/23/15 21:35 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:03 | 151105LA3F |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|
| HSM 270 Trail | 15-10-1995-17-F | 10/23/15 21:55 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:05 | 151105LA3F |
|---------------|-----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|------------|
| FDHSM 210 Trail | 15-10-1995-18-F | 10/23/15 20:35 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:08 | 151105LA3F |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|------------|
| FDHSM 230 Trail | 15-10-1995-19-F | 10/23/15 21:01 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:10 | 151105LA3F |
|-----------------|-----------------|----------------|---------|------------|----------|----------------|------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|--------|----------|-----------|------|------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| FDHSM 231 Trail | 15-10-1995-20-F | 10/23/15 21:21 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:12 | 151105LA3F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 210 Lead | 15-10-1995-21-F | 10/23/15 15:25 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:14 | 151105LA3F |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 230 Lead | 15-10-1995-22-F | 10/23/15 15:41 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:32 | 151105LA4F |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 231 Lead | 15-10-1995-23-F | 10/23/15 16:02 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:35 | 151105LA4F |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 240 Lead | 15-10-1995-24-F | 10/23/15 15:30 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:21 | 151105LA4F |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 250 Lead | 15-10-1995-25-F | 10/23/15 16:21 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:37 | 151105LA4F |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 260 Lead | 15-10-1995-26-F | 10/23/15 15:58 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:39 | 151105LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 270 Lead | 15-10-1995-27-F | 10/23/15 16:15 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:41 | 151105LA4F |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 210 Peak | 15-10-1995-28-F | 10/23/15 17:02 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:43 | 151105LA4F |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 230 Peak | 15-10-1995-29-F | 10/23/15 17:16 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:45 | 151105LA4F |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 231 Peak | 15-10-1995-30-F | 10/23/15 17:06 | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:48 | 151105LA4F |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 240 Peak | 15-10-1995-31-F | 10/23/15 17:34 | Aqueous | Mercury 04 | 11/05/15 | 11/06/15 14:31 | 151105LA4F |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 7470A Filt.
Method: EPA 7470A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 260 Peak | 15-10-1995-33-F | 10/23/15 17:25 | Aqueous | Mercury 04 | 11/05/15 | 11/06/15 14:33 | 151105LA4F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| HSM 270 Peak | 15-10-1995-34-F | 10/23/15 17:45 | Aqueous | Mercury 04 | 11/05/15 | 11/06/15 14:36 | 151105LA4F |
|---------------------|------------------------|-----------------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| Method Blank | 099-15-763-654 | N/A | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:14 | 151105LA3F |
|---------------------|-----------------------|------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

| | | | | | | | |
|---------------------|-----------------------|------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| Method Blank | 099-15-763-655 | N/A | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:17 | 151105LA4F |
|---------------------|-----------------------|------------|----------------|-------------------|-----------------|-----------------------|-------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Mercury | ND | 0.000500 | 0.0000453 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-10-1995-1-I | 10/23/15 15:01 | Aqueous | GC 44 | 10/29/15 | 10/30/15 19:34 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 103 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 100 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Peak | 15-10-1995-2-I | 10/23/15 15:16 | Aqueous | GC 44 | 10/29/15 | 10/30/15 19:49 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 106 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 105 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Trail | 15-10-1995-3-I | 10/23/15 16:14 | Aqueous | GC 44 | 10/29/15 | 10/30/15 20:03 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 95 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 99 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-10-1995-4-I | 10/23/15 16:29 | Aqueous | GC 44 | 10/29/15 | 10/30/15 20:17 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 107 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 102 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Trail | 15-10-1995-5-I | 10/23/15 17:09 | Aqueous | GC 44 | 10/29/15 | 10/30/15 20:32 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 103 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 101 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Trail | 15-10-1995-6-I | 10/23/15 16:50 | Aqueous | GC 44 | 10/29/15 | 10/30/15 20:46 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 97 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 91 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-10-1995-7-I | 10/23/15 17:34 | Aqueous | GC 44 | 10/29/15 | 10/30/15 21:00 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 106 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 103 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 260 Trail | 15-10-1995-8-I | 10/23/15 16:50 | Aqueous | GC 44 | 10/29/15 | 10/30/15 21:14 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 106 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 93 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 270 Trail | 15-10-1995-9-I | 10/23/15 17:34 | Aqueous | GC 44 | 10/29/15 | 10/30/15 21:29 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 105 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 93 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-10-1995-11-I | 10/23/15 20:35 | Aqueous | GC 44 | 10/29/15 | 10/30/15 21:43 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 84 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 76 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Trail | 15-10-1995-12-I | 10/23/15 21:01 | Aqueous | GC 44 | 10/29/15 | 10/30/15 21:57 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 102 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 92 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 231 Trail | 15-10-1995-13-I | 10/23/15 21:21 | Aqueous | GC 44 | 10/29/15 | 10/30/15 22:12 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 112 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 94 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-10-1995-14-I | 10/23/15 20:51 | Aqueous | GC 44 | 10/29/15 | 10/30/15 22:26 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 111 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 96 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-10-1995-15-I | 10/23/15 21:06 | Aqueous | GC 44 | 10/29/15 | 10/30/15 22:40 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 105 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 90 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Trail | 15-10-1995-16-I | 10/23/15 21:35 | Aqueous | GC 44 | 10/29/15 | 10/30/15 22:55 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 107 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 99 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Trail | 15-10-1995-17-I | 10/23/15 21:55 | Aqueous | GC 44 | 10/29/15 | 10/30/15 23:09 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 108 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 96 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 210 Trail | 15-10-1995-18-I | 10/23/15 20:35 | Aqueous | GC 44 | 10/29/15 | 10/30/15 23:23 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 75 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 75 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 230 Trail | 15-10-1995-19-I | 10/23/15 21:01 | Aqueous | GC 44 | 10/29/15 | 10/30/15 23:37 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 98 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 98 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 231 Trail | 15-10-1995-20-I | 10/23/15 21:21 | Aqueous | GC 44 | 10/29/15 | 10/30/15 23:52 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 109 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 100 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 210 Lead | 15-10-1995-21-I | 10/23/15 15:25 | Aqueous | GC 44 | 10/29/15 | 10/31/15 00:06 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 100 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 107 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Lead | 15-10-1995-22-I | 10/23/15 15:41 | Aqueous | GC 44 | 10/29/15 | 10/31/15 01:03 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | 0.090 | 0.10 | 0.028 | 1.00 | J |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 95 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 102 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 231 Lead | 15-10-1995-23-I | 10/23/15 16:02 | Aqueous | GC 44 | 10/29/15 | 10/31/15 01:17 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 103 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 101 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Lead | 15-10-1995-24-I | 10/23/15 15:30 | Aqueous | GC 44 | 10/29/15 | 10/31/15 01:32 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 95 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 96 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 250 Lead | 15-10-1995-25-I | 10/23/15 16:21 | Aqueous | GC 44 | 10/29/15 | 10/31/15 01:46 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 106 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 103 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Lead | 15-10-1995-26-I | 10/23/15 15:58 | Aqueous | GC 44 | 10/29/15 | 10/31/15 02:00 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 107 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 107 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 270 Lead | 15-10-1995-27-I | 10/23/15 16:15 | Aqueous | GC 44 | 10/29/15 | 10/31/15 02:15 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 104 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 104 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 210 Peak | 15-10-1995-28-I | 10/23/15 17:02 | Aqueous | GC 44 | 10/29/15 | 10/31/15 02:29 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 112 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 111 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Peak | 15-10-1995-29-I | 10/23/15 17:16 | Aqueous | GC 44 | 10/29/15 | 10/31/15 02:43 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 99 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 97 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Peak | 15-10-1995-30-I | 10/23/15 17:06 | Aqueous | GC 44 | 10/29/15 | 10/31/15 02:58 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 108 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 103 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Peak | 15-10-1995-31-I | 10/23/15 17:34 | Aqueous | GC 44 | 10/29/15 | 10/31/15 03:12 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 103 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 97 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Peak | 15-10-1995-33-I | 10/23/15 17:25 | Aqueous | GC 44 | 10/29/15 | 10/31/15 03:26 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 110 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 109 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 270 Peak | 15-10-1995-34-I | 10/23/15 17:45 | Aqueous | GC 44 | 10/29/15 | 10/31/15 03:40 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|--------------------|---------------|-----------|------------|-----------|-------------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 108 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 109 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-851 | N/A | Aqueous | GC 44 | 10/29/15 | 10/30/15 13:57 | 151029L12 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 100 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 106 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-529-852 | N/A | Aqueous | GC 44 | 10/29/15 | 10/30/15 14:40 | 151029L13 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------------|--------|------|-------|------|------------|
| Alpha-BHC | ND | 0.10 | 0.028 | 1.00 | |
| Gamma-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Beta-BHC | ND | 0.10 | 0.030 | 1.00 | |
| Heptachlor | ND | 0.10 | 0.026 | 1.00 | |
| Delta-BHC | ND | 0.10 | 0.029 | 1.00 | |
| Aldrin | ND | 0.10 | 0.027 | 1.00 | |
| Heptachlor Epoxide | ND | 0.10 | 0.025 | 1.00 | |
| Endosulfan I | ND | 0.10 | 0.028 | 1.00 | |
| Dieldrin | ND | 0.10 | 0.029 | 1.00 | |
| 4,4'-DDE | ND | 0.10 | 0.027 | 1.00 | |
| Endrin | ND | 0.10 | 0.031 | 1.00 | |
| Endrin Aldehyde | ND | 0.10 | 0.026 | 1.00 | |
| 4,4'-DDD | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan II | ND | 0.10 | 0.027 | 1.00 | |
| 4,4'-DDT | ND | 0.10 | 0.027 | 1.00 | |
| Endosulfan Sulfate | ND | 0.10 | 0.029 | 1.00 | |
| Methoxychlor | ND | 0.10 | 0.025 | 1.00 | |
| Chlordane | ND | 1.0 | 0.33 | 1.00 | |
| Toxaphene | ND | 2.0 | 0.59 | 1.00 | |
| Endrin Ketone | ND | 0.10 | 0.024 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 107 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 93 | 50-135 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-10-1995-1-I | 10/23/15 15:01 | Aqueous | GC 31 | 10/29/15 | 10/30/15 15:10 | 151029L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 111 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 87 | 50-135 | |

| HCS 270 Peak | 15-10-1995-2-I | 10/23/15 15:16 | Aqueous | GC 31 | 10/29/15 | 10/30/15 15:29 | 151029L14 |
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|--------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 116 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 94 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Trail | 15-10-1995-3-I | 10/23/15 16:14 | Aqueous | GC 31 | 10/29/15 | 10/30/15 15:48 | 151029L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 109 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 85 | 50-135 | |

| | | | | | | | |
|---------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
| HCS 240 Trail | 15-10-1995-4-I | 10/23/15 16:29 | Aqueous | GC 31 | 10/29/15 | 10/30/15 16:07 | 151029L14 |
|---------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 116 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 85 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Trail | 15-10-1995-5-I | 10/23/15 17:09 | Aqueous | GC 31 | 10/29/15 | 10/30/15 16:26 | 151029L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 113 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 87 | 50-135 | |

| HCS 260 Trail | 15-10-1995-6-I | 10/23/15 16:50 | Aqueous | GC 31 | 10/29/15 | 10/30/15 16:45 | 151029L14 |
|---------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|---------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 145 | 50-135 | 2,7 |
| 2,4,5,6-Tetrachloro-m-Xylene | 110 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-10-1995-7-I | 10/23/15 17:34 | Aqueous | GC 31 | 10/29/15 | 10/30/15 17:04 | 151029L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 118 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 95 | 50-135 | |

| | | | | | | | |
|-----------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
| FDHCS 260 Trail | 15-10-1995-8-I | 10/23/15 16:50 | Aqueous | GC 31 | 10/29/15 | 10/30/15 17:23 | 151029L14 |
|-----------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 114 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 81 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|-----------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| FDHCS 270 Trail | 15-10-1995-9-I | 10/23/15 17:34 | Aqueous | GC 31 | 10/29/15 | 10/30/15 17:42 | 151029L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 114 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 83 | 50-135 | |

| | | | | | | | |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 210 Trail | 15-10-1995-11-I | 10/23/15 20:35 | Aqueous | GC 31 | 10/29/15 | 10/30/15 18:01 | 151029L14 |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 102 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 67 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Trail | 15-10-1995-12-I | 10/23/15 21:01 | Aqueous | GC 31 | 10/29/15 | 10/30/15 18:20 | 151029L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 113 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 84 | 50-135 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Trail | 15-10-1995-13-I | 10/23/15 21:21 | Aqueous | GC 31 | 10/29/15 | 10/30/15 18:39 | 151029L14 |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 124 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 83 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-10-1995-14-I | 10/23/15 20:51 | Aqueous | GC 31 | 10/29/15 | 10/30/15 18:58 | 151029L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 126 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 85 | 50-135 | |

| HSM 250 Trail | 15-10-1995-15-I | 10/23/15 21:06 | Aqueous | GC 31 | 10/29/15 | 10/30/15 19:18 | 151029L14 |
|---------------|-----------------|----------------|---------|-------|----------|----------------|-----------|
|---------------|-----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 117 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 81 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Trail | 15-10-1995-16-I | 10/23/15 21:35 | Aqueous | GC 31 | 10/29/15 | 10/30/15 19:56 | 151029L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 122 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 91 | 50-135 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 270 Trail | 15-10-1995-17-I | 10/23/15 21:55 | Aqueous | GC 31 | 10/29/15 | 10/30/15 20:15 | 151029L14 |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 120 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 92 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 210 Trail | 15-10-1995-18-I | 10/23/15 20:35 | Aqueous | GC 31 | 10/29/15 | 10/30/15 20:34 | 151029L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 92 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 70 | 50-135 | |

| | | | | | | | |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 230 Trail | 15-10-1995-19-I | 10/23/15 21:01 | Aqueous | GC 31 | 10/29/15 | 10/30/15 20:53 | 151029L14 |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 111 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 89 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 231 Trail | 15-10-1995-20-I | 10/23/15 21:21 | Aqueous | GC 31 | 10/29/15 | 10/30/15 21:12 | 151029L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 124 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 95 | 50-135 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 210 Lead | 15-10-1995-21-I | 10/23/15 15:25 | Aqueous | GC 31 | 10/29/15 | 10/30/15 21:31 | 151029L14 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 125 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 100 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Lead | 15-10-1995-22-I | 10/23/15 15:41 | Aqueous | GC 31 | 10/29/15 | 10/30/15 21:50 | 151029L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 112 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 96 | 50-135 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Lead | 15-10-1995-23-I | 10/23/15 16:02 | Aqueous | GC 31 | 10/29/15 | 10/30/15 22:09 | 151029L15 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 125 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 103 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Lead | 15-10-1995-24-I | 10/23/15 15:30 | Aqueous | GC 31 | 10/29/15 | 10/30/15 22:28 | 151029L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 112 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 95 | 50-135 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 250 Lead | 15-10-1995-25-I | 10/23/15 16:21 | Aqueous | GC 31 | 10/29/15 | 10/30/15 22:47 | 151029L15 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 125 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 97 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Lead | 15-10-1995-26-I | 10/23/15 15:58 | Aqueous | GC 31 | 10/29/15 | 10/30/15 23:06 | 151029L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 122 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 93 | 50-135 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 270 Lead | 15-10-1995-27-I | 10/23/15 16:15 | Aqueous | GC 31 | 10/29/15 | 10/30/15 23:25 | 151029L15 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 128 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 99 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 210 Peak | 15-10-1995-28-I | 10/23/15 17:02 | Aqueous | GC 31 | 10/29/15 | 10/30/15 23:44 | 151029L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 127 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 98 | 50-135 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Peak | 15-10-1995-29-I | 10/23/15 17:16 | Aqueous | GC 31 | 10/29/15 | 10/31/15 00:03 | 151029L15 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 113 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 89 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Peak | 15-10-1995-30-I | 10/23/15 17:06 | Aqueous | GC 31 | 10/29/15 | 10/31/15 00:22 | 151029L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 124 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 97 | 50-135 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Peak | 15-10-1995-31-I | 10/23/15 17:34 | Aqueous | GC 31 | 10/29/15 | 10/31/15 00:41 | 151029L15 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 119 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 95 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Peak | 15-10-1995-33-I | 10/23/15 17:25 | Aqueous | GC 31 | 10/29/15 | 10/31/15 01:00 | 151029L15 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 123 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 100 | 50-135 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 270 Peak | 15-10-1995-34-I | 10/23/15 17:45 | Aqueous | GC 31 | 10/29/15 | 10/31/15 01:19 | 151029L15 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------------|-----------------|-----------------------|-------------------|
| Decachlorobiphenyl | 123 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 102 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-12-533-1101 | N/A | Aqueous | GC 31 | 10/29/15 | 10/30/15 13:53 | 151029L14 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 116 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 86 | 50-135 | |

| | | | | | | | |
|--------------|-----------------|-----|---------|-------|----------|----------------|-----------|
| Method Blank | 099-12-533-1102 | N/A | Aqueous | GC 31 | 10/29/15 | 10/30/15 14:51 | 151029L15 |
|--------------|-----------------|-----|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|--------------|--------|-----|------|------|------------|
| Aroclor-1016 | ND | 1.0 | 0.29 | 1.00 | |
| Aroclor-1221 | ND | 1.0 | 0.28 | 1.00 | |
| Aroclor-1232 | ND | 1.0 | 0.25 | 1.00 | |
| Aroclor-1242 | ND | 1.0 | 0.18 | 1.00 | |
| Aroclor-1248 | ND | 1.0 | 0.20 | 1.00 | |
| Aroclor-1254 | ND | 1.0 | 0.23 | 1.00 | |
| Aroclor-1260 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1262 | ND | 1.0 | 0.26 | 1.00 | |
| Aroclor-1268 | ND | 1.0 | 0.21 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------------|----------|----------------|------------|
| Decachlorobiphenyl | 119 | 50-135 | |
| 2,4,5,6-Tetrachloro-m-Xylene | 83 | 50-135 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-10-1995-1-I | 10/23/15 15:01 | Aqueous | GC 26 | 10/30/15 | 10/31/15 06:51 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 74 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Peak | 15-10-1995-2-I | 10/23/15 15:16 | Aqueous | GC 26 | 10/30/15 | 10/31/15 07:35 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 69 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Trail | 15-10-1995-3-I | 10/23/15 16:14 | Aqueous | GC 26 | 10/30/15 | 10/31/15 08:19 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 77 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-10-1995-4-I | 10/23/15 16:29 | Aqueous | GC 26 | 10/30/15 | 10/31/15 09:03 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 70 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Trail | 15-10-1995-5-I | 10/23/15 17:09 | Aqueous | GC 26 | 10/30/15 | 10/31/15 09:47 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 76 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Trail | 15-10-1995-6-I | 10/23/15 16:50 | Aqueous | GC 26 | 10/30/15 | 10/31/15 10:31 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 77 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-10-1995-7-I | 10/23/15 17:34 | Aqueous | GC 26 | 10/30/15 | 10/31/15 11:16 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 73 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 260 Trail | 15-10-1995-8-I | 10/23/15 16:50 | Aqueous | GC 26 | 10/30/15 | 10/31/15 12:00 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 75 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 270 Trail | 15-10-1995-9-I | 10/23/15 17:34 | Aqueous | GC 26 | 10/30/15 | 10/31/15 12:44 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 74 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-10-1995-11-I | 10/23/15 20:35 | Aqueous | GC 26 | 10/30/15 | 10/31/15 13:28 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 76 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Trail | 15-10-1995-12-I | 10/23/15 21:01 | Aqueous | GC 26 | 10/30/15 | 10/31/15 14:12 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 79 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 231 Trail | 15-10-1995-13-I | 10/23/15 21:21 | Aqueous | GC 26 | 10/30/15 | 10/31/15 14:56 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 76 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-10-1995-14-I | 10/23/15 20:51 | Aqueous | GC 26 | 10/30/15 | 10/31/15 15:40 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 79 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-10-1995-15-I | 10/23/15 21:06 | Aqueous | GC 26 | 10/30/15 | 10/31/15 16:24 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 72 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Trail | 15-10-1995-16-I | 10/23/15 21:35 | Aqueous | GC 26 | 10/30/15 | 10/31/15 17:09 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 67 | 30-130 | |

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RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Trail | 15-10-1995-17-I | 10/23/15 21:55 | Aqueous | GC 26 | 10/30/15 | 10/31/15 18:37 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 77 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 210 Trail | 15-10-1995-18-I | 10/23/15 20:35 | Aqueous | GC 26 | 10/30/15 | 10/31/15 19:21 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 81 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 230 Trail | 15-10-1995-19-I | 10/23/15 21:01 | Aqueous | GC 26 | 10/30/15 | 10/31/15 20:05 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 81 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 231 Trail | 15-10-1995-20-I | 10/23/15 21:21 | Aqueous | GC 26 | 10/30/15 | 10/31/15 20:49 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 70 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 210 Lead | 15-10-1995-21-I | 10/23/15 15:25 | Aqueous | GC 26 | 10/30/15 | 10/31/15 21:33 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 74 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Lead | 15-10-1995-22-I | 10/23/15 15:41 | Aqueous | GC 26 | 10/30/15 | 10/31/15 22:17 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 71 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 231 Lead | 15-10-1995-23-I | 10/23/15 16:02 | Aqueous | GC 26 | 10/30/15 | 10/31/15 23:02 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 78 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Lead | 15-10-1995-24-I | 10/23/15 15:30 | Aqueous | GC 26 | 10/30/15 | 10/31/15 23:46 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 74 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 250 Lead | 15-10-1995-25-I | 10/23/15 16:21 | Aqueous | GC 26 | 10/30/15 | 11/01/15 00:30 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 73 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Lead | 15-10-1995-26-I | 10/23/15 15:58 | Aqueous | GC 26 | 10/30/15 | 11/01/15 01:14 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 74 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 270 Lead | 15-10-1995-27-I | 10/23/15 16:15 | Aqueous | GC 26 | 10/30/15 | 11/01/15 01:58 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 66 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 210 Peak | 15-10-1995-28-I | 10/23/15 17:02 | Aqueous | GC 26 | 10/30/15 | 11/01/15 02:42 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 78 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Peak | 15-10-1995-29-I | 10/23/15 17:16 | Aqueous | GC 26 | 10/30/15 | 11/01/15 03:26 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 76 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 231 Peak | 15-10-1995-30-I | 10/23/15 17:06 | Aqueous | GC 26 | 10/30/15 | 11/01/15 04:10 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 72 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 240 Peak | 15-10-1995-31-I | 10/23/15 17:34 | Aqueous | GC 26 | 10/30/15 | 11/01/15 04:54 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 71 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 260 Peak | 15-10-1995-33-I | 10/23/15 17:25 | Aqueous | GC 26 | 10/30/15 | 11/01/15 06:22 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 68 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|--------------|-----------------|---------------------------|------------------|
| HSM 270 Peak | 15-10-1995-34-I | 10/23/15 17:45 | Aqueous | GC 26 | 10/30/15 | 11/01/15 07:07 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------|---------------|-----------|------------|-----------|-------------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------|-----------------|-----------------------|-------------------|
| Tributylphosphate | 74 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-121 | N/A | Aqueous | GC 26 | 10/30/15 | 10/30/15 19:49 | 151030L02 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 68 | 30-130 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A
Units: mg/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-963-122 | N/A | Aqueous | GC 26 | 10/30/15 | 10/30/15 22:02 | 151030L03 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------|--------|--------|--------|------|------------|
| Azinphos Methyl | ND | 0.0050 | 0.0029 | 1.00 | |
| Bolstar | ND | 0.0050 | 0.0029 | 1.00 | |
| Chlorpyrifos | ND | 0.0050 | 0.0024 | 1.00 | |
| Coumaphos | ND | 0.0050 | 0.0024 | 1.00 | |
| Diazinon | ND | 0.0050 | 0.0029 | 1.00 | |
| Dichlorvos | ND | 0.0050 | 0.0036 | 1.00 | |
| Disulfoton | ND | 0.010 | 0.0026 | 1.00 | |
| Ethoprop | ND | 0.0050 | 0.0025 | 1.00 | |
| Fensulfothion | ND | 0.0050 | 0.0029 | 1.00 | |
| Fenthion | ND | 0.0050 | 0.0026 | 1.00 | |
| Merphos | ND | 0.0050 | 0.0026 | 1.00 | |
| Methyl Parathion | ND | 0.0050 | 0.0030 | 1.00 | |
| Mevinphos | ND | 0.0050 | 0.0027 | 1.00 | |
| Naled | ND | 0.040 | 0.019 | 1.00 | |
| Phorate | ND | 0.0050 | 0.0025 | 1.00 | |
| Ronnel | ND | 0.0050 | 0.0032 | 1.00 | |
| Stirophos | ND | 0.020 | 0.0088 | 1.00 | |
| Tokuthion | ND | 0.0050 | 0.0028 | 1.00 | |
| Trichloronate | ND | 0.0050 | 0.0021 | 1.00 | |
| Demeton-o/s | ND | 0.0050 | 0.0028 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------|----------|----------------|------------|
| Tributylphosphate | 72 | 30-130 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-10-1995-1-I | 10/23/15 15:01 | Aqueous | GC 40 | 10/30/15 | 11/05/15 22:04 | 151030L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 92 | 0-123 | |

| HCS 270 Peak | 15-10-1995-2-I | 10/23/15 15:16 | Aqueous | GC 40 | 10/30/15 | 11/05/15 22:27 | 151030L19 |
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|
|--------------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 82 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Trail | 15-10-1995-3-I | 10/23/15 16:14 | Aqueous | GC 40 | 10/30/15 | 11/05/15 22:50 | 151030L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 134 | 0-123 | 2,7 |

| HCS 240 Trail | 15-10-1995-4-I | 10/23/15 16:29 | Aqueous | GC 40 | 10/30/15 | 11/05/15 23:13 | 151030L19 |
|---------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
|---------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 101 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Trail | 15-10-1995-5-I | 10/23/15 17:09 | Aqueous | GC 40 | 10/30/15 | 11/05/15 23:36 | 151030L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 122 | 0-123 | |

| | | | | | | | |
|---------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|
| HCS 260 Trail | 15-10-1995-6-I | 10/23/15 16:50 | Aqueous | GC 40 | 10/30/15 | 11/05/15 23:59 | 151030L19 |
|---------------|----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 113 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-10-1995-7-I | 10/23/15 17:34 | Aqueous | GC 40 | 10/30/15 | 11/06/15 00:22 | 151030L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 112 | 0-123 | |

| | | | | | | | |
|-----------------|----------------|----------------|---------|-------|----------|----------------|-----------|
| FDHCS 260 Trail | 15-10-1995-8-I | 10/23/15 16:50 | Aqueous | GC 40 | 10/30/15 | 11/06/15 00:45 | 151030L19 |
|-----------------|----------------|----------------|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 105 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 270 Trail | 15-10-1995-9-I | 10/23/15 17:34 | Aqueous | GC 40 | 10/30/15 | 11/06/15 01:08 | 151030L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 100 | 0-123 | |

| HSM 210 Trail | 15-10-1995-11-I | 10/23/15 20:35 | Aqueous | GC 40 | 10/30/15 | 11/06/15 01:31 | 151030L19 |
|---------------|-----------------|-------------------|---------|-------|----------|-------------------|-----------|
|---------------|-----------------|-------------------|---------|-------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 115 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Trail | 15-10-1995-12-I | 10/23/15 21:01 | Aqueous | GC 40 | 10/30/15 | 11/06/15 01:54 | 151030L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 123 | 0-123 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Trail | 15-10-1995-13-I | 10/23/15 21:21 | Aqueous | GC 40 | 10/30/15 | 11/06/15 02:17 | 151030L19 |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 110 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Trail | 15-10-1995-14-I | 10/23/15 20:51 | Aqueous | GC 40 | 10/30/15 | 11/06/15 02:40 | 151030L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 106 | 0-123 | |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 250 Trail | 15-10-1995-15-I | 10/23/15 21:06 | Aqueous | GC 40 | 10/30/15 | 11/06/15 03:03 | 151030L19 |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 116 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Trail | 15-10-1995-16-I | 10/23/15 21:35 | Aqueous | GC 40 | 10/30/15 | 11/06/15 03:49 | 151030L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 125 | 0-123 | 2,7 |

| | | | | | | | |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 270 Trail | 15-10-1995-17-I | 10/23/15 21:55 | Aqueous | GC 40 | 10/30/15 | 11/06/15 04:13 | 151030L19 |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 117 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 210 Trail | 15-10-1995-18-I | 10/23/15 20:35 | Aqueous | GC 40 | 10/30/15 | 11/06/15 04:35 | 151030L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 145 | 0-123 | 2,7 |

| | | | | | | | |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 230 Trail | 15-10-1995-19-I | 10/23/15 21:01 | Aqueous | GC 40 | 10/30/15 | 11/06/15 04:58 | 151030L20 |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 113 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| FDHSM 231 Trail | 15-10-1995-20-I | 10/23/15 21:21 | Aqueous | GC 40 | 10/30/15 | 11/06/15 05:22 | 151030L20 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 112 | 0-123 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 210 Lead | 15-10-1995-21-I | 10/23/15 15:25 | Aqueous | GC 40 | 10/30/15 | 11/06/15 05:44 | 151030L20 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 98 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Lead | 15-10-1995-22-I | 10/23/15 15:41 | Aqueous | GC 40 | 10/30/15 | 11/06/15 06:08 | 151030L20 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 113 | 0-123 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Lead | 15-10-1995-23-I | 10/23/15 16:02 | Aqueous | GC 40 | 10/30/15 | 11/06/15 06:31 | 151030L20 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 106 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Lead | 15-10-1995-24-I | 10/23/15 15:30 | Aqueous | GC 40 | 10/30/15 | 11/06/15 06:54 | 151030L20 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 116 | 0-123 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 250 Lead | 15-10-1995-25-I | 10/23/15 16:21 | Aqueous | GC 40 | 10/30/15 | 11/06/15 07:17 | 151030L20 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 52 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Lead | 15-10-1995-26-I | 10/23/15 15:58 | Aqueous | GC 40 | 10/30/15 | 11/06/15 07:40 | 151030L20 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 112 | 0-123 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 270 Lead | 15-10-1995-27-I | 10/23/15 16:15 | Aqueous | GC 40 | 10/30/15 | 11/06/15 08:03 | 151030L20 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 105 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 210 Peak | 15-10-1995-28-I | 10/23/15 17:02 | Aqueous | GC 40 | 10/30/15 | 11/06/15 08:26 | 151030L20 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPD | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 84 | 0-123 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 230 Peak | 15-10-1995-29-I | 10/23/15 17:16 | Aqueous | GC 40 | 10/30/15 | 11/06/15 08:49 | 151030L20 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCPD | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 99 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 231 Peak | 15-10-1995-30-I | 10/23/15 17:06 | Aqueous | GC 40 | 10/30/15 | 11/06/15 09:12 | 151030L20 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 109 | 0-123 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 240 Peak | 15-10-1995-31-I | 10/23/15 17:34 | Aqueous | GC 40 | 10/30/15 | 11/06/15 09:35 | 151030L20 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 116 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 260 Peak | 15-10-1995-33-I | 10/23/15 17:25 | Aqueous | GC 40 | 10/30/15 | 11/06/15 09:58 | 151030L20 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 100 | 0-123 | |

| | | | | | | | |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|
| HSM 270 Peak | 15-10-1995-34-I | 10/23/15 17:45 | Aqueous | GC 40 | 10/30/15 | 11/06/15 10:21 | 151030L20 |
|---------------------|------------------------|-----------------------|----------------|--------------|-----------------|-----------------------|------------------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-------------------|---------------|-----------|------------|-----------|-------------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|-------------------------------|-----------------|-----------------------|-------------------|
| 2,4-Dichlorophenylacetic acid | 128 | 0-123 | 2,7 |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-034-667 | N/A | Aqueous | GC 40 | 10/30/15 | 11/05/15 20:32 | 151030L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 91 | 0-123 | |

| | | | | | | | |
|--------------|----------------|-----|---------|-------|----------|----------------|-----------|
| Method Blank | 095-01-034-668 | N/A | Aqueous | GC 40 | 10/30/15 | 11/05/15 21:41 | 151030L20 |
|--------------|----------------|-----|---------|-------|----------|----------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-------------------|--------|------|------|------|------------|
| Dalapon | ND | 13 | 3.7 | 1.00 | |
| Dicamba | ND | 0.50 | 0.17 | 1.00 | |
| MCP | ND | 500 | 170 | 1.00 | |
| MCPA | ND | 500 | 170 | 1.00 | |
| Dichlorprop | ND | 5.0 | 1.8 | 1.00 | |
| 2,4-D | ND | 5.0 | 1.8 | 1.00 | |
| 2,4,5-TP (Silvex) | ND | 0.50 | 0.22 | 1.00 | |
| 2,4,5-T | ND | 0.50 | 0.18 | 1.00 | |
| 2,4-DB | ND | 5.0 | 1.5 | 1.00 | |
| Dinoseb | ND | 2.5 | 0.95 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|-------------------------------|----------|----------------|------------|
| 2,4-Dichlorophenylacetic acid | 85 | 0-123 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 1 of 102

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-10-1995-1-H | 10/23/15 15:01 | Aqueous | GC/MS SS | 10/30/15 | 10/31/15 17:19 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 84 | 33-120 | |
| 2-Fluorophenol | 55 | 24-120 | |
| Nitrobenzene-d5 | 92 | 38-120 | |
| p-Terphenyl-d14 | 90 | 41-137 | |
| Phenol-d6 | 34 | 16-120 | |
| 2,4,6-Tribromophenol | 104 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Peak | 15-10-1995-2-H | 10/23/15 15:16 | Aqueous | GC/MS SS | 10/30/15 | 11/02/15 11:19 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 68 | 33-120 | |
| 2-Fluorophenol | 43 | 24-120 | |
| Nitrobenzene-d5 | 69 | 38-120 | |
| p-Terphenyl-d14 | 69 | 41-137 | |
| Phenol-d6 | 27 | 16-120 | |
| 2,4,6-Tribromophenol | 66 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 7 of 102

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Trail | 15-10-1995-3-H | 10/23/15 16:14 | Aqueous | GC/MS SS | 10/30/15 | 11/02/15 11:39 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 91 | 33-120 | |
| 2-Fluorophenol | 55 | 24-120 | |
| Nitrobenzene-d5 | 95 | 38-120 | |
| p-Terphenyl-d14 | 95 | 41-137 | |
| Phenol-d6 | 35 | 16-120 | |
| 2,4,6-Tribromophenol | 90 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-10-1995-4-G | 10/23/15 16:29 | Aqueous | GC/MS SS | 10/30/15 | 11/02/15 11:58 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 72 | 33-120 | |
| 2-Fluorophenol | 41 | 24-120 | |
| Nitrobenzene-d5 | 75 | 38-120 | |
| p-Terphenyl-d14 | 70 | 41-137 | |
| Phenol-d6 | 26 | 16-120 | |
| 2,4,6-Tribromophenol | 62 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Trail | 15-10-1995-5-H | 10/23/15 17:09 | Aqueous | GC/MS SS | 10/30/15 | 10/31/15 18:38 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 87 | 33-120 | |
| 2-Fluorophenol | 57 | 24-120 | |
| Nitrobenzene-d5 | 97 | 38-120 | |
| p-Terphenyl-d14 | 91 | 41-137 | |
| Phenol-d6 | 36 | 16-120 | |
| 2,4,6-Tribromophenol | 104 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Trail | 15-10-1995-6-H | 10/23/15 16:50 | Aqueous | GC/MS SS | 10/30/15 | 10/31/15 18:58 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 74 | 33-120 | |
| 2-Fluorophenol | 51 | 24-120 | |
| Nitrobenzene-d5 | 83 | 38-120 | |
| p-Terphenyl-d14 | 78 | 41-137 | |
| Phenol-d6 | 32 | 16-120 | |
| 2,4,6-Tribromophenol | 90 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-10-1995-7-H | 10/23/15 17:34 | Aqueous | GC/MS SS | 10/30/15 | 10/31/15 19:17 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 87 | 33-120 | |
| 2-Fluorophenol | 60 | 24-120 | |
| Nitrobenzene-d5 | 99 | 38-120 | |
| p-Terphenyl-d14 | 89 | 41-137 | |
| Phenol-d6 | 38 | 16-120 | |
| 2,4,6-Tribromophenol | 103 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|-----------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| FDHCS 260 Trail | 15-10-1995-8-H | 10/23/15 16:50 | Aqueous | GC/MS SS | 10/30/15 | 10/31/15 19:37 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 80 | 33-120 | |
| 2-Fluorophenol | 56 | 24-120 | |
| Nitrobenzene-d5 | 89 | 38-120 | |
| p-Terphenyl-d14 | 81 | 41-137 | |
| Phenol-d6 | 34 | 16-120 | |
| 2,4,6-Tribromophenol | 93 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 270 Trail | 15-10-1995-9-H | 10/23/15 17:34 | Aqueous | GC/MS SS | 10/30/15 | 10/31/15 19:57 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 82 | 33-120 | |
| 2-Fluorophenol | 54 | 24-120 | |
| Nitrobenzene-d5 | 91 | 38-120 | |
| p-Terphenyl-d14 | 82 | 41-137 | |
| Phenol-d6 | 33 | 16-120 | |
| 2,4,6-Tribromophenol | 94 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 210 Trail | 15-10-1995-11-H | 10/23/15 20:35 | Aqueous | GC/MS SS | 10/30/15 | 10/31/15 20:17 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.9 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 9.9 | 2.9 | 1.00 | |
| Aniline | ND | 9.9 | 1.5 | 1.00 | |
| Anthracene | ND | 9.9 | 3.0 | 1.00 | |
| Azobenzene | ND | 9.9 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.9 | 4.6 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.9 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.9 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.9 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.9 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.9 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.9 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.9 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.9 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.9 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.9 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.9 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 9.9 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.9 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.9 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.9 | 2.6 | 1.00 | |
| Chrysene | ND | 9.9 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.9 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.9 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.9 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.9 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.9 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.9 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.9 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.9 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.9 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.9 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.9 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.9 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.9 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.9 | 3.1 | 1.00 | |
| Fluorene | ND | 9.9 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.9 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 9.9 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 9.9 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.9 | 2.1 | 1.00 | |
| Isophorone | ND | 9.9 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.9 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.9 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 9.9 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 9.9 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.9 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.9 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.9 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.9 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.9 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.9 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.9 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 9.9 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 9.9 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 9.9 | 4.6 | 1.00 | |
| Phenanthrene | ND | 9.9 | 2.9 | 1.00 | |
| Phenol | ND | 9.9 | 2.0 | 1.00 | |
| Pyrene | ND | 9.9 | 2.9 | 1.00 | |
| Pyridine | ND | 9.9 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.9 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.9 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.9 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 88 | 33-120 | |
| 2-Fluorophenol | 59 | 24-120 | |
| Nitrobenzene-d5 | 96 | 38-120 | |
| p-Terphenyl-d14 | 89 | 41-137 | |
| Phenol-d6 | 38 | 16-120 | |
| 2,4,6-Tribromophenol | 101 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Trail | 15-10-1995-12-H | 10/23/15 21:01 | Aqueous | GC/MS SS | 10/30/15 | 10/31/15 20:36 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 77 | 33-120 | |
| 2-Fluorophenol | 53 | 24-120 | |
| Nitrobenzene-d5 | 88 | 38-120 | |
| p-Terphenyl-d14 | 78 | 41-137 | |
| Phenol-d6 | 40 | 16-120 | |
| 2,4,6-Tribromophenol | 89 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 231 Trail | 15-10-1995-13-H | 10/23/15 21:21 | Aqueous | GC/MS SS | 10/30/15 | 10/31/15 20:56 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 35 of 102

| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 85 | 33-120 | |
| 2-Fluorophenol | 57 | 24-120 | |
| Nitrobenzene-d5 | 94 | 38-120 | |
| p-Terphenyl-d14 | 86 | 41-137 | |
| Phenol-d6 | 36 | 16-120 | |
| 2,4,6-Tribromophenol | 98 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-10-1995-14-H | 10/23/15 20:51 | Aqueous | GC/MS SS | 10/30/15 | 10/31/15 21:16 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 67 | 33-120 | |
| 2-Fluorophenol | 50 | 24-120 | |
| Nitrobenzene-d5 | 79 | 38-120 | |
| p-Terphenyl-d14 | 72 | 41-137 | |
| Phenol-d6 | 32 | 16-120 | |
| 2,4,6-Tribromophenol | 78 | 27-159 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-10-1995-15-H | 10/23/15 21:06 | Aqueous | GC/MS SS | 10/30/15 | 11/02/15 12:18 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 58 | 33-120 | |
| 2-Fluorophenol | 40 | 24-120 | |
| Nitrobenzene-d5 | 62 | 38-120 | |
| p-Terphenyl-d14 | 60 | 41-137 | |
| Phenol-d6 | 24 | 16-120 | |
| 2,4,6-Tribromophenol | 53 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Trail | 15-10-1995-16-H | 10/23/15 21:35 | Aqueous | GC/MS SS | 10/30/15 | 11/02/15 12:38 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 71 | 33-120 | |
| 2-Fluorophenol | 48 | 24-120 | |
| Nitrobenzene-d5 | 77 | 38-120 | |
| p-Terphenyl-d14 | 73 | 41-137 | |
| Phenol-d6 | 29 | 16-120 | |
| 2,4,6-Tribromophenol | 68 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Trail | 15-10-1995-17-H | 10/23/15 21:55 | Aqueous | GC/MS SS | 10/30/15 | 11/02/15 12:57 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 84 | 33-120 | |
| 2-Fluorophenol | 54 | 24-120 | |
| Nitrobenzene-d5 | 89 | 38-120 | |
| p-Terphenyl-d14 | 86 | 41-137 | |
| Phenol-d6 | 33 | 16-120 | |
| 2,4,6-Tribromophenol | 80 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| FDHSM 210 Trail | 15-10-1995-18-H | 10/23/15 20:35 | Aqueous | GC/MS SS | 10/30/15 | 11/02/15 13:17 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 86 | 33-120 | |
| 2-Fluorophenol | 57 | 24-120 | |
| Nitrobenzene-d5 | 94 | 38-120 | |
| p-Terphenyl-d14 | 83 | 41-137 | |
| Phenol-d6 | 33 | 16-120 | |
| 2,4,6-Tribromophenol | 81 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| FDHSM 230 Trail | 15-10-1995-19-H | 10/23/15 21:01 | Aqueous | GC/MS SS | 10/30/15 | 11/02/15 13:37 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 54 of 102

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 80 | 33-120 | |
| 2-Fluorophenol | 59 | 24-120 | |
| Nitrobenzene-d5 | 90 | 38-120 | |
| p-Terphenyl-d14 | 88 | 41-137 | |
| Phenol-d6 | 38 | 16-120 | |
| 2,4,6-Tribromophenol | 81 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|------------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| FDHSM 231 Trail | 15-10-1995-20-H | 10/23/15 21:21 | Aqueous | GC/MS SS | 10/30/15 | 11/02/15 13:57 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 80 | 33-120 | |
| 2-Fluorophenol | 51 | 24-120 | |
| Nitrobenzene-d5 | 86 | 38-120 | |
| p-Terphenyl-d14 | 83 | 41-137 | |
| Phenol-d6 | 31 | 16-120 | |
| 2,4,6-Tribromophenol | 74 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 210 Lead | 15-10-1995-21-H | 10/23/15 15:25 | Aqueous | GC/MS SS | 10/30/15 | 11/02/15 14:16 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 83 | 33-120 | |
| 2-Fluorophenol | 55 | 24-120 | |
| Nitrobenzene-d5 | 91 | 38-120 | |
| p-Terphenyl-d14 | 86 | 41-137 | |
| Phenol-d6 | 34 | 16-120 | |
| 2,4,6-Tribromophenol | 78 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 230 Lead | 15-10-1995-22-H | 10/23/15 15:41 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 17:15 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 80 | 33-120 | |
| 2-Fluorophenol | 57 | 24-120 | |
| Nitrobenzene-d5 | 82 | 38-120 | |
| p-Terphenyl-d14 | 78 | 41-137 | |
| Phenol-d6 | 35 | 16-120 | |
| 2,4,6-Tribromophenol | 96 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|-----------------|-----------------|---------------------------|------------------|
| HSM 231 Lead | 15-10-1995-23-H | 10/23/15 16:02 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 17:35 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 73 | 33-120 | |
| 2-Fluorophenol | 51 | 24-120 | |
| Nitrobenzene-d5 | 71 | 38-120 | |
| p-Terphenyl-d14 | 77 | 41-137 | |
| Phenol-d6 | 32 | 16-120 | |
| 2,4,6-Tribromophenol | 89 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 240 Lead | 15-10-1995-24-H | 10/23/15 15:30 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 17:54 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 62 | 33-120 | |
| 2-Fluorophenol | 47 | 24-120 | |
| Nitrobenzene-d5 | 60 | 38-120 | |
| p-Terphenyl-d14 | 66 | 41-137 | |
| Phenol-d6 | 29 | 16-120 | |
| 2,4,6-Tribromophenol | 79 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|-----------------|-----------------|---------------------------|------------------|
| HSM 250 Lead | 15-10-1995-25-H | 10/23/15 16:21 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 18:12 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.8 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 9.8 | 2.8 | 1.00 | |
| Aniline | ND | 9.8 | 1.5 | 1.00 | |
| Anthracene | ND | 9.8 | 3.0 | 1.00 | |
| Azobenzene | ND | 9.8 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.4 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.8 | 4.6 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.8 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.8 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.8 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.8 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.8 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.8 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.8 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.8 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.8 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.8 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.8 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.8 | 2.6 | 1.00 | |
| Chrysene | ND | 9.8 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.8 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.8 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.8 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.8 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.8 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.8 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.8 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.8 | 3.0 | 1.00 | |
| Fluorene | ND | 9.8 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.8 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.8 | 1.00 | |
| Hexachloroethane | ND | 9.8 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.8 | 2.1 | 1.00 | |
| Isophorone | ND | 9.8 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.8 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.8 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.8 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.8 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 9.8 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 9.8 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.8 | 4.6 | 1.00 | |
| Phenanthrene | ND | 9.8 | 2.9 | 1.00 | |
| Phenol | ND | 9.8 | 2.0 | 1.00 | |
| Pyrene | ND | 9.8 | 2.9 | 1.00 | |
| Pyridine | ND | 9.8 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 63 | 33-120 | |
| 2-Fluorophenol | 43 | 24-120 | |
| Nitrobenzene-d5 | 60 | 38-120 | |
| p-Terphenyl-d14 | 70 | 41-137 | |
| Phenol-d6 | 27 | 16-120 | |
| 2,4,6-Tribromophenol | 81 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 73 of 102

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 260 Lead | 15-10-1995-26-H | 10/23/15 15:58 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 18:31 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 67 | 33-120 | |
| 2-Fluorophenol | 43 | 24-120 | |
| Nitrobenzene-d5 | 63 | 38-120 | |
| p-Terphenyl-d14 | 70 | 41-137 | |
| Phenol-d6 | 26 | 16-120 | |
| 2,4,6-Tribromophenol | 86 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Lead | 15-10-1995-27-H | 10/23/15 16:15 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 18:50 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|-----|-----|------|------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 77 | 33-120 | |
| 2-Fluorophenol | 51 | 24-120 | |
| Nitrobenzene-d5 | 72 | 38-120 | |
| p-Terphenyl-d14 | 83 | 41-137 | |
| Phenol-d6 | 31 | 16-120 | |
| 2,4,6-Tribromophenol | 99 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 210 Peak | 15-10-1995-28-H | 10/23/15 17:02 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 19:09 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.8 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 9.8 | 2.8 | 1.00 | |
| Aniline | ND | 9.8 | 1.5 | 1.00 | |
| Anthracene | ND | 9.8 | 3.0 | 1.00 | |
| Azobenzene | ND | 9.8 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.4 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.8 | 4.6 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.8 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.8 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.8 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.8 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.8 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.8 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.8 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.8 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.8 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.8 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.8 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.8 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.8 | 2.6 | 1.00 | |
| Chrysene | ND | 9.8 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.8 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.8 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.8 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.8 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.8 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.8 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.8 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.8 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.8 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.8 | 3.0 | 1.00 | |
| Fluorene | ND | 9.8 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.8 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.8 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.8 | 1.00 | |
| Hexachloroethane | ND | 9.8 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.8 | 2.1 | 1.00 | |
| Isophorone | ND | 9.8 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 9.8 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.8 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.8 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.8 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.8 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.8 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.8 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 9.8 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 9.8 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.8 | 4.6 | 1.00 | |
| Phenanthrene | ND | 9.8 | 2.9 | 1.00 | |
| Phenol | ND | 9.8 | 2.0 | 1.00 | |
| Pyrene | ND | 9.8 | 2.9 | 1.00 | |
| Pyridine | ND | 9.8 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.8 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.8 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 78 | 33-120 | |
| 2-Fluorophenol | 62 | 24-120 | |
| Nitrobenzene-d5 | 79 | 38-120 | |
| p-Terphenyl-d14 | 79 | 41-137 | |
| Phenol-d6 | 51 | 16-120 | |
| 2,4,6-Tribromophenol | 97 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 230 Peak | 15-10-1995-29-H | 10/23/15 17:16 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 19:27 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 74 | 33-120 | |
| 2-Fluorophenol | 53 | 24-120 | |
| Nitrobenzene-d5 | 73 | 38-120 | |
| p-Terphenyl-d14 | 76 | 41-137 | |
| Phenol-d6 | 33 | 16-120 | |
| 2,4,6-Tribromophenol | 91 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|-----------------|-----------------|---------------------------|------------------|
| HSM 231 Peak | 15-10-1995-30-H | 10/23/15 17:06 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 19:46 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.7 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.7 | 2.8 | 1.00 | |
| Aniline | ND | 9.7 | 1.5 | 1.00 | |
| Anthracene | ND | 9.7 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.7 | 2.6 | 1.00 | |
| Benzidine | ND | 49 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.7 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.7 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.7 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.7 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.7 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 49 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.7 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.7 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.7 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.7 | 3.1 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.7 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.7 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.7 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.7 | 2.6 | 1.00 | |
| Chrysene | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.7 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.7 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.7 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 9.7 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.7 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.7 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.7 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 49 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 49 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.7 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.7 | 3.0 | 1.00 | |
| Fluorene | ND | 9.7 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.7 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.7 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.7 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.7 | 2.1 | 1.00 | |
| Isophorone | ND | 9.7 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.7 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.7 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.7 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.7 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.7 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.7 | 2.7 | 1.00 | |
| Naphthalene | ND | 9.7 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.7 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.7 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.7 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.7 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.7 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.7 | 2.8 | 1.00 | |
| Phenol | ND | 9.7 | 2.0 | 1.00 | |
| Pyrene | ND | 9.7 | 2.9 | 1.00 | |
| Pyridine | ND | 9.7 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.7 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.7 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 73 | 33-120 | |
| 2-Fluorophenol | 52 | 24-120 | |
| Nitrobenzene-d5 | 72 | 38-120 | |
| p-Terphenyl-d14 | 75 | 41-137 | |
| Phenol-d6 | 35 | 16-120 | |
| 2,4,6-Tribromophenol | 88 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|------------------|
| HSM 240 Peak | 15-10-1995-31-H | 10/23/15 17:34 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 20:05 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 76 | 33-120 | |
| 2-Fluorophenol | 50 | 24-120 | |
| Nitrobenzene-d5 | 75 | 38-120 | |
| p-Terphenyl-d14 | 75 | 41-137 | |
| Phenol-d6 | 30 | 16-120 | |
| 2,4,6-Tribromophenol | 91 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|-----------------|-----------------|---------------------------|------------------|
| HSM 260 Peak | 15-10-1995-33-H | 10/23/15 17:25 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 20:23 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.5 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.5 | 2.8 | 1.00 | |
| Aniline | ND | 9.5 | 1.4 | 1.00 | |
| Anthracene | ND | 9.5 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.5 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.2 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.5 | 4.4 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.5 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.5 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.5 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.5 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.5 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.5 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.3 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.5 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.5 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.5 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.5 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.5 | 2.6 | 1.00 | |
| 2-Chlorophenol | ND | 9.5 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.5 | 2.5 | 1.00 | |
| Chrysene | ND | 9.5 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.5 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.5 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.5 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.5 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.5 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.5 | 2.6 | 1.00 | |
| Dimethyl Phthalate | ND | 9.5 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.5 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.5 | 2.2 | 1.00 | |
| Fluoranthene | ND | 9.5 | 3.0 | 1.00 | |
| Fluorene | ND | 9.5 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.5 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.5 | 2.9 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.6 | 1.00 | |
| Hexachloroethane | ND | 9.5 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.5 | 2.0 | 1.00 | |
| Isophorone | ND | 9.5 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.5 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.5 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.5 | 2.2 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.5 | 3.0 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.5 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.5 | 2.7 | 1.00 | |
| 4-Nitroaniline | ND | 9.5 | 2.0 | 1.00 | |
| 3-Nitroaniline | ND | 9.5 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.5 | 2.1 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.5 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.5 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.5 | 4.4 | 1.00 | |
| Phenanthrene | ND | 9.5 | 2.8 | 1.00 | |
| Phenol | ND | 9.5 | 2.0 | 1.00 | |
| Pyrene | ND | 9.5 | 2.8 | 1.00 | |
| Pyridine | ND | 9.5 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.5 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.5 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 74 | 33-120 | |
| 2-Fluorophenol | 48 | 24-120 | |
| Nitrobenzene-d5 | 69 | 38-120 | |
| p-Terphenyl-d14 | 80 | 41-137 | |
| Phenol-d6 | 31 | 16-120 | |
| 2,4,6-Tribromophenol | 94 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|-----------------|-----------------|---------------------------|------------------|
| HSM 270 Peak | 15-10-1995-34-H | 10/23/15 17:45 | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 20:42 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|------------------------------|---------------|-----------|------------|-----------|-------------------|
| Acenaphthene | ND | 9.6 | 2.7 | 1.00 | |
| Acenaphthylene | ND | 9.6 | 2.8 | 1.00 | |
| Aniline | ND | 9.6 | 1.4 | 1.00 | |
| Anthracene | ND | 9.6 | 2.9 | 1.00 | |
| Azobenzene | ND | 9.6 | 2.5 | 1.00 | |
| Benzidine | ND | 48 | 6.3 | 1.00 | |
| Benzo (a) Anthracene | ND | 9.6 | 4.5 | 1.00 | |
| Benzo (a) Pyrene | ND | 9.6 | 2.3 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 9.6 | 2.2 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 9.6 | 2.4 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 9.6 | 3.1 | 1.00 | |
| Benzoic Acid | ND | 48 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 9.6 | 2.1 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 9.6 | 2.4 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 24 | 2.4 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 9.6 | 3.1 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 9.6 | 3.0 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4-Chloroaniline | ND | 9.6 | 1.9 | 1.00 | |
| 2-Chloronaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Chlorophenol | ND | 9.6 | 2.2 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 9.6 | 2.6 | 1.00 | |
| Chrysene | ND | 9.6 | 2.7 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 9.6 | 2.8 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 9.6 | 2.4 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 9.6 | 2.4 | 1.00 | |
| Dibenzofuran | ND | 9.6 | 2.7 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 9.6 | 2.9 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 9.6 | 2.8 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 24 | 2.5 | 1.00 | |
| 2,4-Dichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|-----|-----|------|------------|
| Diethyl Phthalate | ND | 9.6 | 2.7 | 1.00 | |
| Dimethyl Phthalate | ND | 9.6 | 2.5 | 1.00 | |
| 2,4-Dimethylphenol | ND | 9.6 | 2.3 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 48 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 48 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 9.6 | 2.2 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 9.6 | 2.3 | 1.00 | |
| Fluoranthene | ND | 9.6 | 3.0 | 1.00 | |
| Fluorene | ND | 9.6 | 2.6 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 9.6 | 2.8 | 1.00 | |
| Hexachlorobenzene | ND | 9.6 | 3.0 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 24 | 6.7 | 1.00 | |
| Hexachloroethane | ND | 9.6 | 2.9 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 9.6 | 2.1 | 1.00 | |
| Isophorone | ND | 9.6 | 2.4 | 1.00 | |
| 2-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 1-Methylnaphthalene | ND | 9.6 | 2.7 | 1.00 | |
| 2-Methylphenol | ND | 9.6 | 2.0 | 1.00 | |
| 3/4-Methylphenol | ND | 9.6 | 2.1 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 9.6 | 2.3 | 1.00 | |
| N-Nitrosodimethylamine | ND | 9.6 | 3.1 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 9.6 | 2.6 | 1.00 | |
| Naphthalene | ND | 9.6 | 2.8 | 1.00 | |
| 4-Nitroaniline | ND | 9.6 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| 2-Nitroaniline | ND | 9.6 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 24 | 2.9 | 1.00 | |
| 4-Nitrophenol | ND | 9.6 | 1.5 | 1.00 | |
| 2-Nitrophenol | ND | 9.6 | 2.5 | 1.00 | |
| Pentachlorophenol | ND | 9.6 | 4.5 | 1.00 | |
| Phenanthrene | ND | 9.6 | 2.8 | 1.00 | |
| Phenol | ND | 9.6 | 2.0 | 1.00 | |
| Pyrene | ND | 9.6 | 2.9 | 1.00 | |
| Pyridine | ND | 9.6 | 2.9 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 9.6 | 2.7 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 9.6 | 2.4 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 69 | 33-120 | |
| 2-Fluorophenol | 47 | 24-120 | |
| Nitrobenzene-d5 | 66 | 38-120 | |
| p-Terphenyl-d14 | 77 | 41-137 | |
| Phenol-d6 | 29 | 16-120 | |
| 2,4,6-Tribromophenol | 91 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-4133 | N/A | Aqueous | GC/MS SS | 10/30/15 | 10/31/15 16:20 | 151030L04 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 68 | 33-120 | |
| 2-Fluorophenol | 74 | 24-120 | |
| Nitrobenzene-d5 | 75 | 38-120 | |
| p-Terphenyl-d14 | 69 | 41-137 | |
| Phenol-d6 | 74 | 16-120 | |
| 2,4,6-Tribromophenol | 77 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

Page 100 of 102

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 095-01-003-4132 | N/A | Aqueous | GC/MS TT | 10/30/15 | 10/31/15 21:38 | 151030L05 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|------------------------------|--------|----|-----|------|------------|
| Acenaphthene | ND | 10 | 2.8 | 1.00 | |
| Acenaphthylene | ND | 10 | 2.9 | 1.00 | |
| Aniline | ND | 10 | 1.5 | 1.00 | |
| Anthracene | ND | 10 | 3.0 | 1.00 | |
| Azobenzene | ND | 10 | 2.6 | 1.00 | |
| Benzidine | ND | 50 | 6.5 | 1.00 | |
| Benzo (a) Anthracene | ND | 10 | 4.7 | 1.00 | |
| Benzo (a) Pyrene | ND | 10 | 2.4 | 1.00 | |
| Benzo (b) Fluoranthene | ND | 10 | 2.3 | 1.00 | |
| Benzo (g,h,i) Perylene | ND | 10 | 2.5 | 1.00 | |
| Benzo (k) Fluoranthene | ND | 10 | 3.2 | 1.00 | |
| Benzoic Acid | ND | 50 | 12 | 1.00 | |
| Benzyl Alcohol | ND | 10 | 2.2 | 1.00 | |
| Bis(2-Chloroethoxy) Methane | ND | 10 | 2.5 | 1.00 | |
| Bis(2-Chloroethyl) Ether | ND | 25 | 2.5 | 1.00 | |
| Bis(2-Chloroisopropyl) Ether | ND | 10 | 3.2 | 1.00 | |
| Bis(2-Ethylhexyl) Phthalate | ND | 10 | 3.2 | 1.00 | |
| 4-Bromophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Butyl Benzyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| 4-Chloro-3-Methylphenol | ND | 10 | 2.4 | 1.00 | |
| 4-Chloroaniline | ND | 10 | 2.0 | 1.00 | |
| 2-Chloronaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Chlorophenol | ND | 10 | 2.3 | 1.00 | |
| 4-Chlorophenyl-Phenyl Ether | ND | 10 | 2.7 | 1.00 | |
| Chrysene | ND | 10 | 2.8 | 1.00 | |
| Di-n-Butyl Phthalate | ND | 10 | 2.9 | 1.00 | |
| Di-n-Octyl Phthalate | ND | 10 | 2.5 | 1.00 | |
| Dibenz (a,h) Anthracene | ND | 10 | 2.5 | 1.00 | |
| Dibenzofuran | ND | 10 | 2.8 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 10 | 3.0 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 10 | 3.1 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 10 | 2.9 | 1.00 | |
| 3,3'-Dichlorobenzidine | ND | 25 | 2.6 | 1.00 | |
| 2,4-Dichlorophenol | ND | 10 | 2.5 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|----------------------------|--------|----|-----|------|------------|
| Diethyl Phthalate | ND | 10 | 2.8 | 1.00 | |
| Dimethyl Phthalate | ND | 10 | 2.6 | 1.00 | |
| 2,4-Dimethylphenol | ND | 10 | 2.4 | 1.00 | |
| 4,6-Dinitro-2-Methylphenol | ND | 50 | 14 | 1.00 | |
| 2,4-Dinitrophenol | ND | 50 | 13 | 1.00 | |
| 2,4-Dinitrotoluene | ND | 10 | 2.3 | 1.00 | |
| 2,6-Dinitrotoluene | ND | 10 | 2.4 | 1.00 | |
| Fluoranthene | ND | 10 | 3.1 | 1.00 | |
| Fluorene | ND | 10 | 2.7 | 1.00 | |
| Hexachloro-1,3-Butadiene | ND | 10 | 2.9 | 1.00 | |
| Hexachlorobenzene | ND | 10 | 3.1 | 1.00 | |
| Hexachlorocyclopentadiene | ND | 25 | 6.9 | 1.00 | |
| Hexachloroethane | ND | 10 | 3.0 | 1.00 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 10 | 2.1 | 1.00 | |
| Isophorone | ND | 10 | 2.5 | 1.00 | |
| 2-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 1-Methylnaphthalene | ND | 10 | 2.8 | 1.00 | |
| 2-Methylphenol | ND | 10 | 2.1 | 1.00 | |
| 3/4-Methylphenol | ND | 10 | 2.2 | 1.00 | |
| N-Nitroso-di-n-propylamine | ND | 10 | 2.4 | 1.00 | |
| N-Nitrosodimethylamine | ND | 10 | 3.2 | 1.00 | |
| N-Nitrosodiphenylamine | ND | 10 | 2.8 | 1.00 | |
| Naphthalene | ND | 10 | 2.9 | 1.00 | |
| 4-Nitroaniline | ND | 10 | 2.1 | 1.00 | |
| 3-Nitroaniline | ND | 10 | 2.3 | 1.00 | |
| 2-Nitroaniline | ND | 10 | 2.2 | 1.00 | |
| Nitrobenzene | ND | 25 | 3.0 | 1.00 | |
| 4-Nitrophenol | ND | 10 | 1.6 | 1.00 | |
| 2-Nitrophenol | ND | 10 | 2.6 | 1.00 | |
| Pentachlorophenol | ND | 10 | 4.6 | 1.00 | |
| Phenanthrene | ND | 10 | 2.9 | 1.00 | |
| Phenol | ND | 10 | 2.1 | 1.00 | |
| Pyrene | ND | 10 | 3.0 | 1.00 | |
| Pyridine | ND | 10 | 3.0 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 10 | 2.8 | 1.00 | |
| 2,4,6-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |
| 2,4,5-Trichlorophenol | ND | 10 | 2.5 | 1.00 | |

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C
Units: ug/L

Project: EAA 27122

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| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|----------------------|-----------------|-----------------------|-------------------|
| 2-Fluorobiphenyl | 69 | 33-120 | |
| 2-Fluorophenol | 71 | 24-120 | |
| Nitrobenzene-d5 | 62 | 38-120 | |
| p-Terphenyl-d14 | 64 | 41-137 | |
| Phenol-d6 | 67 | 16-120 | |
| 2,4,6-Tribromophenol | 74 | 27-159 | |


Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Peak | 15-10-1995-1-B | 10/23/15 15:01 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 13:50 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|----------|----------------|------------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| 1,4-Bromofluorobenzene | 89 | 80-120 | | | |
| Dibromofluoromethane | 91 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 86 | 75-135 | | | |
| Toluene-d8 | 100 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Peak | 15-10-1995-2-B | 10/23/15 15:16 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 17:04 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 87 | 80-120 | |
| Dibromofluoromethane | 90 | 78-126 | |
| 1,2-Dichloroethane-d4 | 82 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 210 Trail | 15-10-1995-3-B | 10/23/15 16:14 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 17:32 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 87 | 80-120 | |
| Dibromofluoromethane | 90 | 78-126 | |
| 1,2-Dichloroethane-d4 | 84 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 240 Trail | 15-10-1995-4-B | 10/23/15 16:29 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 18:00 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 85 | 80-120 | |
| Dibromofluoromethane | 91 | 78-126 | |
| 1,2-Dichloroethane-d4 | 83 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 250 Trail | 15-10-1995-5-B | 10/23/15 17:09 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 18:28 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 86 | 80-120 | |
| Dibromofluoromethane | 92 | 78-126 | |
| 1,2-Dichloroethane-d4 | 85 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 260 Trail | 15-10-1995-6-B | 10/23/15 16:50 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 18:56 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|----------|----------------|------------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| 1,4-Bromofluorobenzene | 83 | 80-120 | | | |
| Dibromofluoromethane | 91 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 85 | 75-135 | | | |
| Toluene-d8 | 101 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 13 of 72

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HCS 270 Trail | 15-10-1995-7-B | 10/23/15 17:34 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 19:23 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 14 of 72

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 87 | 80-120 | |
| Dibromofluoromethane | 90 | 78-126 | |
| 1,2-Dichloroethane-d4 | 85 | 75-135 | |
| Toluene-d8 | 103 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 260 Trail | 15-10-1995-8-B | 10/23/15 16:50 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 19:51 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 87 | 80-120 | |
| Dibromofluoromethane | 90 | 78-126 | |
| 1,2-Dichloroethane-d4 | 85 | 75-135 | |
| Toluene-d8 | 104 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHCS 270 Trail | 15-10-1995-9-B | 10/23/15 17:34 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 20:19 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 84 | 80-120 | |
| Dibromofluoromethane | 93 | 78-126 | |
| 1,2-Dichloroethane-d4 | 88 | 75-135 | |
| Toluene-d8 | 109 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|----------------|-----------------|-----------------------|-------------------|
| TB15 | 15-10-1995-10-A | 10/23/15 00:00 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 16:37 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 90 | 80-120 | |
| Dibromofluoromethane | 92 | 78-126 | |
| 1,2-Dichloroethane-d4 | 85 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Trail | 15-10-1995-11-B | 10/23/15 20:35 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 20:46 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 84 | 80-120 | |
| Dibromofluoromethane | 95 | 78-126 | |
| 1,2-Dichloroethane-d4 | 90 | 75-135 | |
| Toluene-d8 | 102 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Trail | 15-10-1995-12-B | 10/23/15 21:01 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 21:14 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|----------|----------------|------------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| 1,4-Bromofluorobenzene | 85 | 80-120 | | | |
| Dibromofluoromethane | 92 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 84 | 75-135 | | | |
| Toluene-d8 | 104 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 231 Trail | 15-10-1995-13-B | 10/23/15 21:21 | Aqueous | GC/MS O | 10/30/15 | 10/30/15 21:41 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 85 | 80-120 | |
| Dibromofluoromethane | 96 | 78-126 | |
| 1,2-Dichloroethane-d4 | 88 | 75-135 | |
| Toluene-d8 | 101 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 27 of 72

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 240 Trail | 15-10-1995-14-A | 10/23/15 20:51 | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 14:34 | 151030L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 90 | 80-120 | |
| Dibromofluoromethane | 101 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 250 Trail | 15-10-1995-15-A | 10/23/15 21:06 | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 15:02 | 151030L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 91 | 80-120 | |
| Dibromofluoromethane | 100 | 78-126 | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | |
| Toluene-d8 | 96 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Trail | 15-10-1995-16-A | 10/23/15 21:35 | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 15:30 | 151030L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 91 | 80-120 | |
| Dibromofluoromethane | 102 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Trail | 15-10-1995-17-A | 10/23/15 21:55 | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 12:17 | 151030L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 34 of 72

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 92 | 80-120 | |
| Dibromofluoromethane | 100 | 78-126 | |
| 1,2-Dichloroethane-d4 | 101 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 210 Trail | 15-10-1995-18-A | 10/23/15 20:35 | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 15:57 | 151030L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 89 | 80-120 | |
| Dibromofluoromethane | 102 | 78-126 | |
| 1,2-Dichloroethane-d4 | 105 | 75-135 | |
| Toluene-d8 | 96 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 230 Trail | 15-10-1995-19-A | 10/23/15 21:01 | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 16:25 | 151030L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|----------|----------------|------------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| 1,4-Bromofluorobenzene | 90 | 80-120 | | | |
| Dibromofluoromethane | 101 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 104 | 75-135 | | | |
| Toluene-d8 | 96 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| FDHSM 231 Trail | 15-10-1995-20-A | 10/23/15 21:21 | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 16:53 | 151030L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 89 | 80-120 | |
| Dibromofluoromethane | 102 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 96 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 210 Lead | 15-10-1995-21-A | 10/23/15 15:25 | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 17:20 | 151030L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 92 | 80-120 | |
| Dibromofluoromethane | 102 | 78-126 | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 230 Lead | 15-10-1995-22-A | 10/23/15 15:41 | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 17:12 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | 1.1 | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 105 | 78-126 | |
| 1,2-Dichloroethane-d4 | 101 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 231 Lead | 15-10-1995-23-A | 10/23/15 16:02 | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 13:39 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|----------|----------------|------------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| 1,4-Bromofluorobenzene | 98 | 80-120 | | | |
| Dibromofluoromethane | 104 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | | | |
| Toluene-d8 | 99 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 47 of 72

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 240 Lead | 15-10-1995-24-A | 10/23/15 15:30 | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 17:38 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | 0.38 | 1.0 | 0.37 | 1.00 | J |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 104 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 250 Lead | 15-10-1995-25-A | 10/23/15 16:21 | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 18:05 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|-----------------|-----------------------|-------------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | |
| 1,4-Bromofluorobenzene | 98 | 80-120 | | | |
| Dibromofluoromethane | 104 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | | | |
| Toluene-d8 | 99 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 260 Lead | 15-10-1995-26-A | 10/23/15 15:58 | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 18:31 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|----------|----------------|------------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| 1,4-Bromofluorobenzene | 98 | 80-120 | | | |
| Dibromofluoromethane | 105 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | | | |
| Toluene-d8 | 100 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 270 Lead | 15-10-1995-27-A | 10/23/15 16:15 | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 18:58 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|----------|----------------|------------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| 1,4-Bromofluorobenzene | 98 | 80-120 | | | |
| Dibromofluoromethane | 105 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | | | |
| Toluene-d8 | 100 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 55 of 72

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 210 Peak | 15-10-1995-28-A | 10/23/15 17:02 | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 19:25 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 56 of 72

| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|----------|----------------|------------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| 1,4-Bromofluorobenzene | 98 | 80-120 | | | |
| Dibromofluoromethane | 107 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | | | |
| Toluene-d8 | 100 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 230 Peak | 15-10-1995-29-A | 10/23/15 17:16 | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 19:51 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|-----------------|-----------------------|-------------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | |
| 1,4-Bromofluorobenzene | 99 | 80-120 | | | |
| Dibromofluoromethane | 105 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | | | |
| Toluene-d8 | 100 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|---------------------------|----------------|-----------------|-----------------|---------------------------|-------------------|
| HSM 231 Peak | 15-10-1995-30-A | 10/23/15 17:06 | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 20:18 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 105 | 78-126 | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 240 Peak | 15-10-1995-31-A | 10/23/15 17:34 | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 20:44 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Parameter | Result | RL | MDL | DF | Qualifiers |
|---------------------------------------|--------|------|------|------|------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | 0.26 | 1.0 | 0.24 | 1.00 | J |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------------|----------|----------------|------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 105 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 63 of 72

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| HSM 260 Peak | 15-10-1995-33-A | 10/23/15 17:25 | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 21:11 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 98 | 80-120 | |
| Dibromofluoromethane | 106 | 78-126 | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|------------------------|-----------------------|----------------|-----------------|-----------------|-----------------------|-------------------|
| HSM 270 Peak | 15-10-1995-34-A | 10/23/15 17:45 | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 21:37 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|-----------|-------------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|-----------------|-----------------------|-------------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | |
| 1,4-Bromofluorobenzene | 98 | 80-120 | | | |
| Dibromofluoromethane | 106 | 78-126 | | | |
| 1,2-Dichloroethane-d4 | 103 | 75-135 | | | |
| Toluene-d8 | 99 | 80-120 | | | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-18638 | N/A | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 11:50 | 151030L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 90 | 80-120 | |
| Dibromofluoromethane | 100 | 78-126 | |
| 1,2-Dichloroethane-d4 | 102 | 75-135 | |
| Toluene-d8 | 97 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

Page 69 of 72

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-18641 | N/A | Aqueous | GC/MS O | 10/30/15 | 10/30/15 13:14 | 151030L006 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 89 | 80-120 | |
| Dibromofluoromethane | 90 | 78-126 | |
| 1,2-Dichloroethane-d4 | 87 | 75-135 | |
| Toluene-d8 | 100 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-001-18642 | N/A | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 12:46 | 151030L010 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------------------------|--------|------|------|------|------------|
| Acetone | ND | 20 | 10 | 1.00 | |
| Benzene | ND | 0.50 | 0.14 | 1.00 | |
| Bromobenzene | ND | 1.0 | 0.30 | 1.00 | |
| Bromochloromethane | ND | 1.0 | 0.48 | 1.00 | |
| Bromodichloromethane | ND | 1.0 | 0.21 | 1.00 | |
| Bromoform | ND | 1.0 | 0.50 | 1.00 | |
| Bromomethane | ND | 10 | 3.9 | 1.00 | |
| 2-Butanone | ND | 10 | 2.2 | 1.00 | |
| n-Butylbenzene | ND | 1.0 | 0.23 | 1.00 | |
| sec-Butylbenzene | ND | 1.0 | 0.25 | 1.00 | |
| tert-Butylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Carbon Disulfide | ND | 10 | 0.41 | 1.00 | |
| Carbon Tetrachloride | ND | 0.50 | 0.23 | 1.00 | |
| Chlorobenzene | ND | 1.0 | 0.17 | 1.00 | |
| Chloroethane | ND | 5.0 | 2.3 | 1.00 | |
| Chloroform | ND | 1.0 | 0.46 | 1.00 | |
| Chloromethane | ND | 10 | 1.8 | 1.00 | |
| 2-Chlorotoluene | ND | 1.0 | 0.24 | 1.00 | |
| 4-Chlorotoluene | ND | 1.0 | 0.13 | 1.00 | |
| Dibromochloromethane | ND | 1.0 | 0.25 | 1.00 | |
| 1,2-Dibromo-3-Chloropropane | ND | 5.0 | 1.2 | 1.00 | |
| 1,2-Dibromoethane | ND | 1.0 | 0.36 | 1.00 | |
| Dibromomethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.46 | 1.00 | |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.40 | 1.00 | |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.43 | 1.00 | |
| Dichlorodifluoromethane | ND | 1.0 | 0.46 | 1.00 | |
| 1,1-Dichloroethane | ND | 1.0 | 0.28 | 1.00 | |
| 1,2-Dichloroethane | ND | 0.50 | 0.24 | 1.00 | |
| 1,1-Dichloroethene | ND | 1.0 | 0.43 | 1.00 | |
| c-1,2-Dichloroethene | ND | 1.0 | 0.48 | 1.00 | |
| t-1,2-Dichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| 1,2-Dichloropropane | ND | 1.0 | 0.42 | 1.00 | |
| 1,3-Dichloropropane | ND | 1.0 | 0.30 | 1.00 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B
Units: ug/L

Project: EAA 27122

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| <u>Parameter</u> | <u>Result</u> | <u>RL</u> | <u>MDL</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------------------|---------------|-----------|------------|-----------|-------------------|
| 2,2-Dichloropropane | ND | 1.0 | 0.36 | 1.00 | |
| 1,1-Dichloropropene | ND | 1.0 | 0.46 | 1.00 | |
| c-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| t-1,3-Dichloropropene | ND | 0.50 | 0.25 | 1.00 | |
| Ethylbenzene | ND | 1.0 | 0.14 | 1.00 | |
| 2-Hexanone | ND | 10 | 2.1 | 1.00 | |
| Isopropylbenzene | ND | 1.0 | 0.58 | 1.00 | |
| p-Isopropyltoluene | ND | 1.0 | 0.16 | 1.00 | |
| Methylene Chloride | ND | 10 | 0.64 | 1.00 | |
| 4-Methyl-2-Pentanone | ND | 10 | 4.4 | 1.00 | |
| Naphthalene | ND | 10 | 2.5 | 1.00 | |
| n-Propylbenzene | ND | 1.0 | 0.17 | 1.00 | |
| Styrene | ND | 1.0 | 0.17 | 1.00 | |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.40 | 1.00 | |
| 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.41 | 1.00 | |
| Tetrachloroethene | ND | 1.0 | 0.39 | 1.00 | |
| Toluene | ND | 1.0 | 0.24 | 1.00 | |
| 1,2,3-Trichlorobenzene | ND | 1.0 | 0.51 | 1.00 | |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.50 | 1.00 | |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.30 | 1.00 | |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane | ND | 10 | 0.78 | 1.00 | |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.38 | 1.00 | |
| Trichloroethene | ND | 1.0 | 0.37 | 1.00 | |
| Trichlorofluoromethane | ND | 10 | 1.7 | 1.00 | |
| 1,2,3-Trichloropropane | ND | 5.0 | 0.64 | 1.00 | |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.36 | 1.00 | |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.28 | 1.00 | |
| Vinyl Acetate | ND | 10 | 2.8 | 1.00 | |
| Vinyl Chloride | ND | 0.50 | 0.30 | 1.00 | |
| p/m-Xylene | ND | 1.0 | 0.30 | 1.00 | |
| o-Xylene | ND | 1.0 | 0.23 | 1.00 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 1.0 | 0.31 | 1.00 | |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------------|-----------------|-----------------------|-------------------|
| 1,4-Bromofluorobenzene | 97 | 80-120 | |
| Dibromofluoromethane | 103 | 78-126 | |
| 1,2-Dichloroethane-d4 | 98 | 75-135 | |
| Toluene-d8 | 99 | 80-120 | |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

10/27/15

Work Order:

15-10-1995

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 260 Peak | 15-10-1995-1 | 10/23/15 15:01 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.050 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 203 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 203 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 305 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 8.7 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.48 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 10 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

| | | | |
|---------------------|---------------------|-----------------------|----------------|
| HCS 270 Peak | 15-10-1995-2 | 10/23/15 15:16 | Aqueous |
|---------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.055 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 204 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 204 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 285 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 14 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.56 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.70 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

10/27/15

Work Order:

15-10-1995

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 210 Trail | 15-10-1995-3 | 10/23/15 16:14 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.16 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 102 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 102 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 125 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 9.5 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.50 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.98 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

| | | | |
|----------------------|---------------------|-----------------------|----------------|
| HCS 240 Trail | 15-10-1995-4 | 10/23/15 16:29 | Aqueous |
|----------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.031 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 230 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 230 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 295 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.4 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.54 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.1 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

10/27/15

Work Order:

15-10-1995

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 250 Trail | 15-10-1995-5 | 10/23/15 17:09 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.046 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 208 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 208 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 235 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.7 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.54 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.70 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

| | | | |
|----------------------|---------------------|-----------------------|----------------|
| HCS 260 Trail | 15-10-1995-6 | 10/23/15 16:50 | Aqueous |
|----------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.040 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 219 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 219 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 275 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 5.7 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.66 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.84 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

10/27/15

Work Order:

15-10-1995

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|---------------------|-----------------------|----------------|
| HCS 270 Trail | 15-10-1995-7 | 10/23/15 17:34 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.050 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 206 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 206 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 315 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 2.4 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.68 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.98 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

| | | | |
|------------------------|---------------------|-----------------------|----------------|
| FDHCS 260 Trail | 15-10-1995-8 | 10/23/15 16:50 | Aqueous |
|------------------------|---------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.039 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 220 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 220 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 285 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 2.8 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.68 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

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Project: EAA 27122

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|------------------------|---------------------|-----------------------|----------------|
| FDHCS 270 Trail | 15-10-1995-9 | 10/23/15 17:34 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.038 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 206 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 206 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 295 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 4.4 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.69 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.84 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

| HSM 210 Trail | 15-10-1995-11 | 10/23/15 20:35 | Aqueous |
|----------------------|----------------------|-----------------------|----------------|
|----------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.026 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 228 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 228 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Solids, Total Dissolved (24) | 345 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 2.0 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.43 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.56 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|----------------------|-----------------------|----------------|
| HSM 230 Trail | 15-10-1995-12 | 10/23/15 21:01 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.039 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 228 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 228 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 290 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.36 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.84 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

| HSM 231 Trail | 15-10-1995-13 | 10/23/15 21:21 | Aqueous |
|----------------------|----------------------|-----------------------|----------------|
|----------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.034 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 245 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 245 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Solids, Total Dissolved (24) | 235 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.48 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.84 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|----------------------|-----------------------|----------------|
| HSM 240 Trail | 15-10-1995-14 | 10/23/15 20:51 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.027 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 248 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 248 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 270 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.49 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

| | | | |
|----------------------|----------------------|-----------------------|----------------|
| HSM 250 Trail | 15-10-1995-15 | 10/23/15 21:06 | Aqueous |
|----------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.024 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 162 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 162 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 315 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.8 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.55 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.49 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|----------------------|-----------------------|----------------|
| HSM 260 Trail | 15-10-1995-16 | 10/23/15 21:35 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.033 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 148 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 148 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 305 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 3.7 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.68 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

| | | | |
|----------------------|----------------------|-----------------------|----------------|
| HSM 270 Trail | 15-10-1995-17 | 10/23/15 21:55 | Aqueous |
|----------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.038 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 142 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 142 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 270 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 6.5 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.63 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|------------------------|----------------------|-----------------------|----------------|
| FDHSM 210 Trail | 15-10-1995-18 | 10/23/15 20:35 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.036 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 233 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 233 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Solids, Total Dissolved (24) | 370 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 4.7 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.39 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

| | | | |
|------------------------|----------------------|-----------------------|----------------|
| FDHSM 230 Trail | 15-10-1995-19 | 10/23/15 21:01 | Aqueous |
|------------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.058 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 222 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 222 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 310 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 4.0 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.10 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.84 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|------------------------|----------------------|-----------------------|----------------|
| FDHSM 231 Trail | 15-10-1995-20 | 10/23/15 21:21 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.031 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 245 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 245 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Solids, Total Dissolved (24) | 295 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 1.1 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.28 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 7.4 | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

| HSM 210 Lead | 15-10-1995-21 | 10/23/15 15:25 | Aqueous |
|---------------------|----------------------|-----------------------|----------------|
|---------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.028 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 203 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 203 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 345 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 9.9 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.20 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.5 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.1 | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|----------------------|-----------------------|----------------|
| HSM 230 Lead | 15-10-1995-22 | 10/23/15 15:41 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-----------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.12 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 123 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 123 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 215 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 161 | 1.00 | 0.829 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.14 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 2.0 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 18 | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | 5.0 | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

| | | | |
|---------------------|----------------------|-----------------------|----------------|
| HSM 231 Lead | 15-10-1995-23 | 10/23/15 16:02 | Aqueous |
|---------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.
(313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|-------------------------------------|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.061 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO3) (24) | 203 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO3) (24) | 203 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO3) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 285 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 9.2 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.34 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 1.3 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 2.3 | 2.5 | 1.2 | 5.00 | J | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|----------------------|-----------------------|----------------|
| HSM 240 Lead | 15-10-1995-24 | 10/23/15 15:30 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.092 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 208 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 208 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 280 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 137 | 1.00 | 0.829 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.40 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.77 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

| | | | |
|---------------------|----------------------|-----------------------|----------------|
| HSM 250 Lead | 15-10-1995-25 | 10/23/15 16:21 | Aqueous |
|---------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.073 | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 203 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 203 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 390 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 12 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.38 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.77 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 3.4 | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|----------------------|-----------------------|----------------|
| HSM 260 Lead | 15-10-1995-26 | 10/23/15 15:58 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.037 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 246 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 246 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 315 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 26 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.52 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.35 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

| | | | |
|---------------------|----------------------|-----------------------|----------------|
| HSM 270 Lead | 15-10-1995-27 | 10/23/15 16:15 | Aqueous |
|---------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.045 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 201 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 201 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 300 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 49 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.68 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

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|----------------------|----------------------|-----------------------|----------------|
| HSM 210 Peak | 15-10-1995-28 | 10/23/15 17:02 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.020 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 156 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 156 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 330 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 4.9 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.41 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 3.1 | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

| | | | |
|---------------------|----------------------|-----------------------|----------------|
| HSM 230 Peak | 15-10-1995-29 | 10/23/15 17:16 | Aqueous |
|---------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.035 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 248 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 248 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 330 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.26 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 5.7 | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

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|----------------------|----------------------|-----------------------|----------------|
| HSM 231 Peak | 15-10-1995-30 | 10/23/15 17:06 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.025 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 150 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 150 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 320 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 5.3 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.37 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

| | | | |
|---------------------|----------------------|-----------------------|----------------|
| HSM 240 Peak | 15-10-1995-31 | 10/23/15 17:34 | Aqueous |
|---------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.020 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 144 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 144 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 315 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 4.3 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.43 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.42 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

10/27/15

Work Order:

15-10-1995

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|----------------------|-----------------------|----------------|
| HSM 260 Peak | 15-10-1995-33 | 10/23/15 17:25 | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.038 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 176 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 176 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 280 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 15 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.51 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.70 | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

| | | | |
|---------------------|----------------------|-----------------------|----------------|
| HSM 270 Peak | 15-10-1995-34 | 10/23/15 17:45 | Aqueous |
|---------------------|----------------------|-----------------------|----------------|

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
 (313) - The reporting limit is elevated resulting from matrix interference.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|----------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | 0.035 | 0.050 | 0.020 | 1.00 | J | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | 190 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | 190 | 5.00 | 0.848 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | 310 | 1.00 | 0.870 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | 16 | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| pH (24) | 7.55 | 0.01 | 0.01 | 1.00 | BV,BU | pH units | N/A | 10/27/15 | SM 4500 H+ B |
| Total Kjeldahl Nitrogen (24) | 0.28 | 0.50 | 0.28 | 1.00 | J | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | 3.6 | 2.5 | 1.2 | 5.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24, 313) | ND | 2.5 | 1.2 | 5.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Analytical Report

SWCA Environmental Consultants

6200 UTSA Blvd., Suite 102

San Antonio, TX 78249-1618

Project: EAA 27122

Date Received:

10/27/15

Work Order:

15-10-1995

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| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix |
|----------------------|-------------------|---------------------|----------------|
| Method Blank | | N/A | Aqueous |

Comment(s): (24) - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Results | RL | MDL | DF | Qualifiers | Units | Date Prepared | Date Analyzed | Method |
|--|---------|-------|-------|------|------------|-------|---------------|---------------|-----------------|
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Phosphorus, Total (24) | ND | 0.050 | 0.020 | 1.00 | | mg/L | N/A | 11/10/15 | EPA 365.1 |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Alkalinity, Total (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Bicarbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/06/15 | SM 2320B |
| Carbonate (as CaCO ₃) (24) | ND | 1.0 | 0.85 | 1.00 | | mg/L | N/A | 11/05/15 | SM 2320B |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.87 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Dissolved (24) | ND | 1.0 | 0.87 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 C |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| Solids, Total Suspended (24) | ND | 1.0 | 0.83 | 1.00 | | mg/L | 10/30/15 | 10/30/15 | SM 2540 D |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Total Kjeldahl Nitrogen (24) | ND | 0.50 | 0.28 | 1.00 | | mg/L | 11/05/15 | 11/05/15 | SM 4500 N Org B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/09/15 | 11/10/15 | SM 5310 B |
| Carbon, Total Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/10/15 | 11/11/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/05/15 | 11/10/15 | SM 5310 B |
| Carbon, Dissolved Organic (24) | ND | 0.50 | 0.24 | 1.00 | | mg/L | 11/05/15 | 11/11/15 | SM 5310 B |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-10-2300-2 | Sample | Aqueous | IC 15 | N/A | 10/31/15 14:13 | 151031S01 |
| 15-10-2300-2 | Matrix Spike | Aqueous | IC 15 | N/A | 10/31/15 16:04 | 151031S01 |
| 15-10-2300-2 | Matrix Spike Duplicate | Aqueous | IC 15 | N/A | 10/31/15 16:24 | 151031S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.1851 | 2.500 | 2.506 | 93 | 2.415 | 89 | 80-120 | 4 | 0-20 | |
| Chloride | 84.26 | 50.00 | 142.4 | 116 | 139.8 | 111 | 80-120 | 2 | 0-20 | |
| Bromide | 0.2843 | 5.000 | 5.444 | 103 | 5.216 | 99 | 80-120 | 4 | 0-20 | |
| Nitrate (as N) | ND | 5.000 | 5.080 | 102 | 4.833 | 97 | 80-120 | 5 | 0-20 | |
| Sulfate | 153.4 | 50.00 | 232.8 | 159 | 229.6 | 152 | 80-120 | 1 | 0-20 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
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San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM 210 Lead | Sample | Aqueous | IC 15 | N/A | 11/01/15 03:50 | 151031S02 |
| HSM 210 Lead | Matrix Spike | Aqueous | IC 15 | N/A | 11/01/15 04:08 | 151031S02 |
| HSM 210 Lead | Matrix Spike Duplicate | Aqueous | IC 15 | N/A | 11/01/15 04:26 | 151031S02 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Fluoride | 0.2071 | 2.500 | 2.526 | 93 | 2.512 | 92 | 80-120 | 1 | 0-20 | |
| Chloride | 27.20 | 50.00 | 82.19 | 110 | 82.11 | 110 | 80-120 | 0 | 0-20 | |
| Bromide | 0.1739 | 5.000 | 5.249 | 102 | 5.247 | 101 | 80-120 | 0 | 0-20 | |
| Nitrate (as N) | 0.6244 | 5.000 | 5.715 | 102 | 5.718 | 102 | 80-120 | 0 | 0-20 | |
| Sulfate | 33.40 | 50.00 | 89.37 | 112 | 89.25 | 112 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

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San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM 210 Trail | Sample | Aqueous | ACA 1 | N/A | 11/10/15 16:37 | 151110S01 |
| HSM 210 Trail | Matrix Spike | Aqueous | ACA 1 | N/A | 11/10/15 16:37 | 151110S01 |
| HSM 210 Trail | Matrix Spike Duplicate | Aqueous | ACA 1 | N/A | 11/10/15 16:37 | 151110S01 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | ND | 0.2000 | 0.1771 | 89 | 0.1776 | 89 | 90-110 | 0 | 0-25 | 3 |

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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
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Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM 270 Peak | Sample | Aqueous | ACA 1 | N/A | 11/10/15 16:37 | 151110S02 |
| HSM 270 Peak | Matrix Spike | Aqueous | ACA 1 | N/A | 11/10/15 16:37 | 151110S02 |
| HSM 270 Peak | Matrix Spike Duplicate | Aqueous | ACA 1 | N/A | 11/10/15 16:37 | 151110S02 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Phosphorus, Total | ND | 0.2000 | 0.2390 | 119 | 0.2396 | 120 | 90-110 | 0 | 0-25 | 3 |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

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Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| FDHSM 231 Trail | Sample | Aqueous | TOC 8 | 11/10/15 | 11/11/15 12:55 | F1110TOCS2 | | | | |
| FDHSM 231 Trail | Matrix Spike | Aqueous | TOC 8 | 11/10/15 | 11/11/15 12:55 | F1110TOCS2 | | | | |
| FDHSM 231 Trail | Matrix Spike Duplicate | Aqueous | TOC 8 | 11/10/15 | 11/11/15 12:55 | F1110TOCS2 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | 7.350 | 50.00 | 55.00 | 95 | 59.00 | 103 | 31-145 | 7 | 0-23 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

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Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HCS 260 Peak | Sample | Aqueous | TOC 11 | 11/09/15 | 11/10/15 09:15 | F1109TOCS2 | | | | |
| HCS 260 Peak | Matrix Spike | Aqueous | TOC 11 | 11/09/15 | 11/10/15 09:15 | F1109TOCS2 | | | | |
| HCS 260 Peak | Matrix Spike Duplicate | Aqueous | TOC 11 | 11/09/15 | 11/10/15 09:15 | F1109TOCS2 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Total Organic | ND | 10.00 | 2.130 | 21 | 3.210 | 32 | 31-145 | 40 | 0-23 | 3,4 |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
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San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|---------------------|-----------------|------------|---------------|-------------------|
| HCS 260 Peak | Sample | Aqueous | TOC 11 | 11/05/15 | 11/10/15 23:32 | F1110DOCS1 | | | | |
| HCS 260 Peak | Matrix Spike | Aqueous | TOC 11 | 11/05/15 | 11/10/15 23:32 | F1110DOCS1 | | | | |
| HCS 260 Peak | Matrix Spike Duplicate | Aqueous | TOC 11 | 11/05/15 | 11/10/15 23:32 | F1110DOCS1 | | | | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Carbon, Dissolved Organic | ND | 50.00 | 41.80 | 84 | 43.05 | 86 | 31-145 | 3 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
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Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|---------------------|-----------------|------------|---------------|-------------------|
| FDHSM 231 Trail | Sample | Aqueous | TOC 11 | 11/05/15 | 11/11/15 09:56 | F1110DOCS2 | | | | |
| FDHSM 231 Trail | Matrix Spike | Aqueous | TOC 11 | 11/05/15 | 11/11/15 09:56 | F1110DOCS2 | | | | |
| FDHSM 231 Trail | Matrix Spike Duplicate | Aqueous | TOC 11 | 11/05/15 | 11/11/15 09:56 | F1110DOCS2 | | | | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Carbon, Dissolved Organic | ND | 50.00 | 47.90 | 96 | 46.70 | 93 | 31-145 | 3 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

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Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HCS 260 Peak | Sample | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:11 | 151028SA8 |
| HCS 260 Peak | Matrix Spike | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:13 | 151028SA8 |
| HCS 260 Peak | Matrix Spike Duplicate | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:15 | 151028SA8 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Calcium | 66.37 | 0.5000 | 71.03 | 4X | 69.31 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 13.90 | 0.5000 | 15.64 | 4X | 14.69 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 1.819 | 5.000 | 6.498 | 94 | 6.482 | 93 | 83-131 | 0 | 0-7 | |
| Sodium | 10.82 | 5.000 | 16.16 | 107 | 15.96 | 103 | 73-127 | 1 | 0-9 | |
| Strontium | 0.5503 | 0.5000 | 1.065 | 103 | 1.051 | 100 | 81-123 | 1 | 0-6 | |
| Silicon | 4.922 | 0.5000 | 5.851 | 4X | 5.514 | 4X | 24-180 | 4X | 0-15 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | | Instrument | Date Prepared | Date Analyzed | | MS/MSD Batch Number | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|------------------|-----------------|------------|---------------------|-------------------|
| HSM 230 Lead | Sample | | Aqueous | | ICP 7300 | 10/28/15 | 10/30/15 12:30 | | 151028SA9 | |
| HSM 230 Lead | Matrix Spike | | Aqueous | | ICP 7300 | 10/28/15 | 10/30/15 12:33 | | 151028SA9 | |
| HSM 230 Lead | Matrix Spike Duplicate | | Aqueous | | ICP 7300 | 10/28/15 | 10/30/15 12:35 | | 151028SA9 | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Calcium | 44.03 | 0.5000 | 44.59 | 4X | 42.77 | 4X | 77-113 | 4X | 0-11 | Q |
| Magnesium | 7.440 | 0.5000 | 7.888 | 4X | 7.617 | 4X | 56-140 | 4X | 0-11 | Q |
| Potassium | 4.132 | 5.000 | 9.456 | 106 | 9.409 | 106 | 83-131 | 1 | 0-7 | |
| Sodium | 7.862 | 5.000 | 13.38 | 110 | 13.18 | 106 | 73-127 | 2 | 0-9 | |
| Strontium | 0.2264 | 0.5000 | 0.7637 | 107 | 0.7338 | 101 | 81-123 | 4 | 0-6 | |
| Silicon | 2.714 | 0.5000 | 3.248 | 4X | 3.178 | 4X | 24-180 | 4X | 0-15 | Q |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|---------------------------|------------------------|-------------|----------|------------|---------------|----------------|---------------------|-----|--------|------------|
| HCS 270 Peak | Sample | | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 21:39 | 151028SA4 | | | |
| HCS 270 Peak | Matrix Spike | | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 21:22 | 151028SA4 | | | |
| HCS 270 Peak | Matrix Spike Duplicate | | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 21:25 | 151028SA4 | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.09641 | 96 | 0.09743 | 97 | 85-133 | 1 | 0-11 | |
| Arsenic | ND | 0.1000 | 0.09269 | 93 | 0.09385 | 94 | 73-127 | 1 | 0-11 | |
| Barium | 0.05104 | 0.1000 | 0.1604 | 109 | 0.1623 | 111 | 74-128 | 1 | 0-10 | |
| Beryllium | ND | 0.1000 | 0.1010 | 101 | 0.09967 | 100 | 56-122 | 1 | 0-11 | |
| Cadmium | ND | 0.1000 | 0.09287 | 93 | 0.09426 | 94 | 84-114 | 1 | 0-8 | |
| Chromium | ND | 0.1000 | 0.1090 | 109 | 0.1102 | 110 | 73-133 | 1 | 0-11 | |
| Copper | ND | 0.1000 | 0.1006 | 101 | 0.09996 | 100 | 72-108 | 1 | 0-10 | |
| Lead | ND | 0.1000 | 0.1086 | 109 | 0.1075 | 108 | 79-121 | 1 | 0-10 | |
| Nickel | 0.001523 | 0.1000 | 0.1011 | 100 | 0.09944 | 98 | 68-122 | 2 | 0-10 | |
| Selenium | ND | 0.1000 | 0.08024 | 80 | 0.07796 | 78 | 59-125 | 3 | 0-12 | |
| Silver | ND | 0.05000 | 0.05385 | 108 | 0.05348 | 107 | 68-128 | 1 | 0-14 | |
| Thallium | ND | 0.1000 | 0.1074 | 107 | 0.1064 | 106 | 73-121 | 1 | 0-11 | |
| Zinc | 0.07542 | 0.1000 | 0.09585 | 20 | 0.09292 | 18 | 43-145 | 3 | 0-39 | 3 |
| Aluminum | ND | 0.1000 | 0.1259 | 126 | 0.1218 | 122 | 47-161 | 3 | 0-24 | |
| Iron | 0.05482 | 5.100 | 5.803 | 113 | 5.744 | 112 | 27-201 | 1 | 0-24 | |
| Manganese | 0.003253 | 0.1000 | 0.1062 | 103 | 0.1067 | 103 | 72-126 | 0 | 0-42 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HSM 231 Lead | Sample | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 22:11 | 151028SA5 | | | | |
| HSM 231 Lead | Matrix Spike | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 21:43 | 151028SA5 | | | | |
| HSM 231 Lead | Matrix Spike Duplicate | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 21:57 | 151028SA5 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Antimony | ND | 0.1000 | 0.1037 | 104 | 0.09573 | 96 | 85-133 | 8 | 0-11 | |
| Arsenic | ND | 0.1000 | 0.09556 | 96 | 0.09159 | 92 | 73-127 | 4 | 0-11 | |
| Barium | 0.03620 | 0.1000 | 0.1541 | 118 | 0.1432 | 107 | 74-128 | 7 | 0-10 | |
| Beryllium | ND | 0.1000 | 0.1033 | 103 | 0.09545 | 95 | 56-122 | 8 | 0-11 | |
| Cadmium | ND | 0.1000 | 0.09864 | 99 | 0.09155 | 92 | 84-114 | 7 | 0-8 | |
| Chromium | ND | 0.1000 | 0.1149 | 115 | 0.1050 | 105 | 73-133 | 9 | 0-11 | |
| Copper | ND | 0.1000 | 0.1028 | 103 | 0.09710 | 97 | 72-108 | 6 | 0-10 | |
| Lead | ND | 0.1000 | 0.1124 | 112 | 0.1055 | 105 | 79-121 | 6 | 0-10 | |
| Nickel | 0.001766 | 0.1000 | 0.1034 | 102 | 0.09686 | 95 | 68-122 | 7 | 0-10 | |
| Selenium | ND | 0.1000 | 0.08353 | 84 | 0.07190 | 72 | 59-125 | 15 | 0-12 | 4 |
| Silver | ND | 0.05000 | 0.05602 | 112 | 0.05324 | 106 | 68-128 | 5 | 0-14 | |
| Thallium | ND | 0.1000 | 0.1099 | 110 | 0.1040 | 104 | 73-121 | 6 | 0-11 | |
| Zinc | 0.02899 | 0.1000 | 0.1064 | 77 | 0.09627 | 67 | 43-145 | 10 | 0-39 | |
| Aluminum | ND | 0.1000 | 0.1217 | 122 | 0.1172 | 117 | 47-161 | 4 | 0-24 | |
| Iron | 0.05717 | 5.100 | 6.018 | 117 | 5.666 | 110 | 27-201 | 6 | 0-24 | |
| Manganese | 0.004124 | 0.1000 | 0.1125 | 108 | 0.1046 | 101 | 72-126 | 7 | 0-42 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HCS 210 Trail | Sample | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:18 | 151105SA3 |
| HCS 210 Trail | Matrix Spike | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:21 | 151105SA3 |
| HCS 210 Trail | Matrix Spike Duplicate | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 20:23 | 151105SA3 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | ND | 0.01000 | 0.01017 | 102 | 0.01024 | 102 | 55-133 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
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San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| HSM 240 Lead | Sample | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:21 | 151105SA4 |
| HSM 240 Lead | Matrix Spike | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:28 | 151105SA4 |
| HSM 240 Lead | Matrix Spike Duplicate | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:30 | 151105SA4 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Mercury | ND | 0.01000 | 0.01000 | 100 | 0.01010 | 101 | 55-133 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | |
|-----------------------------|------------------------|-------------|----------|------------|---------------|----------------|---------------------|-----|--------|------------|
| HCS 260 Peak | Sample | | Aqueous | GC/MS O | 10/30/15 | 10/30/15 13:50 | 151030S001 | | | |
| HCS 260 Peak | Matrix Spike | | Aqueous | GC/MS O | 10/30/15 | 10/30/15 14:18 | 151030S001 | | | |
| HCS 260 Peak | Matrix Spike Duplicate | | Aqueous | GC/MS O | 10/30/15 | 10/30/15 14:46 | 151030S001 | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 53.60 | 107 | 56.41 | 113 | 74-122 | 5 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 58.43 | 117 | 61.88 | 124 | 60-144 | 6 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 51.13 | 102 | 54.37 | 109 | 73-120 | 6 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 42.67 | 85 | 45.23 | 90 | 80-122 | 6 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 47.56 | 95 | 51.07 | 102 | 70-120 | 7 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 45.51 | 91 | 46.13 | 92 | 64-142 | 1 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 55.28 | 111 | 60.09 | 120 | 52-136 | 8 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 56.19 | 112 | 60.86 | 122 | 77-125 | 8 | 0-24 | |
| Toluene | ND | 50.00 | 53.54 | 107 | 56.90 | 114 | 72-126 | 6 | 0-23 | |
| Trichloroethene | ND | 50.00 | 54.14 | 108 | 57.22 | 114 | 74-128 | 6 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 44.95 | 90 | 50.05 | 100 | 67-133 | 11 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 118.9 | 119 | 128.5 | 128 | 63-129 | 8 | 0-25 | |
| o-Xylene | ND | 50.00 | 57.31 | 115 | 61.88 | 124 | 62-128 | 8 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 42.03 | 84 | 43.92 | 88 | 68-134 | 4 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HSM 270 Trail | Sample | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 12:17 | 151030S008 | | | | |
| HSM 270 Trail | Matrix Spike | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 10:01 | 151030S008 | | | | |
| HSM 270 Trail | Matrix Spike Duplicate | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 10:28 | 151030S008 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 51.24 | 102 | 48.56 | 97 | 74-122 | 5 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 53.78 | 108 | 50.71 | 101 | 60-144 | 6 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 54.76 | 110 | 51.42 | 103 | 73-120 | 6 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 50.66 | 101 | 47.77 | 96 | 80-122 | 6 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 55.07 | 110 | 52.49 | 105 | 70-120 | 5 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 48.38 | 97 | 45.58 | 91 | 64-142 | 6 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 55.20 | 110 | 53.18 | 106 | 52-136 | 4 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 57.34 | 115 | 53.66 | 107 | 77-125 | 7 | 0-24 | |
| Toluene | ND | 50.00 | 51.89 | 104 | 49.10 | 98 | 72-126 | 6 | 0-23 | |
| Trichloroethene | ND | 50.00 | 52.86 | 106 | 50.48 | 101 | 74-128 | 5 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 51.14 | 102 | 47.42 | 95 | 67-133 | 8 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 116.4 | 116 | 108.8 | 109 | 63-129 | 7 | 0-25 | |
| o-Xylene | ND | 50.00 | 56.12 | 112 | 52.77 | 106 | 62-128 | 6 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 43.07 | 86 | 41.12 | 82 | 68-134 | 5 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|-----------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| HSM 231 Lead | Sample | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 13:39 | 151030S009 | | | | |
| HSM 231 Lead | Matrix Spike | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 14:06 | 151030S009 | | | | |
| HSM 231 Lead | Matrix Spike Duplicate | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 14:32 | 151030S009 | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Benzene | ND | 50.00 | 46.39 | 93 | 46.36 | 93 | 74-122 | 0 | 0-21 | |
| Carbon Tetrachloride | ND | 50.00 | 55.77 | 112 | 56.04 | 112 | 60-144 | 0 | 0-21 | |
| Chlorobenzene | ND | 50.00 | 46.55 | 93 | 46.25 | 93 | 73-120 | 1 | 0-22 | |
| 1,2-Dibromoethane | ND | 50.00 | 52.05 | 104 | 51.51 | 103 | 80-122 | 1 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 49.74 | 99 | 49.53 | 99 | 70-120 | 0 | 0-26 | |
| 1,2-Dichloroethane | ND | 50.00 | 47.11 | 94 | 47.53 | 95 | 64-142 | 1 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 32.74 | 65 | 32.76 | 66 | 52-136 | 0 | 0-21 | |
| Ethylbenzene | ND | 50.00 | 45.70 | 91 | 44.80 | 90 | 77-125 | 2 | 0-24 | |
| Toluene | ND | 50.00 | 45.99 | 92 | 45.93 | 92 | 72-126 | 0 | 0-23 | |
| Trichloroethene | ND | 50.00 | 45.34 | 91 | 45.06 | 90 | 74-128 | 1 | 0-22 | |
| Vinyl Chloride | ND | 50.00 | 29.82 | 60 | 29.42 | 59 | 67-133 | 1 | 0-20 | 3 |
| p/m-Xylene | ND | 100.0 | 86.13 | 86 | 85.41 | 85 | 63-129 | 1 | 0-25 | |
| o-Xylene | ND | 50.00 | 43.86 | 88 | 43.46 | 87 | 62-128 | 1 | 0-24 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 49.43 | 99 | 50.03 | 100 | 68-134 | 1 | 0-21 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|---------------------|--------------------|------------------|------------------|-----------------|-----------------------|
| HCS 270 Peak | Sample | Aqueous | ICP/MS 03 | 10/28/15 00:00 | 10/30/15 21:39 | 151028SA4 |
| HCS 270 Peak | PDS | Aqueous | ICP/MS 03 | 10/28/15 00:00 | 10/30/15 21:29 | 151028SA4 |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>PDS Conc.</u> | <u>PDS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Antimony | ND | 0.1000 | 0.09399 | 94 | 75-125 | |
| Arsenic | ND | 0.1000 | 0.08828 | 88 | 75-125 | |
| Barium | 0.05104 | 0.1000 | 0.1462 | 95 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.09487 | 95 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.08708 | 87 | 75-125 | |
| Chromium | ND | 0.1000 | 0.1012 | 101 | 75-125 | |
| Copper | ND | 0.1000 | 0.09431 | 94 | 75-125 | |
| Lead | ND | 0.1000 | 0.1004 | 100 | 75-125 | |
| Nickel | 0.001523 | 0.1000 | 0.09544 | 94 | 75-125 | |
| Selenium | ND | 0.1000 | 0.07303 | 73 | 75-125 | 5 |
| Silver | ND | 0.05000 | 0.04148 | 83 | 75-125 | |
| Thallium | ND | 0.1000 | 0.09872 | 99 | 75-125 | |
| Zinc | 0.07542 | 0.1000 | 0.1660 | 91 | 75-125 | |
| Aluminum | ND | 0.1000 | 0.1086 | 109 | 75-125 | |
| Iron | 0.05482 | 5.100 | 5.246 | 102 | 75-125 | |
| Manganese | 0.003253 | 0.1000 | 0.1021 | 99 | 75-125 | |

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|---------------------|--------------------|------------------|-----------------------|-----------------------|-----------------------|
| HSM 231 Lead | Sample | Aqueous | ICP/MS 03 | 10/28/15 00:00 | 10/30/15 22:11 | 151028SA5 |
| HSM 231 Lead | PDS | Aqueous | ICP/MS 03 | 10/28/15 00:00 | 10/30/15 22:00 | 151028SA5 |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>PDS Conc.</u> | <u>PDS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Antimony | ND | 0.1000 | 0.09520 | 95 | 75-125 | |
| Arsenic | ND | 0.1000 | 0.08681 | 87 | 75-125 | |
| Barium | 0.03620 | 0.1000 | 0.1335 | 97 | 75-125 | |
| Beryllium | ND | 0.1000 | 0.09342 | 93 | 75-125 | |
| Cadmium | ND | 0.1000 | 0.08872 | 89 | 75-125 | |
| Chromium | ND | 0.1000 | 0.1003 | 100 | 75-125 | |
| Copper | ND | 0.1000 | 0.09336 | 93 | 75-125 | |
| Lead | ND | 0.1000 | 0.1015 | 101 | 75-125 | |
| Nickel | 0.001766 | 0.1000 | 0.09390 | 92 | 75-125 | |
| Selenium | ND | 0.1000 | 0.07272 | 73 | 75-125 | 5 |
| Silver | ND | 0.05000 | 0.04229 | 85 | 75-125 | |
| Thallium | ND | 0.1000 | 0.09942 | 99 | 75-125 | |
| Zinc | 0.02899 | 0.1000 | 0.1097 | 81 | 75-125 | |
| Aluminum | ND | 0.1000 | 0.1051 | 105 | 75-125 | |
| Iron | 0.05717 | 5.100 | 5.245 | 102 | 75-125 | |
| Manganese | 0.004124 | 0.1000 | 0.1029 | 99 | 75-125 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---|------------------|---------------------|------------------|---------------|----------------|------------------------|
| HCS 260 Peak | Sample | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:45 | F1105ALKD2 |
| HCS 260 Peak | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:45 | F1105ALKD2 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Alkalinity, Total (as CaCO ₃) | | 203.0 | 202.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-11-0235-1 | Sample | Aqueous | PH1/BUR03 | N/A | 11/05/15 16:55 | F1105ALKD3 |
| 15-11-0235-1 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 11/05/15 16:55 | F1105ALKD3 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|---|---------------------|------------------|------------|---------------|-------------------|
| Alkalinity, Total (as CaCO ₃) | 203.0 | 204.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---|-------------------------|---------------------|------------------|---------------|-----------------------|------------------------|
| FDHSM 231 Trail | Sample | Aqueous | PH1/BUR03 | N/A | 11/06/15 19:35 | F1106ALKD2 |
| FDHSM 231 Trail | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 11/06/15 19:35 | F1106ALKD2 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Alkalinity, Total (as CaCO ₃) | | 245.0 | 245.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|-------------------------------------|------------------|---------------------|------------------|---------------|----------------|------------------------|
| HCS 260 Peak | Sample | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:45 | F1105HCOD2 |
| HCS 260 Peak | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:45 | F1105HCOD2 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Bicarbonate (as CaCO ₃) | | 203.0 | 202.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|-------------------------------------|------------------|---------------------|------------------|---------------|----------------|------------------------|
| 15-11-0235-1 | Sample | Aqueous | PH1/BUR03 | N/A | 11/05/15 16:55 | F1105HCOD3 |
| 15-11-0235-1 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 11/05/15 16:55 | F1105HCOD3 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Bicarbonate (as CaCO ₃) | | 203.0 | 204.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|-------------------------------------|------------------|---------------------|------------------|---------------|----------------|------------------------|
| FDHSM 231 Trail | Sample | Aqueous | PH1/BUR03 | N/A | 11/06/15 19:35 | F1106HCOD2 |
| FDHSM 231 Trail | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 11/06/15 19:35 | F1106HCOD2 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Bicarbonate (as CaCO ₃) | | 245.0 | 245.0 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| HCS 260 Peak | Sample | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:45 | F1105CO3D2 |
| HCS 260 Peak | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:45 | F1105CO3D2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-11-0235-1 | Sample | Aqueous | PH1/BUR03 | N/A | 11/05/15 16:55 | F1105CO3D3 |
| 15-11-0235-1 | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 11/05/15 16:55 | F1105CO3D3 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------------|---------------------|------------------|------------|---------------|-------------------|
| Carbonate (as CaCO ₃) | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|-----------------------------------|-------------------------|---------------------|------------------|---------------|-----------------------|------------------------|
| FDHSM 231 Trail | Sample | Aqueous | PH1/BUR03 | N/A | 11/06/15 19:35 | F1106CO3D2 |
| FDHSM 231 Trail | Sample Duplicate | Aqueous | PH1/BUR03 | N/A | 11/06/15 19:35 | F1106CO3D2 |
| <u>Parameter</u> | | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Carbonate (as CaCO ₃) | | ND | ND | N/A | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|----------------|------------|-----------------------|-----------------------|------------------------|
| HSM 210 Trail | Sample | Aqueous | N/A | 10/30/15 00:00 | 10/30/15 16:00 | F1030TDSD3 |
| HSM 210 Trail | Sample Duplicate | Aqueous | N/A | 10/30/15 00:00 | 10/30/15 16:00 | F1030TDSD3 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Dissolved | 345.0 | 360.0 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|----------------|------------|-----------------------|-----------------------|------------------------|
| HSM 230 Lead | Sample | Aqueous | N/A | 10/30/15 00:00 | 10/30/15 17:00 | F1030TDSD5 |
| HSM 230 Lead | Sample Duplicate | Aqueous | N/A | 10/30/15 00:00 | 10/30/15 17:00 | F1030TDSD5 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Dissolved | 215.0 | 210.0 | 2 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| HCS 260 Peak | Sample | Aqueous | N/A | 10/30/15 00:00 | 10/30/15 15:00 | F1030TSSD5 |
| HCS 260 Peak | Sample Duplicate | Aqueous | N/A | 10/30/15 00:00 | 10/30/15 15:00 | F1030TSSD5 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Suspended | 8.700 | 9.100 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|----------------|----------------|------------------------|
| 15-10-2139-2 | Sample | Aqueous | N/A | 10/30/15 00:00 | 10/30/15 17:00 | F1030TSSD6 |
| 15-10-2139-2 | Sample Duplicate | Aqueous | N/A | 10/30/15 00:00 | 10/30/15 17:00 | F1030TSSD6 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Solids, Total Suspended | 740.0 | 720.0 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| 15-10-2000-5 | Sample | Aqueous | PH 1 | N/A | 10/27/15 20:04 | F1027PHD2 |
| 15-10-2000-5 | Sample Duplicate | Aqueous | PH 1 | N/A | 10/27/15 20:04 | F1027PHD2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|------------------|---------------------|------------------|------------|---------------|-------------------|
| pH | 8.080 | 8.050 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 4500 H+ B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|------------------|---------|------------|---------------|----------------|------------------------|
| FDHSM 230 Trail | Sample | Aqueous | PH 1 | N/A | 10/27/15 20:25 | F1027PHD3 |
| FDHSM 230 Trail | Sample Duplicate | Aqueous | PH 1 | N/A | 10/27/15 20:25 | F1027PHD3 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|------------------|---------------------|------------------|------------|---------------|-------------------|
| pH | 7.100 | 7.110 | 0 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 4500 N Org B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|----------------|--------------|-----------------------|-----------------------|------------------------|
| HSM 230 Trail | Sample | Aqueous | BUR05 | 11/05/15 00:00 | 11/05/15 16:05 | F1105TKND1 |
| HSM 230 Trail | Sample Duplicate | Aqueous | BUR05 | 11/05/15 00:00 | 11/05/15 16:05 | F1105TKND1 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Total Kjeldahl Nitrogen | 0.8400 | 0.7700 | 9 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 4500 N Org B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | Duplicate Batch Number |
|---------------------------|-------------------------|----------------|--------------|-----------------------|-----------------------|------------------------|
| HSM 240 Lead | Sample | Aqueous | BUR05 | 11/05/15 00:00 | 11/05/15 17:45 | F1105TKND2 |
| HSM 240 Lead | Sample Duplicate | Aqueous | BUR05 | 11/05/15 00:00 | 11/05/15 17:45 | F1105TKND2 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>DUP Conc.</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-------------------------|---------------------|------------------|------------|---------------|-------------------|
| Total Kjeldahl Nitrogen | 0.7700 | 0.7000 | 10 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 099-12-906-6212 | LCS | Aqueous | IC 15 | N/A | 10/31/15 12:22 | 151031L01 |
| 099-12-906-6212 | LCSD | Aqueous | IC 15 | N/A | 10/31/15 12:40 | 151031L01 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|----------------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| Fluoride | 2.500 | 2.552 | 102 | 2.557 | 102 | 90-110 | 0 | 0-15 | |
| Chloride | 50.00 | 50.97 | 102 | 50.92 | 102 | 90-110 | 0 | 0-15 | |
| Bromide | 5.000 | 5.118 | 102 | 5.125 | 103 | 90-110 | 0 | 0-15 | |
| Nitrate (as N) | 5.000 | 5.075 | 102 | 5.071 | 101 | 90-110 | 0 | 0-15 | |
| Sulfate | 50.00 | 50.51 | 101 | 50.45 | 101 | 90-110 | 0 | 0-15 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 300.0

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|--------------------|------------------------|------------------|-----------------|-------------------|
| 099-12-906-6215 | LCS | Aqueous | IC 15 | N/A | 10/31/15 22:02 | 151031L02 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Fluoride | | 2.500 | 2.341 | 94 | 90-110 | |
| Chloride | | 50.00 | 47.53 | 95 | 90-110 | |
| Bromide | | 5.000 | 4.818 | 96 | 90-110 | |
| Nitrate (as N) | | 5.000 | 4.744 | 95 | 90-110 | |
| Sulfate | | 50.00 | 47.20 | 94 | 90-110 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-739-134 | LCS | Aqueous | ACA 1 | N/A | 11/10/15 16:37 | 151110L01 | | | |
| 099-12-739-134 | LCSD | Aqueous | ACA 1 | N/A | 11/10/15 16:37 | 151110L01 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Phosphorus, Total | 0.2000 | 0.1937 | 97 | 0.1984 | 99 | 90-110 | 2 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: EPA 365.1

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-739-135 | LCS | Aqueous | ACA 1 | N/A | 11/10/15 16:37 | 151110L02 | | | |
| 099-12-739-135 | LCSD | Aqueous | ACA 1 | N/A | 11/10/15 16:37 | 151110L02 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Phosphorus, Total | 0.2000 | 0.1929 | 96 | 0.1949 | 97 | 90-110 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 099-15-859-851 | LCS | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:45 | F1105ALKB2 |
| 099-15-859-851 | LCSD | Aqueous | PH1/BUR03 | N/A | 11/05/15 14:45 | F1105ALKB2 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|---|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| Alkalinity, Total (as CaCO ₃) | 100.0 | 100.0 | 100 | 100.0 | 100 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-859-853 | LCS | Aqueous | PH1/BUR03 | N/A | 11/05/15 16:55 | F1105ALKB3 | | | |
| 099-15-859-853 | LCSD | Aqueous | PH1/BUR03 | N/A | 11/05/15 16:55 | F1105ALKB3 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 100.0 | 100.0 | 100 | 100.0 | 100 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
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San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2320B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|------------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-859-852 | LCS | Aqueous | PH1/BUR03 | N/A | 11/06/15 19:35 | F1106ALKB2 | | | |
| 099-15-859-852 | LCSD | Aqueous | PH1/BUR03 | N/A | 11/06/15 19:35 | F1106ALKB2 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Alkalinity, Total (as CaCO3) | 100.0 | 99.00 | 99 | 99.00 | 99 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4822 | LCS | Aqueous | N/A | 10/30/15 | 10/30/15 16:00 | F1030TDSL3 | | | |
| 099-12-180-4822 | LCSD | Aqueous | N/A | 10/30/15 | 10/30/15 16:00 | F1030TDSL3 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 80.00 | 80 | 85.00 | 85 | 80-120 | 6 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2540 C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-180-4823 | LCS | Aqueous | N/A | 10/30/15 | 10/30/15 17:00 | F1030TDSL5 | | | |
| 099-12-180-4823 | LCSD | Aqueous | N/A | 10/30/15 | 10/30/15 17:00 | F1030TDSL5 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Dissolved | 100.0 | 110.0 | 110 | 105.0 | 105 | 80-120 | 5 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-09-010-7388 | LCS | Aqueous | N/A | 10/30/15 | 10/30/15 15:00 | F1030TSSL5 | | | |
| 099-09-010-7388 | LCSD | Aqueous | N/A | 10/30/15 | 10/30/15 15:00 | F1030TSSL5 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Solids, Total Suspended | 100.0 | 105.0 | 105 | 109.0 | 109 | 80-120 | 4 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 2540 D

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------------|------------|---------------|-------------------|
| 099-09-010-7387 | LCS | Aqueous | N/A | 10/30/15 | 10/30/15 17:00 | F1030TSSL6 | | | |
| 099-09-010-7387 | LCSD | Aqueous | N/A | 10/30/15 | 10/30/15 17:00 | F1030TSSL6 | | | |
| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Solids, Total Suspended | 100.0 | 106.0 | 106 | 105.0 | 105 | 80-120 | 1 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 099-05-097-5779 | LCS | Aqueous | TOC 8 | 11/10/15 | 11/11/15 12:55 | F1110TOCL2 |
| 099-05-097-5779 | LCSD | Aqueous | TOC 8 | 11/10/15 | 11/11/15 12:55 | F1110TOCL2 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| Carbon, Total Organic | 10.00 | 10.80 | 108 | 11.10 | 111 | 80-120 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 099-05-097-5778 | LCS | Aqueous | TOC 11 | 11/09/15 | 11/10/15 09:15 | F1109TOCL2 |
| 099-05-097-5778 | LCSD | Aqueous | TOC 11 | 11/09/15 | 11/10/15 09:15 | F1109TOCL2 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| Carbon, Total Organic | 10.00 | 9.490 | 95 | 9.230 | 92 | 80-120 | 3 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-115-1467 | LCS | Aqueous | TOC 11 | 11/05/15 | 11/10/15 23:32 | F1110DOCL1 | | | |
| 099-05-115-1467 | LCSD | Aqueous | TOC 11 | 11/05/15 | 11/10/15 23:32 | F1110DOCL1 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Dissolved Organic | 10.00 | 9.720 | 97 | 9.740 | 97 | 80-120 | 0 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: N/A
Method: SM 5310 B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-05-115-1468 | LCS | Aqueous | TOC 11 | 11/05/15 | 11/11/15 09:56 | F1110DOCL2 | | | |
| 099-05-115-1468 | LCSD | Aqueous | TOC 11 | 11/05/15 | 11/11/15 09:56 | F1110DOCL2 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Carbon, Dissolved Organic | 10.00 | 10.40 | 104 | 10.20 | 102 | 80-120 | 2 | 0-20 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-15-683-1458 | LCS | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 19:04 | 151028LA8F |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Calcium | | 0.5000 | 0.5160 | 103 | 80-120 | |
| Magnesium | | 0.5000 | 0.5054 | 101 | 80-120 | |
| Potassium | | 5.000 | 4.722 | 94 | 80-120 | |
| Sodium | | 5.000 | 4.855 | 97 | 80-120 | |
| Strontium | | 0.5000 | 0.4955 | 99 | 80-120 | |
| Silicon | | 0.5000 | 0.4764 | 95 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

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6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6010B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------|-------------|-----------------|---------------|----------------|------------------|
| 099-15-683-1442 | LCS | Aqueous | ICP 7300 | 10/28/15 | 10/30/15 12:22 | 151028LA9F |
| Parameter | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | Qualifiers |
| Calcium | | 0.5000 | 0.4977 | 100 | 80-120 | |
| Magnesium | | 0.5000 | 0.5121 | 102 | 80-120 | |
| Potassium | | 5.000 | 5.251 | 105 | 80-120 | |
| Sodium | | 5.000 | 5.479 | 110 | 80-120 | |
| Strontium | | 0.5000 | 0.5265 | 105 | 80-120 | |
| Silicon | | 0.5000 | 0.5240 | 105 | 80-120 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

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San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-963 | LCS | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 21:15 | 151028LA4F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.09884 | 99 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.09837 | 98 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.09643 | 96 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.1005 | 101 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.09867 | 99 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.09479 | 95 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.1012 | 101 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.09665 | 97 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.09814 | 98 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.08975 | 90 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.04447 | 89 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.09317 | 93 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.09933 | 99 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.09923 | 99 | 80-120 | 73-127 | |
| Iron | 5.100 | 5.206 | 102 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.1009 | 101 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-15-693-964 | LCS | Aqueous | ICP/MS 03 | 10/28/15 | 10/30/15 21:18 | 151028LA5F |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Antimony | 0.1000 | 0.09827 | 98 | 80-120 | 73-127 | |
| Arsenic | 0.1000 | 0.09624 | 96 | 80-120 | 73-127 | |
| Barium | 0.1000 | 0.09508 | 95 | 80-120 | 73-127 | |
| Beryllium | 0.1000 | 0.1005 | 101 | 80-120 | 73-127 | |
| Cadmium | 0.1000 | 0.09656 | 97 | 80-120 | 73-127 | |
| Chromium | 0.1000 | 0.09415 | 94 | 80-120 | 73-127 | |
| Copper | 0.1000 | 0.09892 | 99 | 80-120 | 73-127 | |
| Lead | 0.1000 | 0.09651 | 97 | 80-120 | 73-127 | |
| Nickel | 0.1000 | 0.09432 | 94 | 80-120 | 73-127 | |
| Selenium | 0.1000 | 0.08904 | 89 | 80-120 | 73-127 | |
| Silver | 0.05000 | 0.04447 | 89 | 80-120 | 73-127 | |
| Thallium | 0.1000 | 0.09366 | 94 | 80-120 | 73-127 | |
| Zinc | 0.1000 | 0.09761 | 98 | 80-120 | 73-127 | |
| Aluminum | 0.1000 | 0.09823 | 98 | 80-120 | 73-127 | |
| Iron | 5.100 | 5.085 | 100 | 80-120 | 73-127 | |
| Manganese | 0.1000 | 0.09653 | 97 | 80-120 | 73-127 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| 099-15-763-654 | LCS | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:19 | 151105LA3F |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.01000 | 0.009980 | 100 | 80-120 | |

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Quality Control - LCS

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6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 7470A Filt.
Method: EPA 7470A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|----------------|-------------------|-----------------|-----------------------|-------------------|
| 099-15-763-655 | LCS | Aqueous | Mercury 04 | 11/05/15 | 11/05/15 21:19 | 151105LA4F |

| <u>Parameter</u> | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|--------------------|------------------------|------------------|-----------------|-------------------|
| Mercury | 0.01000 | 0.009980 | 100 | 80-120 | |

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Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-529-851 | LCS | Aqueous | | GC 44 | 10/29/15 | 10/30/15 13:29 | 151029L12 | | | |
| 099-12-529-851 | LCSD | Aqueous | | GC 44 | 10/29/15 | 10/30/15 13:43 | 151029L12 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Alpha-BHC | 0.5000 | 0.5212 | 104 | 0.5246 | 105 | 50-135 | 36-149 | 1 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.5451 | 109 | 0.5510 | 110 | 50-135 | 36-149 | 1 | 0-25 | |
| Beta-BHC | 0.5000 | 0.5091 | 102 | 0.5149 | 103 | 50-135 | 36-149 | 1 | 0-25 | |
| Heptachlor | 0.5000 | 0.4684 | 94 | 0.4774 | 95 | 50-135 | 36-149 | 2 | 0-25 | |
| Delta-BHC | 0.5000 | 0.5177 | 104 | 0.5188 | 104 | 50-135 | 36-149 | 0 | 0-25 | |
| Aldrin | 0.5000 | 0.4176 | 84 | 0.4293 | 86 | 50-135 | 36-149 | 3 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.4869 | 97 | 0.4884 | 98 | 50-135 | 36-149 | 0 | 0-25 | |
| Endosulfan I | 0.5000 | 0.4903 | 98 | 0.4896 | 98 | 50-135 | 36-149 | 0 | 0-25 | |
| Dieldrin | 0.5000 | 0.5074 | 101 | 0.5082 | 102 | 50-135 | 36-149 | 0 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.4968 | 99 | 0.5034 | 101 | 50-135 | 36-149 | 1 | 0-25 | |
| Endrin | 0.5000 | 0.6128 | 123 | 0.6133 | 123 | 50-135 | 36-149 | 0 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.4507 | 90 | 0.4471 | 89 | 50-135 | 36-149 | 1 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.5030 | 101 | 0.5058 | 101 | 50-135 | 36-149 | 1 | 0-25 | |
| Endosulfan II | 0.5000 | 0.4986 | 100 | 0.5004 | 100 | 50-135 | 36-149 | 0 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.5245 | 105 | 0.5191 | 104 | 50-135 | 36-149 | 1 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.5056 | 101 | 0.5078 | 102 | 50-135 | 36-149 | 0 | 0-25 | |
| Methoxychlor | 0.5000 | 0.5595 | 112 | 0.5592 | 112 | 50-135 | 36-149 | 0 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8081A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-529-852 | LCS | Aqueous | | GC 44 | 10/29/15 | 10/30/15 14:12 | 151029L13 | | | |
| 099-12-529-852 | LCSD | Aqueous | | GC 44 | 10/29/15 | 10/30/15 14:26 | 151029L13 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Alpha-BHC | 0.5000 | 0.5198 | 104 | 0.5532 | 111 | 50-135 | 36-149 | 6 | 0-25 | |
| Gamma-BHC | 0.5000 | 0.5467 | 109 | 0.5851 | 117 | 50-135 | 36-149 | 7 | 0-25 | |
| Beta-BHC | 0.5000 | 0.5146 | 103 | 0.5382 | 108 | 50-135 | 36-149 | 4 | 0-25 | |
| Heptachlor | 0.5000 | 0.4826 | 97 | 0.4994 | 100 | 50-135 | 36-149 | 3 | 0-25 | |
| Delta-BHC | 0.5000 | 0.5174 | 103 | 0.5375 | 108 | 50-135 | 36-149 | 4 | 0-25 | |
| Aldrin | 0.5000 | 0.4432 | 89 | 0.4496 | 90 | 50-135 | 36-149 | 1 | 0-25 | |
| Heptachlor Epoxide | 0.5000 | 0.4976 | 100 | 0.5185 | 104 | 50-135 | 36-149 | 4 | 0-25 | |
| Endosulfan I | 0.5000 | 0.4965 | 99 | 0.5141 | 103 | 50-135 | 36-149 | 3 | 0-25 | |
| Dieldrin | 0.5000 | 0.5167 | 103 | 0.5388 | 108 | 50-135 | 36-149 | 4 | 0-25 | |
| 4,4'-DDE | 0.5000 | 0.5024 | 100 | 0.5248 | 105 | 50-135 | 36-149 | 4 | 0-25 | |
| Endrin | 0.5000 | 0.6139 | 123 | 0.6447 | 129 | 50-135 | 36-149 | 5 | 0-25 | |
| Endrin Aldehyde | 0.5000 | 0.4493 | 90 | 0.4706 | 94 | 50-135 | 36-149 | 5 | 0-25 | |
| 4,4'-DDD | 0.5000 | 0.5181 | 104 | 0.5421 | 108 | 50-135 | 36-149 | 5 | 0-25 | |
| Endosulfan II | 0.5000 | 0.5086 | 102 | 0.5294 | 106 | 50-135 | 36-149 | 4 | 0-25 | |
| 4,4'-DDT | 0.5000 | 0.5246 | 105 | 0.5577 | 112 | 50-135 | 36-149 | 6 | 0-25 | |
| Endosulfan Sulfate | 0.5000 | 0.5099 | 102 | 0.5352 | 107 | 50-135 | 36-149 | 5 | 0-25 | |
| Methoxychlor | 0.5000 | 0.5712 | 114 | 0.6055 | 121 | 50-135 | 36-149 | 6 | 0-25 | |

Total number of LCS compounds: 17

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-533-1101 | LCS | Aqueous | GC 31 | 10/29/15 | 10/30/15 13:15 | 151029L14 | | | |
| 099-12-533-1101 | LCSD | Aqueous | GC 31 | 10/29/15 | 10/30/15 13:34 | 151029L14 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Aroclor-1016 | 2.000 | 2.061 | 103 | 2.013 | 101 | 50-135 | 2 | 0-25 | |
| Aroclor-1260 | 2.000 | 2.122 | 106 | 2.189 | 109 | 50-135 | 3 | 0-25 | |

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RPD: Relative Percent Difference. CL: Control Limits



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Quality Control - LCS/LCSD

SWCA Environmental Consultants
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Date Received: 10/27/15
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Preparation: EPA 3510C
Method: EPA 8082

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-12-533-1102 | LCS | Aqueous | GC 31 | 10/29/15 | 10/30/15 14:12 | 151029L15 | | | |
| 099-12-533-1102 | LCSD | Aqueous | GC 31 | 10/29/15 | 10/30/15 14:31 | 151029L15 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Aroclor-1016 | 2.000 | 1.929 | 96 | 1.944 | 97 | 50-135 | 1 | 0-25 | |
| Aroclor-1260 | 2.000 | 2.164 | 108 | 2.242 | 112 | 50-135 | 4 | 0-25 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-963-121 | LCS | Aqueous | | GC 26 | 10/30/15 | 10/30/15 20:33 | 151030L02 | | | |
| 099-15-963-121 | LCSD | Aqueous | | GC 26 | 10/30/15 | 10/30/15 21:17 | 151030L02 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.03548 | 89 | 0.03624 | 91 | 30-130 | 13-147 | 2 | 0-30 | |
| Bolstar | 0.04000 | 0.03344 | 84 | 0.03400 | 85 | 30-130 | 13-147 | 2 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.03030 | 76 | 0.03132 | 78 | 30-130 | 13-147 | 3 | 0-30 | |
| Coumaphos | 0.04000 | 0.03136 | 78 | 0.03214 | 80 | 30-130 | 13-147 | 2 | 0-30 | |
| Diazinon | 0.04000 | 0.03438 | 86 | 0.03575 | 89 | 30-130 | 13-147 | 4 | 0-30 | |
| Disulfoton | 0.04000 | 0.03376 | 84 | 0.03389 | 85 | 30-130 | 13-147 | 0 | 0-30 | |
| Ethoprop | 0.04000 | 0.03390 | 85 | 0.03478 | 87 | 30-130 | 13-147 | 3 | 0-30 | |
| Fensulfothion | 0.04000 | 0.03540 | 88 | 0.03612 | 90 | 30-130 | 13-147 | 2 | 0-30 | |
| Fenthion | 0.04000 | 0.03360 | 84 | 0.03411 | 85 | 30-130 | 13-147 | 2 | 0-30 | |
| Merphos | 0.04000 | 0.02985 | 75 | 0.03243 | 81 | 30-130 | 13-147 | 8 | 0-30 | |
| Methyl Parathion | 0.04000 | 0.03211 | 80 | 0.03305 | 83 | 30-130 | 13-147 | 3 | 0-30 | |
| Phorate | 0.04000 | 0.03878 | 97 | 0.03981 | 100 | 30-130 | 13-147 | 3 | 0-30 | |
| Ronnel | 0.04000 | 0.02918 | 73 | 0.02994 | 75 | 30-130 | 13-147 | 3 | 0-30 | |
| Stirophos | 0.04000 | 0.02682 | 67 | 0.02749 | 69 | 30-130 | 13-147 | 2 | 0-30 | |
| Tokuthion | 0.04000 | 0.02941 | 74 | 0.02983 | 75 | 30-130 | 13-147 | 1 | 0-30 | |
| Trichloronate | 0.04000 | 0.03112 | 78 | 0.03193 | 80 | 30-130 | 13-147 | 3 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8141A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-963-122 | LCS | Aqueous | | GC 26 | 10/30/15 | 10/30/15 22:46 | 151030L03 | | | |
| 099-15-963-122 | LCSD | Aqueous | | GC 26 | 10/30/15 | 10/30/15 23:30 | 151030L03 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Azinphos Methyl | 0.04000 | 0.03277 | 82 | 0.03449 | 86 | 30-130 | 13-147 | 5 | 0-30 | |
| Bolstar | 0.04000 | 0.03204 | 80 | 0.03222 | 81 | 30-130 | 13-147 | 1 | 0-30 | |
| Chlorpyrifos | 0.04000 | 0.02940 | 74 | 0.02958 | 74 | 30-130 | 13-147 | 1 | 0-30 | |
| Coumaphos | 0.04000 | 0.02931 | 73 | 0.03056 | 76 | 30-130 | 13-147 | 4 | 0-30 | |
| Diazinon | 0.04000 | 0.03368 | 84 | 0.03319 | 83 | 30-130 | 13-147 | 1 | 0-30 | |
| Disulfoton | 0.04000 | 0.03242 | 81 | 0.03253 | 81 | 30-130 | 13-147 | 0 | 0-30 | |
| Ethoprop | 0.04000 | 0.03319 | 83 | 0.03292 | 82 | 30-130 | 13-147 | 1 | 0-30 | |
| Fensulfothion | 0.04000 | 0.03365 | 84 | 0.03430 | 86 | 30-130 | 13-147 | 2 | 0-30 | |
| Fenthion | 0.04000 | 0.03205 | 80 | 0.03209 | 80 | 30-130 | 13-147 | 0 | 0-30 | |
| Merphos | 0.04000 | 0.02884 | 72 | 0.03096 | 77 | 30-130 | 13-147 | 7 | 0-30 | |
| Methyl Parathion | 0.04000 | 0.03085 | 77 | 0.03100 | 78 | 30-130 | 13-147 | 0 | 0-30 | |
| Phorate | 0.04000 | 0.03792 | 95 | 0.03742 | 94 | 30-130 | 13-147 | 1 | 0-30 | |
| Ronnel | 0.04000 | 0.02813 | 70 | 0.02833 | 71 | 30-130 | 13-147 | 1 | 0-30 | |
| Stirophos | 0.04000 | 0.02617 | 65 | 0.02631 | 66 | 30-130 | 13-147 | 1 | 0-30 | |
| Tokuthion | 0.04000 | 0.02820 | 70 | 0.02826 | 71 | 30-130 | 13-147 | 0 | 0-30 | |
| Trichloronate | 0.04000 | 0.03008 | 75 | 0.03011 | 75 | 30-130 | 13-147 | 0 | 0-30 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-034-667 | LCS | Aqueous | GC 40 | 10/30/15 | 11/05/15 19:46 | 151030L19 | | | |
| 095-01-034-667 | LCSD | Aqueous | GC 40 | 10/30/15 | 11/05/15 20:09 | 151030L19 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| 2,4-D | 20.00 | 15.48 | 77 | 17.12 | 86 | 30-130 | 10 | 0-30 | |
| 2,4,5-T | 2.000 | 2.010 | 100 | 2.115 | 106 | 30-130 | 5 | 0-30 | |
| 2,4-DB | 20.00 | 17.02 | 85 | 18.42 | 92 | 30-130 | 8 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 8151A
Method: EPA 8151A

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number |
|---------------------------|------|---------|------------|---------------|----------------|-----------------------|
| 095-01-034-668 | LCS | Aqueous | GC 40 | 10/30/15 | 11/05/15 20:55 | 151030L20 |
| 095-01-034-668 | LCSD | Aqueous | GC 40 | 10/30/15 | 11/05/15 21:18 | 151030L20 |

| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|-------------|-----------|-----------|------------|------------|----------|-----|--------|------------|
| 2,4-D | 20.00 | 15.80 | 79 | 17.11 | 86 | 30-130 | 8 | 0-30 | |
| 2,4,5-T | 2.000 | 2.060 | 103 | 2.170 | 108 | 30-130 | 5 | 0-30 | |
| 2,4-DB | 20.00 | 17.16 | 86 | 18.44 | 92 | 30-130 | 7 | 0-30 | |

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RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-4133 | LCS | Aqueous | | GC/MS SS | 10/30/15 | 10/31/15 15:40 | 151030L04 | | | |
| 095-01-003-4133 | LCSD | Aqueous | | GC/MS SS | 10/30/15 | 10/31/15 16:00 | 151030L04 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 200.0 | 153.2 | 77 | 152.6 | 76 | 61-120 | 51-130 | 0 | 0-20 | |
| Acenaphthylene | 200.0 | 150.4 | 75 | 149.8 | 75 | 55-120 | 44-131 | 0 | 0-20 | |
| Butyl Benzyl Phthalate | 200.0 | 169.4 | 85 | 166.7 | 83 | 56-122 | 45-133 | 2 | 0-20 | |
| 4-Chloro-3-Methylphenol | 200.0 | 175.1 | 88 | 172.8 | 86 | 52-120 | 41-131 | 1 | 0-20 | |
| 2-Chlorophenol | 200.0 | 168.7 | 84 | 164.8 | 82 | 47-120 | 35-132 | 2 | 0-20 | |
| 1,4-Dichlorobenzene | 200.0 | 149.5 | 75 | 147.6 | 74 | 36-120 | 22-134 | 1 | 0-20 | |
| Dimethyl Phthalate | 200.0 | 153.6 | 77 | 152.5 | 76 | 60-120 | 50-130 | 1 | 0-20 | |
| 2,4-Dinitrotoluene | 200.0 | 170.9 | 85 | 168.9 | 84 | 61-121 | 51-131 | 1 | 0-20 | |
| Fluorene | 200.0 | 152.9 | 76 | 154.2 | 77 | 67-120 | 58-129 | 1 | 0-20 | |
| N-Nitroso-di-n-propylamine | 200.0 | 160.8 | 80 | 154.3 | 77 | 39-123 | 25-137 | 4 | 0-20 | |
| Naphthalene | 200.0 | 142.2 | 71 | 142.3 | 71 | 54-120 | 43-131 | 0 | 0-20 | |
| 4-Nitrophenol | 200.0 | 167.5 | 84 | 171.7 | 86 | 14-120 | 0-138 | 3 | 0-20 | |
| Pentachlorophenol | 200.0 | 151.2 | 76 | 144.8 | 72 | 31-127 | 15-143 | 4 | 0-20 | |
| Phenol | 200.0 | 164.8 | 82 | 159.9 | 80 | 17-120 | 0-137 | 3 | 0-20 | |
| Pyrene | 200.0 | 155.6 | 78 | 153.2 | 77 | 58-124 | 47-135 | 2 | 0-20 | |
| 1,2,4-Trichlorobenzene | 200.0 | 149.4 | 75 | 147.1 | 74 | 49-120 | 37-132 | 2 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 3510C
Method: EPA 8270C

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 095-01-003-4132 | LCS | Aqueous | | GC/MS TT | 10/30/15 | 10/31/15 21:01 | 151030L05 | | | |
| 095-01-003-4132 | LCSD | Aqueous | | GC/MS TT | 10/30/15 | 10/31/15 21:20 | 151030L05 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | ME CL | RPD | RPD CL | Qualifiers |
| Acenaphthene | 200.0 | 152.3 | 76 | 149.1 | 75 | 61-120 | 51-130 | 2 | 0-20 | |
| Acenaphthylene | 200.0 | 147.1 | 74 | 144.1 | 72 | 55-120 | 44-131 | 2 | 0-20 | |
| Butyl Benzyl Phthalate | 200.0 | 146.7 | 73 | 140.7 | 70 | 56-122 | 45-133 | 4 | 0-20 | |
| 4-Chloro-3-Methylphenol | 200.0 | 147.7 | 74 | 140.6 | 70 | 52-120 | 41-131 | 5 | 0-20 | |
| 2-Chlorophenol | 200.0 | 151.5 | 76 | 151.2 | 76 | 47-120 | 35-132 | 0 | 0-20 | |
| 1,4-Dichlorobenzene | 200.0 | 145.2 | 73 | 144.4 | 72 | 36-120 | 22-134 | 1 | 0-20 | |
| Dimethyl Phthalate | 200.0 | 144.3 | 72 | 141.6 | 71 | 60-120 | 50-130 | 2 | 0-20 | |
| 2,4-Dinitrotoluene | 200.0 | 159.4 | 80 | 156.6 | 78 | 61-121 | 51-131 | 2 | 0-20 | |
| Fluorene | 200.0 | 152.6 | 76 | 148.9 | 74 | 67-120 | 58-129 | 2 | 0-20 | |
| N-Nitroso-di-n-propylamine | 200.0 | 145.4 | 73 | 143.8 | 72 | 39-123 | 25-137 | 1 | 0-20 | |
| Naphthalene | 200.0 | 147.3 | 74 | 143.4 | 72 | 54-120 | 43-131 | 3 | 0-20 | |
| 4-Nitrophenol | 200.0 | 146.4 | 73 | 140.6 | 70 | 14-120 | 0-138 | 4 | 0-20 | |
| Pentachlorophenol | 200.0 | 153.9 | 77 | 150.2 | 75 | 31-127 | 15-143 | 2 | 0-20 | |
| Phenol | 200.0 | 153.6 | 77 | 150.4 | 75 | 17-120 | 0-137 | 2 | 0-20 | |
| Pyrene | 200.0 | 146.5 | 73 | 144.7 | 72 | 58-124 | 47-135 | 1 | 0-20 | |
| 1,2,4-Trichlorobenzene | 200.0 | 144.6 | 72 | 142.8 | 71 | 49-120 | 37-132 | 1 | 0-20 | |

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-18641 | LCS | Aqueous | GC/MS O | 10/30/15 | 10/30/15 11:23 | 151030L006 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 51.34 | 103 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 55.65 | 111 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 53.08 | 106 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 44.29 | 89 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 47.54 | 95 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 43.88 | 88 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 54.03 | 108 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 56.52 | 113 | 80-123 | 73-130 | |
| Toluene | 50.00 | 50.59 | 101 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 51.16 | 102 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 45.35 | 91 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 119.1 | 119 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 57.37 | 115 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 42.34 | 85 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-18638 | LCS | Aqueous | GC/MS JJ | 10/30/15 | 10/30/15 09:22 | 151030L001 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 44.86 | 90 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 45.83 | 92 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 49.75 | 99 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 47.32 | 95 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 50.57 | 101 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 44.90 | 90 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 49.31 | 99 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 51.39 | 103 | 80-123 | 73-130 | |
| Toluene | 50.00 | 46.04 | 92 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 46.01 | 92 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 43.42 | 87 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 103.5 | 104 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 50.82 | 102 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 40.02 | 80 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS

SWCA Environmental Consultants
6200 UTSA Blvd., Suite 102
San Antonio, TX 78249-1618

Date Received: 10/27/15
Work Order: 15-10-1995
Preparation: EPA 5030C
Method: EPA 8260B

Project: EAA 27122

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| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|-----------------------------|-------------|-----------------|------------|---------------|----------------|------------------|
| 099-14-001-18642 | LCS | Aqueous | GC/MS QQ | 10/30/15 | 10/30/15 11:44 | 151030L010 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | %Rec. CL | ME CL | Qualifiers |
| Benzene | 50.00 | 47.31 | 95 | 80-120 | 73-127 | |
| Carbon Tetrachloride | 50.00 | 56.26 | 113 | 67-139 | 55-151 | |
| Chlorobenzene | 50.00 | 47.56 | 95 | 78-120 | 71-127 | |
| 1,2-Dibromoethane | 50.00 | 52.71 | 105 | 80-120 | 73-127 | |
| 1,2-Dichlorobenzene | 50.00 | 50.77 | 102 | 63-129 | 52-140 | |
| 1,2-Dichloroethane | 50.00 | 47.64 | 95 | 70-130 | 60-140 | |
| 1,1-Dichloroethene | 50.00 | 33.34 | 67 | 66-126 | 56-136 | |
| Ethylbenzene | 50.00 | 46.63 | 93 | 80-123 | 73-130 | |
| Toluene | 50.00 | 46.98 | 94 | 80-120 | 73-127 | |
| Trichloroethene | 50.00 | 45.97 | 92 | 80-122 | 73-129 | |
| Vinyl Chloride | 50.00 | 37.64 | 75 | 70-130 | 60-140 | |
| p/m-Xylene | 100.0 | 88.45 | 88 | 75-123 | 67-131 | |
| o-Xylene | 50.00 | 45.28 | 91 | 74-122 | 66-130 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 49.54 | 99 | 69-129 | 59-139 | |

Total number of LCS compounds: 14

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Sample Analysis Summary Report

Work Order: 15-10-1995

Page 1 of 1

| <u>Method</u> | <u>Extraction</u> | <u>Chemist ID</u> | <u>Instrument</u> | <u>Analytical Location</u> |
|-----------------|-------------------|-------------------|-------------------|----------------------------|
| EPA 300.0 | N/A | 1027 | IC 15 | 1 |
| EPA 365.1 | N/A | 735 | ACA 1 | 1 |
| EPA 6010B | EPA 3005A Filt. | 935 | ICP 7300 | 1 |
| EPA 6020 | EPA 3005A Filt. | 598 | ICP/MS 03 | 1 |
| EPA 7470A | EPA 7470A Filt. | 915 | Mercury 04 | 1 |
| EPA 8081A | EPA 3510C | 669 | GC 44 | 1 |
| EPA 8082 | EPA 3510C | 669 | GC 31 | 1 |
| EPA 8141A | EPA 3510C | 960 | GC 26 | 1 |
| EPA 8151A | EPA 8151A | 669 | GC 40 | 1 |
| EPA 8260B | EPA 5030C | 486 | GC/MS JJ | 2 |
| EPA 8260B | EPA 5030C | 486 | GC/MS QQ | 2 |
| EPA 8260B | EPA 5030C | 867 | GC/MS O | 2 |
| EPA 8260B | EPA 5030C | 927 | GC/MS JJ | 2 |
| EPA 8270C | EPA 3510C | 923 | GC/MS SS | 1 |
| EPA 8270C | EPA 3510C | 923 | GC/MS TT | 1 |
| SM 2320B | N/A | 650 | PH1/BUR03 | 1 |
| SM 2320B | N/A | 688 | PH1/BUR03 | 1 |
| SM 2540 C | N/A | 689 | N/A | 1 |
| SM 2540 D | N/A | 689 | N/A | 1 |
| SM 4500 H+ B | N/A | 650 | PH 1 | 1 |
| SM 4500 N Org B | N/A | 685 | BUR05 | 1 |
| SM 5310 B | N/A | 735 | TOC 8 | 1 |
| SM 5310 B | N/A | 735 | TOC 11 | 1 |

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Location 1: 7440 Lincoln Way, Garden Grove, CA 92841

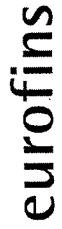
Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841

Glossary of Terms and Qualifiers

Work Order: 15-10-1995

Page 1 of 1

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|---|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| CI | See case narrative. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |
| | Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time. |
| | A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations. |



CHAIN OF CUSTODY RECORD

5

**Jennifer Moreland, Billary Rios,
Philip Pearce**

06/02/14 Revision

24 of 25
MPS# 7748 2142 1489
0263
Mstr# 7748 2141 9066

A7 APVA



2 of 25
MPS# 7748 2141 9272
0263
Mstr# 7748 2141 9066

A7 APVA

4 of 25
MPS# 7748 2141 9640
0263
Mstr# 7748 2141 9066

A7 APVA

3 of 25
MPS# 7748 2141 9342
0263
Mstr# 7748 2141 9066

A7 APVA

25 of 25
MPS# 7748 2142 1684
0263
Mstr# 7748 2141 9066

A7 APVA

22 of 25
MPS# 7748 2142 2007
0263
Mstr# 7748 2141 9066

A7 APVA

13 of 25
MPS# 7748 2142 0839
0263
Mstr# 7748 2141 9066

A7 APVA



19 of 25
MPS# 7748 2142 1055
0263
Mstr# 7748 2141 9066

A7 APVA



TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
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TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
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TUE - 27 OCT 3:00P
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92841
CA-US SNA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

14 of 25
MPS# 7748 2142 0655
0263
Mstr# 7748 2141 9066

A7 APVA

10 of 25
MPS# 7748 2142 0530
0263
Mstr# 7748 2141 9066

A7 APVA

11 of 25
MPS# 7748 2142 0552
0263
Mstr# 7748 2141 9066

A7 APVA

1 of 25
TRK# 7748 2141 9066
0201
MASTER

A7 APVA

1 of 8
TRK# 7748 2288 5010
0201
MASTER

A7 APVA

9 of 25
MPS# 7748 2142 0081
0263
Mstr# 7748 2141 9066

A7 APVA

5 of 25
MPS# 7748 2141 9905
0263
Mstr# 7748 2141 9066

A7 APVA

18 of 25
MPS# 7748 2142 1283
0263
Mstr# 7748 2141 9066

A7 APVA



TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
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CA-US SNA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
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CA-US SNA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
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TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
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CA-US SNA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

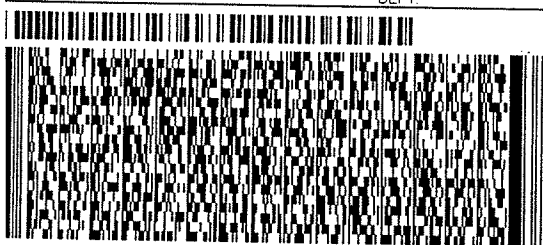
TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

PHILIP C. PEARCE, P.G.
SWCA
6200 UTSA BLVD
STE 102
SAN ANTONIO, TX 78249
UNITED STATES US

SHIP DATE: 2000113
ACTWGT: 70.00 LB
CAD: 8347991/INET3670
DIMS: 24x13x14 IN
BILL SENDER

LAB
EUROFINS CALSCIENCE
7440 LINCOLN WAY

GARDEN GROVE CA 92841
(714) 895-5494 REF: 27122.0201
INV: PO: DEPT:



FedEx
Express



8 of 25
MPS# 7748 2142 0060
0263
Mstr# 7748 2141 9066

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

7 of 25
MPS# 7748 2141 9489
0263
Mstr# 7748 2141 9066

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

8 of 8
MPS# 7748 2288 5866
0263
Mstr# 7748 2288 5010

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

20 of 25
IPS# 7748 2142 1066
0263
Mstr# 7748 2141 9066

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA



6 of 8
MPS# 7748 2288 5616
0263
Mstr# 7748 2288 5010

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

3 of 8
MPS# 7748 2288 5215
0263
Mstr# 7748 2288 5010

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

5 of 8
MPS# 7748 2288 5395
0263
Mstr# 7748 2288 5010

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

4 of 8
MPS# 7748 2288 4860
0263
Mstr# 7748 2288 5010

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841

6 of 25
MPS# 7748 2141 9880
0263
Mstr# 7748 2141 9066

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

2 of 8
MPS# 7748 2288 4734
0263
Mstr# 7748 2288 5010

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

23 of 25
MPS# 7748 2142 2018
0263
Mstr# 7748 2141 9066

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

16 of 25
MPS# 7748 2142 1228
0263
Mstr# 7748 2141 9066

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

15 of 25
MPS# 7748 2142 0482
0263
Mstr# 7748 2141 9066

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

21 of 25
MPS# 7748 2142 1560
0263
Mstr# 7748 2141 9066

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841

17 of 25
MPS# 7748 2142 0986
0263
Mstr# 7748 2141 9066

A7 APVA

TUE - 27 OCT 3:00P
STANDARD OVERNIGHT
ASR
92841
CA-US SNA

SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 2.0 °C (w/ CF): 1.6 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Present but Not Intact☐ Not Present☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact☐ Present but Not Intact☒ Not Present☐ N/A

Checked by: 107

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples

Yes ☒ No ☐ N/A ☐

COC document(s) received complete

Yes ☒ No ☐ N/A ☐☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC

Yes ☒ No ☐ N/A ☐

Sample container label(s) consistent with COC

Yes ☒ No ☐ N/A ☐

Sample container(s) intact and in good condition

Yes ☒ No ☐ N/A ☐

Proper containers for analyses requested

Yes ☒ No ☐ N/A ☐

Sufficient volume/mass for analyses requested

Yes ☒ No ☐ N/A ☐

Samples received within holding time

Yes ☐ No ☒ N/A ☐

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved OxygenYes ☐ No ☒ N/A ☐

Proper preservation chemical(s) noted on COC and/or sample container

Yes ☒ No ☐ N/A ☐

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace

Yes ☒ No ☐ N/A ☐☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation

Yes ☐ No ☐ N/A ☒

CONTAINER TYPE:

(Trip Blank Lot Number: 150909E)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na2} ☒ 100PJ ☐ 100PJ_{na2} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s☐ 500PB ☒ 1AGB ☐ 1AGB_{na2} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5gal ☐ ☐ ☐Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® (____) ☐ TerraCores® (____) ☐Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ Other Matrix (____): ☐ ☐

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄,

Labeled/Checked by: 107

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 2 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 2.1 °C (w/ CF): 1.7 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 1017

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)-

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na2} ☐ 100PJ ☐ 100PJ_{na2} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s
☐ 500PB ☒ 1AGB ☐ 1AGB_{na2} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 25g_{ul} ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 3 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 3.0 °C (w/ CF): 2.6 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact☐ Present but Not Intact☐ Not Present☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact☐ Present but Not Intact☒ Not Present☐ N/A

Checked by: 1017

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples

Yes

No

N/A

☒☐☐

COC document(s) received complete

☒☐☐☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC

☒☐☐

Sample container label(s) consistent with COC

☒☐☐

Sample container(s) intact and in good condition

☒☐☐

Proper containers for analyses requested

☒☐☐

Sufficient volume/mass for analyses requested

☒☐☐

Samples received within holding time

☐☒☐

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen☐☒☐

Proper preservation chemical(s) noted on COC and/or sample container

☒☐☐

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace

☒☐☐☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation

☐☐☒

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☐ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 25g ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄

Labeled/Checked by: 1017

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 4 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 1.5 °C (w/ CF): 1.1 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 107

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)- | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s

☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 25gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 107

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 5 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 1.7 °C (w/ CF): 1.3 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 107

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGBs ☐ 1PB ☐ 1PB_{na} ☒ 25g_{ul} ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 107

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 6 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 3.1 °C (w/ CF): 2.7 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact

☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact

☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 1017

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)- | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☐ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 25gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017

s = H₂SO₄, u = ultra-pure, z₂na = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 7 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 2.2 °C (w/ CF): 1.8 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 107

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)- | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_z ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s

☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 25g ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 107

s = H₂SO₄, u = ultra-pure, z_{na} = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 8 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 2.9 °C (w/ CF): 2.5 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 107

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)-

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s
☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 107

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 9 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 2.9 °C (w/ CF): 2.5 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)- | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☐ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 25gal ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃COO)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 10 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 3.3 °C (w/ CF): 2.9 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 1017

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s
☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 11 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 3.2 °C (w/ CF): 2.8 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 107

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s
☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 25gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 107

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 12 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 3.0 °C (w/ CF): 2.6 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 107

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)-

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☐ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 25gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 107

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 13 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 2.1 °C (w/ CF): 1.7 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s

☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 25g_{ul} ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 14 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.4°C); Temperature (w/o CF): 2.3 °C (w/ CF): 1.9 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 107

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s
☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 25g_{ul} ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 107

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 15 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.4°C); Temperature (w/o CF): 1.4 °C (w/ CF): 1.0 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5gal ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 16 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 2.5 °C (w/ CF): 2.1 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 1017

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)- | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na2} ☐ 100PJ ☐ 100PJ_{na2} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s

☐ 500PB ☒ 1AGB ☐ 1AGB_{na2} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 17 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.4°C); Temperature (w/o CF): 2.6 °C (w/ CF): 2.2 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB
☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s
☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 25gal ☐ _____ ☐ _____
Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____
Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 197

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 18 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.4°C); Temperature (w/o CF): 2.5 °C (w/ CF): 2.1 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)- | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☐ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 25gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 19 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 1.9 °C (w/ CF): 1.5 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOAn₂ ☐ 100PJ ☐ 100PJn₂ ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBn₂ ☒ 1AGBs ☐ 1PB ☐ 1PBn₂ ☒ 25g_{ul} ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® (____) ☐ TerraCores® (____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 20 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 2.0 °C (w/ CF): 1.6 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact

☐ Not Present

☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact

☒ Not Present

☐ N/A

Checked by: 107

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)- | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s

☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 25gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 107

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 21 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 1.9 °C (w/ CF): 1.5 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE: (Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s

☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 25gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 22 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.4°C); Temperature (w/o CF): 1.8 °C (w/ CF): 1.4 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 107

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 25g_{ul} ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 107s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 23 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 2.5 °C (w/ CF): 2.1 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 107

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)- | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s

☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5g/L ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 107

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 24 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.4°C); Temperature (w/o CF): 2.1 °C (w/ CF): 1.7 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)-Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☐ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 2.5gal ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 25 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.4°C); Temperature (w/o CF): 2.3 °C (w/ CF): 1.9 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)- | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☐ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5gal ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 26 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.4°C); Temperature (w/o CF): 2.5 °C (w/ CF): 2.1 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBz_{na} ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☐ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 2.5gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017

s = H₂SO₄, u = ultra-pure, z_{na} = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300



Calscience

WORK ORDER NUMBER: 15-10-1995

SAMPLE RECEIPT CHECKLIST

COOLER 27 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 2-4 °C (w/ CF): 2-0 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/ACOC document(s) received complete ☒ Yes ☐ No ☐ N/A☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)-Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB☐ 125PBz₂na ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☐ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 25g/L ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017s = H₂SO₄, u = ultra-pure, z₂na = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 28 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.4°C); Temperature (w/o CF): 3.0 °C (w/ CF): 2.6 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 107

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)- | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOA_{na} ☒ 100PJ ☐ 100PJ_{na} ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB☐ 125PB_{znna} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AG_J ☐ 500AG_J_s☐ 500PB ☒ 1AGB ☐ 1AGB_{na} ☒ 1AGB_s ☐ 1PB ☐ 1PB_{na} ☒ 25gal ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 107s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300



Calscience

WORK ORDER NUMBER: 15-10-1995

SAMPLE RECEIPT CHECKLIST

COOLER 29 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 2.3 °C (w/ CF): 1.9 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 107

SAMPLE CONDITION:

| | Yes | No | N/A |
|---|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)- | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOAn₂ ☐ 100PJ ☐ 100PJn₂ ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB☐ 125PBznna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☐ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs☐ 500PB ☒ 1AGB ☐ 1AGBn₂ ☒ 1AGBs ☐ 1PB ☐ 1PBn₂ ☒ 25g_{ul} ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (____) ☐ EnCores® (____) ☐ TerraCores® (____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (____): ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄,

Labeled/Checked by: 107

s = H₂SO₄, u = ultra-pure, znna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 30 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF: -0.4°C); Temperature (w/o CF): 2.2 °C (w/ CF): 1.8 °C; ☒ Blank ☐ Sample☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling☐ Sample(s) received at ambient temperature; placed on ice for transport by courierAmbient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|--------------------------|--------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished timeSampler's name indicated on COC ☒ Yes ☐ No ☐ N/ASample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/ASample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/AProper containers for analyses requested ☒ Yes ☐ No ☐ N/ASufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/ASamples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/AProper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved MetalsContainer(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)-Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOAn_a ☐ 100PJ ☐ 100PJn_a ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB☐ 125PBz_{nn} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s☐ 500PB ☒ 1AGB ☐ 1AGBn_a ☒ 1AGB_s ☐ 1PB ☐ 1PBn_a ☒ 25gal ☐ _____ ☐ _____ ☐ _____Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017s = H₂SO₄, u = ultra-pure, z_{nn}a = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 31 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 1.9 °C (w/ CF): 1.5 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 107

SAMPLE CONDITION:

Chain-of-Custody (COC) document(s) received with samples ☒ Yes ☐ No ☐ N/A

COC document(s) received complete ☒ Yes ☐ No ☐ N/A

☐ Sampling date ☐ Sampling time ☐ Matrix ☐ Number of containers

☐ No analysis requested ☐ Not relinquished ☐ No relinquished date ☐ No relinquished time

Sampler's name indicated on COC ☒ Yes ☐ No ☐ N/A

Sample container label(s) consistent with COC ☒ Yes ☐ No ☐ N/A

Sample container(s) intact and in good condition ☒ Yes ☐ No ☐ N/A

Proper containers for analyses requested ☒ Yes ☐ No ☐ N/A

Sufficient volume/mass for analyses requested ☒ Yes ☐ No ☐ N/A

Samples received within holding time ☐ Yes ☒ No ☐ N/A

Aqueous samples for certain analyses received within 15-minute holding time

☒ pH ☐ Residual Chlorine ☐ Dissolved Sulfide ☐ Dissolved Oxygen ☐ Yes ☒ No ☐ N/A

Proper preservation chemical(s) noted on COC and/or sample container ☒ Yes ☐ No ☐ N/A

Unpreserved aqueous sample(s) received for certain analyses

☐ Volatile Organics ☐ Total Metals ☐ Dissolved Metals

Container(s) for certain analysis free of headspace ☒ Yes ☐ No ☐ N/A

☒ Volatile Organics ☐ Dissolved Gases (RSK-175) ☐ Dissolved Oxygen (SM 4500)

☐ Carbon Dioxide (SM 4500) ☐ Ferrous Iron (SM 3500) ☐ Hydrogen Sulfide (Hach)-

Tedlar™ bag(s) free of condensation ☐ Yes ☐ No ☒ N/A

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOAh ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGBh ☐ 125AGBp ☐ 125PB

☐ 125PBznnna ☐ 250AGB ☐ 250CGB ☒ 250CGBs ☐ 250PB ☐ 250PBn ☐ 500AGB ☐ 500AGJ ☐ 500AGJs

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGBs ☐ 1PB ☐ 1PBna ☒ 25g/L ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 107

s = H₂SO₄, u = ultra-pure, znnna = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE RECEIPT CHECKLIST

COOLER 32 OF 32

CLIENT: SWCA

DATE: 10 / 27 / 2015

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)

Thermometer ID: SC2 (CF:-0.4°C); Temperature (w/o CF): 2.0 °C (w/ CF): 1.6 °C; ☒ Blank ☐ Sample

☐ Sample(s) outside temperature criteria (PM/APM contacted by: _____)

☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling

☐ Sample(s) received at ambient temperature; placed on ice for transport by courier

Ambient Temperature: ☐ Air ☐ Filter

Checked by: 15

CUSTODY SEAL:

Cooler ☒ Present and Intact ☐ Present but Not Intact ☐ Not Present ☐ N/A

Checked by: 15

Sample(s) ☐ Present and Intact ☐ Present but Not Intact ☒ Not Present ☐ N/A

Checked by: 1017

SAMPLE CONDITION:

| | Yes | No | N/A |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-of-Custody (COC) document(s) received with samples | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers | | | |
| <input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time | | | |
| Sampler's name indicated on COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and in good condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sufficient volume/mass for analyses requested | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Samples received within holding time | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples for certain analyses received within 15-minute holding time | | | |
| <input checked="" type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper preservation chemical(s) noted on COC and/or sample container | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Unpreserved aqueous sample(s) received for certain analyses | | | |
| <input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals | | | |
| Container(s) for certain analysis free of headspace | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500) | | | |
| <input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach) | | | |
| Tedlar™ bag(s) free of condensation | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: ☐ VOA ☒ VOA_h ☐ VOAna₂ ☐ 100PJ ☐ 100PJna₂ ☐ 125AGB ☐ 125AGB_h ☐ 125AGB_p ☐ 125PB

☐ 125PBz_{na} ☐ 250AGB ☐ 250CGB ☒ 250CGB_s ☐ 250PB ☐ 250PB_n ☐ 500AGB ☐ 500AGJ ☐ 500AGJ_s

☐ 500PB ☒ 1AGB ☐ 1AGBna₂ ☒ 1AGB_s ☐ 1PB ☐ 1PBna ☒ 25gal ☐ _____ ☐ _____ ☐ _____

Solid: ☐ 4ozCGJ ☐ 8ozCGJ ☐ 16ozCGJ ☐ Sleeve (_____) ☐ EnCores® (_____) ☐ TerraCores® (_____) ☐ _____

Air: ☐ Tedlar™ ☐ Canister ☐ Sorbent Tube ☐ PUF ☐ _____ Other Matrix (_____) ☐ _____ ☐ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1017

s = H₂SO₄, u = ultra-pure, z_{na} = Zn(CH₃CO₂)₂ + NaOH

Reviewed by: 300

SAMPLE ANOMALY REPORT

DATE: 10 / 27 / 2015

SAMPLES, CONTAINERS, AND LABELS:

- ☒ Sample(s) NOT RECEIVED but listed on COC
- ☐ Sample(s) received but NOT LISTED on COC
- ☒ Holding time expired (list client or ECI sample ID and analysis)
- ☐ Insufficient sample amount for requested analysis (list analysis)
- ☐ Improper container(s) used (list analysis)
- ☐ Improper preservative used (list analysis)
- ☐ No preservative noted on COC or label (list analysis and notify lab)
- ☐ Sample container(s) not labeled
- ☐ Client sample label(s) illegible (list container type and analysis)
- ☐ Client sample label(s) do not match COC (comment)
 - ☐ Project information
 - ☐ Client sample ID
 - ☐ Sampling date and/or time
 - ☐ Number of container(s)
 - ☐ Requested analysis
- ☐ Sample container(s) compromised (comment)
 - ☐ Broken
 - ☐ Water present in sample container
- ☐ Air sample container(s) compromised (comment)
 - ☐ Flat
 - ☐ Very low in volume
 - ☐ Leaking (not transferred; duplicate bag submitted)
 - ☐ Leaking (transferred into ECI Tedlar™ bags*)
 - ☐ Leaking (transferred into client's Tedlar™ bags*)

* Transferred at client's request.

MISCELLANEOUS: (Describe)

HEADSPACE:

(Containers with bubble > 6 mm or ¼ inch for volatile organic or dissolved gas analysis)

| ECI Sample ID | ECI Container ID | Total Number** | ECI Sample ID | ECI Container ID | Total Number** |
|---------------|------------------|----------------|---------------|------------------|----------------|
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

Comments

(-32)HSM 250 Peak not received

pH and NO₃ received past holding time.

Comments

(Containers with bubble for other analysis)

| ECI Sample ID | ECI Container ID | Total Number** | Requested Analysis |
|---------------|------------------|----------------|--------------------|
| | | | |
| | | | |
| | | | |
| | | | |

Comments: _____

** Record the total number of containers (i.e., vials or bottles) for the affected sample.

Reported by: 1017

Reviewed by: 300



Calscience

Subcontractor Analysis Report

Work Order: 15-10-1995

Page 1 of 1

One or more samples in this work order have tests that were subcontracted. The subcontract report(s) follows.

For subcontracted tests, please reference the laboratory information noted below.

1. ALS - Columbia Analytical Services, Inc. - Kelso, WA CA ELAP 2286, NELAP WA100010
Method 1694 Caffeine

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Analytical Report for Service Request No:

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RE: 15-10-1995

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ALS Group USA, Corp. dba ALS Environmental

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Acronyms

Qualifiers

State Certifications, Accreditations, And Licenses

Case Narrative

Chain of Custody

Steroids and Endocrine Disrupting Compounds

Return to Contents

Acronyms

| | |
|------------|--|
| ASTM | American Society for Testing and Materials |
| A2LA | American Association for Laboratory Accreditation |
| CARB | California Air Resources Board |
| CAS Number | Chemical Abstract Service registry Number |
| CFC | Chlorofluorocarbon |
| CFU | Colony-Forming Unit |
| DEC | Department of Environmental Conservation |
| DEQ | Department of Environmental Quality |
| DHS | Department of Health Services |
| DOE | Department of Ecology |
| DOH | Department of Health |
| EPA | U. S. Environmental Protection Agency |
| ELAP | Environmental Laboratory Accreditation Program |
| GC | Gas Chromatography |
| GC/MS | Gas Chromatography/Mass Spectrometry |
| LOD | Limit of Detection |
| LOQ | Limit of Quantitation |
| LUFT | Leaking Underground Fuel Tank |
| M | Modified |
| MCL | Maximum Contaminant Level is the highest permissible concentration of a substance allowed in drinking water as established by the USEPA. |
| MDL | Method Detection Limit |
| MPN | Most Probable Number |
| MRL | Method Reporting Limit |
| NA | Not Applicable |
| NC | Not Calculated |
| NCASI | National Council of the Paper Industry for Air and Stream Improvement |
| ND | Not Detected |
| NIOSH | National Institute for Occupational Safety and Health |
| PQL | Practical Quantitation Limit |
| RCRA | Resource Conservation and Recovery Act |
| SIM | Selected Ion Monitoring |
| TPH | Total Petroleum Hydrocarbons |
| tr | Trace level is the concentration of an analyte that is less than the PQL but greater than or equal to the MDL. |

Inorganic Data Qualifiers

- * The result is an outlier. See case narrative.
- # The control limit criteria is not applicable. See case narrative.
- B The analyte was found in the associated method blank at a level that is significant relative to the sample result as defined by the DOD or NELAC standards.
- E The result is an estimate amount because the value exceeded the instrument calibration range.
- J The result is an estimated value.
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.
DOD-QSM 4.2 definition : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- i The MRL/MDL or LOQ/LOD is elevated due to a matrix interference.
- X See case narrative.
- Q See case narrative. One or more quality control criteria was outside the limits.
- H The holding time for this test is immediately following sample collection. The samples were analyzed as soon as possible after receipt by the laboratory.

Metals Data Qualifiers

- # The control limit criteria is not applicable. See case narrative.
- J The result is an estimated value.
- E The percent difference for the serial dilution was greater than 10%, indicating a possible matrix interference in the sample.
- M The duplicate injection precision was not met.
- N The Matrix Spike sample recovery is not within control limits. See case narrative.
- S The reported value was determined by the Method of Standard Additions (MSA).
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.
DOD-QSM 4.2 definition : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- W The post-digestion spike for furnace AA analysis is out of control limits, while sample absorbance is less than 50% of spike absorbance.
- i The MRL/MDL or LOQ/LOD is elevated due to a matrix interference.
- X See case narrative.
- + The correlation coefficient for the MSA is less than 0.995.
- Q See case narrative. One or more quality control criteria was outside the limits.

Organic Data Qualifiers

- * The result is an outlier. See case narrative.
- # The control limit criteria is not applicable. See case narrative.
- A A tentatively identified compound, a suspected aldol-condensation product.
- B The analyte was found in the associated method blank at a level that is significant relative to the sample result as defined by the DOD or NELAC standards.
- C The analyte was qualitatively confirmed using GC/MS techniques, pattern recognition, or by comparing to historical data.
- D The reported result is from a dilution.
- E The result is an estimated value.
- J The result is an estimated value.
- N The result is presumptive. The analyte was tentatively identified, but a confirmation analysis was not performed.
- P The GC or HPLC confirmation criteria was exceeded. The relative percent difference is greater than 40% between the two analytical results.
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.
DOD-QSM 4.2 definition : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- i The MRL/MDL or LOQ/LOD is elevated due to a chromatographic interference.
- X See case narrative.
- Q See case narrative. One or more quality control criteria was outside the limits.

Additional Petroleum Hydrocarbon Specific Qualifiers

- F The chromatographic fingerprint of the sample matches the elution pattern of the calibration standard.
- L The chromatographic fingerprint of the sample resembles a petroleum product, but the elution pattern indicates the presence of a greater amount of lighter molecular weight constituents than the calibration standard.
- H The chromatographic fingerprint of the sample resembles a petroleum product, but the elution pattern indicates the presence of a greater amount of heavier molecular weight constituents than the calibration standard.
- O The chromatographic fingerprint of the sample resembles an oil, but does not match the calibration standard.
- Y The chromatographic fingerprint of the sample resembles a petroleum product eluting in approximately the correct carbon range, but the elution pattern does not match the calibration standard.
- Z The chromatographic fingerprint does not resemble a petroleum product.

ALS Group USA Corp. dba ALS Environmental (ALS) - Kelso
State Certifications, Accreditations, and Licenses

| Agency | Web Site | Number |
|--------------------------|---|---------------|
| Alaska DEC UST | http://dec.alaska.gov/applications/eh/ehllabreports/USTLabs.aspx | UST-040 |
| Arizona DHS | http://www.azdhs.gov/lab/license/env.htm | AZ0339 |
| Arkansas - DEQ | http://www.adeq.state.ar.us/techsvs/labcert.htm | 88-0637 |
| California DHS (ELAP) | http://www.cdph.ca.gov/certlic/labs/Pages/ELAP.aspx | 2795 |
| DOD ELAP | http://www.denix.osd.mil/edqw/Accreditation/AccreditedLabs.cfm | L14-51 |
| Florida DOH | http://www.doh.state.fl.us/lab/EnvLabCert/WaterCert.htm | E87412 |
| Hawaii DOH | Not available | - |
| Idaho DHW | http://www.healthandwelfare.idaho.gov/Health/Labs/CertificationDrinkingWaterLabs/tabid/1833/Default.aspx | - |
| ISO 17025 | http://www.pjllabs.com/ | L14-50 |
| Louisiana DEQ | http://www.deq.louisiana.gov/portal/DIVISIONS/PublicParticipationandPermitSupport/LouisianaLaboratoryAccreditationProgram.aspx | 03016 |
| Maine DHS | Not available | WA01276 |
| Michigan DEQ | http://www.michigan.gov/deq/0,1607,7-135-3307_4131_4156---,00.html | 9949 |
| Minnesota DOH | http://www.health.state.mn.us/accreditation | 053-999-457 |
| Montana DPHHS | http://www.dphhs.mt.gov/publichealth/ | CERT0047 |
| Nevada DEP | http://ndep.nv.gov/bsdwlabservice.htm | WA01276 |
| New Jersey DEP | http://www.nj.gov/dep/oqa/ | WA005 |
| North Carolina DWQ | http://www.dwqlab.org/ | 605 |
| Oklahoma DEQ | http://www.deq.state.ok.us/CSDnew/labcert.htm | 9801 |
| Oregon – DEQ (NELAP) | http://public.health.oregon.gov/LaboratoryServices/EnvironmentalLaboratoryAccreditation/Pages/index.aspx | WA100010 |
| South Carolina DHEC | http://www.scdhec.gov/environment/envserv/ | 61002 |
| Texas CEQ | http://www.tceq.texas.gov/field/qa/env_lab_accreditation.html | T104704427 |
| Washington DOE | http://www.ecy.wa.gov/programs/eap/labs/lab-accreditation.html | C544 |
| Wisconsin DNR | http://dnr.wi.gov/ | 998386840 |
| Wyoming (EPA Region 8) | http://www.epa.gov/region8/water/dwhome/wyomingdi.html | - |
| Kelso Laboratory Website | www.alsglobal.com | NA |

Analyses were performed according to our laboratory's NELAP-approved quality assurance program. A complete listing of specific NELAP-certified analytes, can be found in the certification section at www.ALSGlobal.com or at the accreditation bodies web site.

Please refer to the certification and/or accreditation body's web site if samples are submitted for compliance purposes. The states highlighted above, require the analysis be listed on the state certification if used for compliance purposes and if the method/analyte is offered by that state.



Case Narrative


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ALS Environmental—Kelso Laboratory
1317 South 13th Avenue, Kelso, WA 98626
Phone (360)577-7222 Fax (360)636-1068
www.alsglobal.com

ALS ENVIRONMENTAL

| | | | |
|-----------------------|--|-----------------------------|----------|
| Client: | Eurofins Calscience Environmental Laboratory | Service Request No.: | K1512268 |
| Project: | 15-10-1995 | Date Received: | 10/29/15 |
| Sample Matrix: | Water | | |

Case Narrative

All analyses were performed consistent with the quality assurance program of ALS Environmental. This report contains analytical results for samples designated for Tier II data deliverables. When appropriate to the method, method blank results have been reported with each analytical test. Surrogate recoveries have been reported for all applicable organic analyses. Additional quality control analyses reported herein include: Laboratory Control Sample (LCS), and Laboratory/Duplicate Laboratory Control Sample (LCS/DLCS).

Sample Receipt

Thirty-two water samples were received for analysis at ALS Environmental on 10/29/15. The samples were received in good condition and consistent with the accompanying chain of custody form. The samples were stored in a refrigerator at 4°C upon receipt at the laboratory.

Steroids and Endocrine Disrupting Compounds by Method 1694**Elevated Detection Limits:**

Samples HCS 210 Trail, HCS 270 Trail, HSM 230 Trail, FDHSM 230 Trail, HSM 230 Lead and HSM 240 Lead required dilution due to the presence of elevated levels of target analyte. The reporting limits were adjusted to reflect the dilution.

No other anomalies associated with the analysis of these samples were observed.

Approved by _____



Chain of Custody


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Phone (360)577-7222 Fax (360)636-1068
www.alsglobal.com

To: ALS - Kelso

DATE: 10/28/15

PAGE: 1 OF 3

7440 Lincoln Way, Garden Grove, CA 92841-1427 • (714) 895-5494

For courier service / sample drop off information, contact us26_sales@eurofinsus.com or call us.

[illegible]

PAGE: 2 OF 3

[illegible]



Cooler Receipt and Preservation Form

Client / Project: Eurofin Cal Science Service Request K15 12268
Received: 10/29/15 Opened: 10/29/15 By: 16 Unloaded: 10/29/15 By: 16

1. Samples were received via? Mail Fed Ex UPS DHL PDX Courier Hand Delivered
2. Samples were received in: (circle) Cooler Box Envelope Other NA
3. Were custody seals on coolers? NA Y N If yes, how many and where? 1, front
If present, were custody seals intact? Y N If present, were they signed and dated? Y N

| Raw Cooler Temp | Corrected Cooler Temp | Raw Temp Blank | Corrected Temp Blank | Corr. Factor | Thermometer ID | Cooler/COC ID | Tracking Number | NA | Filed |
|-----------------|-----------------------|----------------|----------------------|--------------|----------------|---------------|-----------------|----|-------|
| -6 | -3 | - | - | +3 | 323 | NA | 77485048 1436 | | |
| -7 | -5 | - | - | +2 | 371 | | 77485048 1138 | | |
| -4 | -2 | - | - | +2 | 368 | | 77485048 1697 | | |
| | | | | | | | | | |
| | | | | | | | | | |

4. Packing material: Inserts Baggies Bubble Wrap Gel Packs Wet Ice Dry Ice Sleeves
5. Were custody papers properly filled out (ink, signed, etc.)? NA Y N
6. Did all bottles arrive in good condition (unbroken)? *Indicate in the table below.* NA Y N
7. Were all sample labels complete (i.e analysis, preservation, etc.)? NA Y N
8. Did all sample labels and tags agree with custody papers? *Indicate major discrepancies in the table on page 2.* NA Y N
9. Were appropriate bottles/containers and volumes received for the tests indicated? NA Y N
10. Were the pH-preserved bottles (*see SMO GEN SOP*) received at the appropriate pH? *Indicate in the table below* NA Y N
11. Were VOA vials received without headspace? *Indicate in the table below.* NA Y N
12. Was C12/Res negative? NA Y N

| Sample ID on Bottle | Sample ID on COC | Identified by: |
|---------------------|------------------|----------------|
| | | |
| | | |
| | | |

| Sample ID | Bottle Count Bottle Type | Out of Temp | Head- space | Broke | pH | Reagent | Volume added | Reagent Lot Number | Initials | Time |
|-----------|-----------------------------|----------------|----------------|-------|----|---------|-----------------|-----------------------|----------|------|
| | | | | | | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |

Notes, Discrepancies, & Resolutions: _____



Steroids and Endocrine Disrupting Compounds


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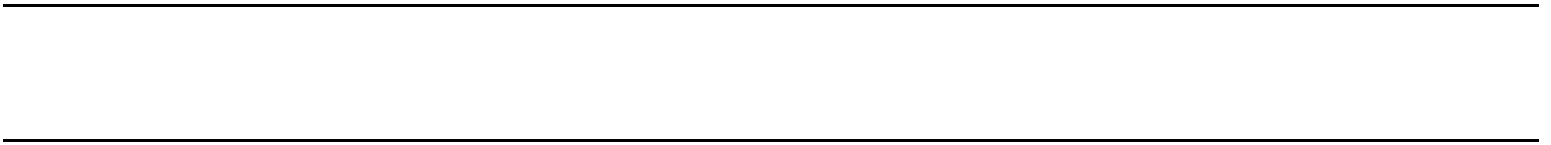





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AMPLIFIED
GEOCHEMICAL
IMAGING, LLC

Laboratory Report

Site: Comal & San Marcos Rivers
February 2015 Sampling

Prepared for:

SWCA Environmental
10245 Little York Road Suite 600
Houston, Texas
UNITED STATES

Prepared on:
March 04, 2015

Project Summary and Objective

Amplified Geochemical Imaging, LLC. (AGI) provided the AGI Environmental Survey used at:

Comal & San Marcos Rivers

February 2015 Sampling

The service provided by AGI included delivery of the required quantity of AGI Universal Samplers, analysis by the method described below for the requested organic compounds, reporting of the data, and contour mapping (as needed).

This report includes results for only the samples noted under the Laboratory Sample Report section. If contour maps are part of the project deliverable, the maps will be prepared and issued under a separate report cover, upon receipt of a usable sitemap (electronic) and compound choices for contouring.

Written/submitted by:

Jim E Whetzel

Project Manager

Reviewed/approved by:

Don D'Apolito

Project Manager

Analytical data approved by:

Fatima Niazi

Chemist

Quality Assurance Statement

The AGI Laboratory, at Amplified Geochemical Imaging's facility in Elkton, MD USA, operates under the guidelines of its ISO Standard 17025 DoD ELAP accreditation, and its Quality Assurance Manual, Operating Procedures, and Methods (SPG-SOP-0462).

For this project, the analytical method, results, and observations reported do [] do not [✓] fall within the scope of AGI's ISO 17025 accreditation.

Screening/Concentration Method

The AGI Universal Samplers are analyzed at AGI's fixed laboratory using thermal desorption-gas chromatography/mass spectrometry (TD-GC/MS) instrumentation following U.S. EPA Method 8260 (SPG-WI-0292) which includes the following:

- **BFB Tuning Frequency:** A BFB tune is analyzed at the start of each analytical run and after every 30 samples.
- **Initial Calibration:** A minimum of a five point calibration curve is analyzed prior to the analysis of samples.
- **Linearity of Target Compounds:** If the RSD of any target analyte is less than or equal to 25% then average response factor can be used for quantitation. If the RSD exceeds 25% for a target compound a regression equation can be used for quantitation.
- **Continuing Calibration Verification:** After every 10 samples, and at the end of each analytical batch, and a second-source Reference Standard is analyzed near the mid point of the calibration curve. The acceptance criteria for all target analytes in the reference standards are +/- 50% of the true value.
- **Method Blank:** Analyzed prior to the analysis of field samples and every 30 samples.

Note: Analyte levels reported for the field-deployed AGI Universal Samplers that exceed trip and method blank levels, and/or the reporting limit, are more likely to have originated from on-site sources.

| | |
|----------------------------|-------------------|
| Media Sampled: | WATER |
| Chemist - sample analysis: | Kelly J Stringham |
| Chemist - data processor: | Kelly J Stringham |
| Chemist - data review: | Fatima Niazi |

Method deviations: No data is available for the initial method blank. Data was inadvertently overwritten. There was no evidence of system contamination observed in any other QC sample.

Please note that data file names ending with R are rerun samples using the second pair of sorbers, in which the original results were not reported. Data file names ending in D are duplicate analysis results for the second set of sorbers from the same sampler, and are reported.

Additional Report Information

- Comments
- Laboratory Sample Report
- Chain of Custody
- Installation and Retrieval Log
- Analytical Results and Key
- Concentration Calculation Method Summary
- Total Ion Chromatograms

Project Specific Comments

None.

Survey period ¹

Samplers were installed on February 3, 2015 and retrieved on February 17, 2015 for an exposure period of 14 days.

Tamper seal intact:

Yes

Date received:

2/18/2015 1:00 PM

By: Lisa Bozette

COC returned:

Yes

Comments:

None

¹ - Installation start to end of retrieval, as reported. See installation and retrieval log for individual deployment and retrieval dates and times (i.e., sampler exposure time).

General Comments

Analytical QA/QC

Laboratory instrumentation consists of gas chromatographs equipped with mass selective detectors, coupled with automated thermal desorption units. Sample preparation involves cutting the tip off the bottom of the AGI Universal Sampler, and transferring one or more "sorbents" to a thermal desorption tube for analysis. The insertion/retrieval cord prevents soil, water and other interferences from coming in contact with the adsorbent. No further sample preparation is required. Any replicate sorbents not consumed in the initial analysis will be discarded fifteen (15) days from the date of the laboratory report.

Data are archived and stored in a secure manner as per AGI's Quality Assurance program (SPG-SOP-0462).

Total petroleum hydrocarbons (TPH), gasoline-range petroleum hydrocarbons (GRPH), and/or diesel range petroleum hydrocarbons (DRPH), when reported, are calculated using the area under the peaks observed in m/z 55 and 57 selected ion chromatograms. Quantitation of the mass values was performed using the response factor for a specific alkane (present in the calibration standards). TPH values include the entire chromatogram and provide estimates for aliphatic hydrocarbon ranges of C4 to C20. GRPH and DRPH include only the relevant regions of the chromatograms and provide estimates for C4 to C10 and C10 to C20 aliphatic hydrocarbons, respectively.

Trip blanks were provided to document potential exposures that were not part of the signal of interest (e.g., impact during sampler shipment, installation and/or retrieval, and storage). The trip blanks are identically manufactured and packaged AGI Universal Samplers to those samplers deployed in the field. The trip blanks remain unopened during all phases of the project. Levels reported on the trip blanks may indicate potential impact to the samplers other than the contaminant source of interest.

Unresolved peak envelopes (UPEs) are represented as a series of compound peaks clustered together around a central gas chromatograph elution time in the total ion chromatogram. UPEs may be indicative of complex fluid mixtures. UPEs observed early in the chromatograms are considered to indicate presence of more volatile fluids, while UPEs observed later in the chromatogram may indicate the presence of less volatile fluids. Multiple UPEs may indicate the presence of multiple complex fluids.

Total ion chromatograms (TICs) are included in the Attachments. The eight-digit serial number of each sampler is incorporated in the TIC identification (e.g., 12345678.D represents AGI Universal Sampler 12345678).

General Comments

Soil Gas Sampling

For soil gas sampling, the AGI Environmental Survey reports mass levels migrating through the open pore spaces of the soil and diffusing through the sampler membrane for sorption by the engineered, hydrophobic adsorbents, housed within the membrane tube. During the migration of the soil gas away from the source to the AGI Universal Sampler, the vapors are subject to a variety of attenuation factors. The soil gas masses reported on the samplers compare favorably with the concentrations reported in the soil or groundwater (e.g., where soil gas levels are reported at greater levels to other sampled locations on the site, the matrix data should reveal the same pattern, and vice versa). However, due to a variety of factors, a perfect comparison between matrix data and soil gas levels can rarely be achieved.

Soil gas concentrations ($\mu\text{g}/\text{m}^3$) are calculated following the method described in the Additional Report Information section.

Soil gas signals reported by this method cannot be correlated specifically to soil adsorbed, groundwater, and /or free-phase contamination. The soil gas signal reported from each AGI Universal Sampler can evolve from all of these sources. Differentiation between soil and groundwater contamination can only be achieved with prior knowledge of the site history (i.e., the site is known to have groundwater contamination only).

Air Sampling

For indoor, outdoor, and crawlspace air sampling, the AGI Environmental Survey reports mass levels present in the air and diffusing through the sampler membrane for sorption by the engineered adsorbents housed within the membrane tube.

Air concentrations ($\mu\text{g}/\text{m}^3$) are calculated following the method described in the Additional Report Information section.

Groundwater and Sediment Porewater Sampling

For groundwater and sediment porewater sampling, the AGI Environmental Survey reports the mass levels of compounds present in the water which, when coming in contact with the sampler membrane, partitions out of solution, and diffuses through the sampler membrane for sorption by the engineered adsorbents.

Water concentrations ($\mu\text{g}/\text{L}$) are calculated using the quantified mass, exposure period and the compound specific uptake rate. The rates were measured under controlled experimental conditions. The uptake rates are corrected for water pressure (depth of the AGI Universal Sampler below the water table), water temperature and the aquifer flow rate. For sediment porewater, the uptake rate is corrected for the reduced volume of water in the sediment, by multiplying the uptake rate by the pore water fraction.

LABORATORY SAMPLE REPORT

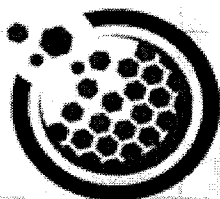
Project: ENV 01315-2

Site Name: Comal & San Marcos Rivers

Module Type: SPG0008

| Module ID | Sample Type | Field ID | |
|----------------------------|--------------------------|---------------------|-------------------|
| 00755224 | FIELD_SAMPLE | HCS410 | |
| 00755225 | FIELD_SAMPLE | HCS420 | |
| 00755226 | FIELD_SAMPLE | HCS430 | |
| 00755227 | FIELD_SAMPLE | HCS440 | |
| 00755228 | FIELD_SAMPLE | FDHCS440 | |
| 00755229 | FIELD_SAMPLE | HCS460 | |
| 00755230 | FIELD_SAMPLE | HSM410 | |
| 00755231 | FIELD_SAMPLE | HSM420 | |
| 00755232 | FIELD_SAMPLE | HSM430 | |
| 00755233 | FIELD_SAMPLE | FDHSM430 | |
| 00755234 | FIELD_SAMPLE | HSM440 | |
| 00755235 | FIELD_SAMPLE | HSM450 | |
| 00755236 | FIELD_SAMPLE | HSM460 | |
| 00755237 | FIELD_SAMPLE | HSM470 | |
| 00755238 | TRIP_BLANK | TB02 | |
| Total # "FIELD SAMPLES" | Total # "TRIP BLANKS" | Total # "UNUSED" | Total # "LOST" |
| 14 | 1 | 0 | 0 |

Duplicate samples: 0



AMPLIFIED GEOCHEMICAL IMAGING, LLC

210 Executive Drive, Suite 1
Newark, DE 19702-3335 USA
ph: +1-302-266-2428
www.agisurveys.net

AGI Universal Passive Sampler Chain of Groundwater Sampling

Production Order #: 01315

Customer Name: SWCA Environmental
Address: 10245 Little York Road Suite 600

Site Name: Comal & San Marcos Rivers
Site Address: 2015 sampling

Houston, Texas 77040
USA

Project Manager:

Serial # of Samplers Shipped
00755224 - 00755241

of Samplers for Installation 17.00
Total Samplers Shipped 18.00
Total Samplers Received 13
Total Samplers Installed 13

of Trip Blanks 1
Pieces
Pieces
Pieces

Returned
00755224 - 00755238

Serial # of Trip Blanks (Client Decides)

00755238

Prepared By:

Lisa Borzette

Verified By:

Jennifer Moreland

Is Concurrent water sampling
planned this monitoring period?

YES

NO

Scheduled Sampling Date:

Installation Performed By:

Name: Jennifer Moreland Brittany Rio
Company: SWCA

Retrieval Performed By:

Name: Jennifer Moreland Brittany Rio
Company: SWCA

Installation Start Date / Time: 2/13/15 9:11

Retrieval Start Date / Time: 2/17/15 10:22

Installation Complete Date / Time: 2/13/15 12:45

Retrieval Complete Date / Time: 2/17/15 12:24

Total Samplers Retrieved:

13

Total Samplers Lost In Field:

0

Total Unused Samplers Returned:

3 extra kept for next install

Relinquished By:

Lisa Borzette

Date/Time

01/26/15
11:00

Company:

agi

Received By:

Jennifer Moreland

Date/Time

01/28/15
13:00

Company:

SWCA

Relinquished By:

Jennifer Moreland

Date/Time

2/17/15
14:45

Company:

SWCA

Received By:

Lisa Borzette

Date/Time

2-18-15
1:00 pm

Company:

agi



AMPLIFIED
GEOCHEMICAL
IMAGING, LLC

210 Executive Drive, Suite 1
Newark, DE USA 19702-3335
ph: 302-266-2428

AGI Project No. ENV 01077
Site Name: Comal & San Marco River
Site Location: New Braunfels and San Marcos, Texas

AGI Soil Gas Sampling
Installation & Retrieval Log

Company Name: SWCA Environmental
Location: Comal and San Marcos Rivers
Samples collected by: Philip Pearce and Jennifer Moreland

* Optional or as needed

| | | | | | WATER QUALITY MONITORING | | | | | OBSERVATIONS/COMMENTS* (e.g., observations, location description, missing, pulled from well, etc. - as needed) | YES / NO | |
|--------------------|-----------|---|---|---|---|---|--|--|---|--|--|--------|
| SAMPLER SERIAL NO. | SAMPLE ID | SAMPLE TYPE (Field Sample, Trip Blank, Field Blank, etc.) | INSTALLATION DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 13:00 | RETRIEVAL DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 15:00 | INSTALLATION DEPTH FROM TOP OF CASING (feet) | DEPTH TO WATER FROM TOP OF CASING (feet) | SCREENED INTERVAL FROM TOP OF CASING (feet) | WATER TEMPERATURE AT MODULE DEPTH (deg C) | FLOW RATE THROUGH SCREEN: HIGH OR LOW (HIGH = flow > than 10 m/day; LOW = flow < 10 m/day) | | EVIDENCE OF LIQUID PETROLEUM HYDROCARBONS? | ODOR ? |
| 00755224 | HCS410 | FIELD_SAMPLE | 2/3/2015 10:02 | 2/17/2015 11:56 | 1.5 | | | 22.4 | High | None | No | No |
| 00755225 | HCS420 | FIELD_SAMPLE | 2/3/2015 10:13 | 2/17/2015 12:01 | 3 | | | 22.4 | High | None | No | No |
| 00755226 | HCS430 | FIELD_SAMPLE | 2/3/2015 9:11 | 2/17/2015 11:25 | 7 | | | 22.4 | High | None | No | No |
| 00755227 | HCS440 | FIELD_SAMPLE | 2/3/2015 10:27 | 2/17/2015 12:12 | 4 | | | 22.4 | High | None | No | No |
| 00755228 | FDHCS440 | FIELD_DUPLICATE | 2/3/2015 10:27 | 2/17/2015 12:12 | 4 | | | 22.4 | High | None | No | No |
| 00755229 | HCS460 | FIELD_SAMPLE | 2/3/2015 10:38 | 2/17/2015 12:24 | 4 | | | 22.4 | High | None | No | No |
| 00755230 | HSM410 | FIELD_SAMPLE | 2/3/2015 11:32 | 2/17/2015 9:45 | 2 | | | 21.2 | High | None | No | No |
| 00755231 | HSM420 | FIELD_SAMPLE | 2/3/2015 11:45 | 2/17/2015 9:55 | 4 | | | 21.2 | High | None | No | No |
| 00755232 | HSM430 | FIELD_SAMPLE | 2/3/2015 11:55 | 2/17/2015 10:01 | 0.5 | | | 21.2 | High | None | No | No |
| 00755233 | FDHSM430 | FIELD_DUPLICATE | 2/3/2015 11:55 | 2/17/2015 10:01 | 0.5 | | | 21.2 | High | None | No | No |
| 00755234 | HSM440 | FIELD_SAMPLE | 2/3/2015 12:06 | 2/17/2015 10:08 | 5 | | | 21.2 | High | None | No | No |
| 00755235 | HSM450 | FIELD_SAMPLE | 2/3/2015 12:22 | 2/17/2015 10:22 | 4 | | | 21.2 | High | None | No | No |
| 00755236 | HSM460 | FIELD_SAMPLE | 2/3/2015 12:33 | 2/17/2015 10:35 | 3 | | | 21.2 | High | None | No | No |
| 00755237 | HSM470 | FIELD_SAMPLE | 2/3/2015 12:45 | 2/17/2015 10:47 | 1.5 | | | 21.2 | High | None | No | No |
| 00755238 | TB02 | TRIP_BLANK | 2/3/2015 9:11 | 2/17/2015 12:24 | NA | | | NA | NA | None | NA | NA |
| | | | | | | | | | | | | |



PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755224 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS410

Installation Date: 2/3/2015 10:02:00AM

Retrieval Date: 2/17/2015 11:56:00AM

Date Analyzed: 2/25/2015 3:12:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.05 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755224 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS410

Installation Date: 2/3/2015 10:02:00AM

Retrieval Date: 2/17/2015 11:56:00AM

Date Analyzed: 2/25/2015 3:12:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.76 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755225 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS420

Installation Date: 2/3/2015 10:13:00AM

Retrieval Date: 2/17/2015 12:01:00PM

Date Analyzed: 2/25/2015 7:50:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.28 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755225 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS420

Installation Date: 2/3/2015 10:13:00AM

Retrieval Date: 2/17/2015 12:01:00PM

Date Analyzed: 2/25/2015 7:50:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.52 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755226 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS430

Installation Date: 2/3/2015 9:11:00AM

Retrieval Date: 2/17/2015 11:25:00AM

Date Analyzed: 2/25/2015 2:40:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.29 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755226 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS430

Installation Date: 2/3/2015 9:11:00AM

Retrieval Date: 2/17/2015 11:25:00AM

Date Analyzed: 2/25/2015 2:40:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 1.42 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755227 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS440

Installation Date: 2/3/2015 10:27:00AM

Retrieval Date: 2/17/2015 12:12:00PM

Date Analyzed: 2/25/2015 7:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.28 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755227 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS440

Installation Date: 2/3/2015 10:27:00AM

Retrieval Date: 2/17/2015 12:12:00PM

Date Analyzed: 2/25/2015 7:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.61 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755228 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: FDHCS440

Installation Date: 2/3/2015 10:27:00AM

Retrieval Date: 2/17/2015 12:12:00PM

Date Analyzed: 2/25/2015 9:23:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.28 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755228 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: FDHCS440

Installation Date: 2/3/2015 10:27:00AM

Retrieval Date: 2/17/2015 12:12:00PM

Date Analyzed: 2/25/2015 9:23:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.76 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755229 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS460

Installation Date: 2/3/2015 10:38:00AM

Retrieval Date: 2/17/2015 12:24:00PM

Date Analyzed: 2/25/2015 6:49:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.18 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755229 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS460

Installation Date: 2/3/2015 10:38:00AM

Retrieval Date: 2/17/2015 12:24:00PM

Date Analyzed: 2/25/2015 6:49:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755230 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM410

Installation Date: 2/3/2015 11:32:00AM

Retrieval Date: 2/17/2015 9:45:00AM

Date Analyzed: 2/25/2015 4:45:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755230 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM410

Installation Date: 2/3/2015 11:32:00AM

Retrieval Date: 2/17/2015 9:45:00AM

Date Analyzed: 2/25/2015 4:45:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.90 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755231 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM420

Installation Date: 2/3/2015 11:45:00AM

Retrieval Date: 2/17/2015 9:55:00AM

Date Analyzed: 2/25/2015 8:52:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.08 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755231 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM420

Installation Date: 2/3/2015 11:45:00AM

Retrieval Date: 2/17/2015 9:55:00AM

Date Analyzed: 2/25/2015 8:52:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755232 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM430

Installation Date: 2/3/2015 11:55:00AM

Retrieval Date: 2/17/2015 10:01:00AM

Date Analyzed: 2/25/2015 9:54:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.13 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755232 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM430

Installation Date: 2/3/2015 11:55:00AM

Retrieval Date: 2/17/2015 10:01:00AM

Date Analyzed: 2/25/2015 9:54:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 1.27 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755233 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: FDHSM430

Installation Date: 2/3/2015 11:55:00AM

Retrieval Date: 2/17/2015 10:01:00AM

Date Analyzed: 2/25/2015 4:14:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.16 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755233 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: FDHSM430

Installation Date: 2/3/2015 11:55:00AM

Retrieval Date: 2/17/2015 10:01:00AM

Date Analyzed: 2/25/2015 4:14:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.99 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755234 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM440

Installation Date: 2/3/2015 12:06:00PM

Retrieval Date: 2/17/2015 10:08:00AM

Date Analyzed: 2/25/2015 10:56:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.13 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



AMPLIFIED GEOCHEMICAL IMAGING, LLC

210 Executive Drive, Suite 1
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ph: +1-302-266-2428
www.agisurveys.net

PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755234 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM440

Installation Date: 2/3/2015 12:06:00PM

Retrieval Date: 2/17/2015 10:08:00AM

Date Analyzed: 2/25/2015 10:56:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755235 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM450

Installation Date: 2/3/2015 12:22:00PM

Retrieval Date: 2/17/2015 10:22:00AM

Date Analyzed: 2/25/2015 10:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.05 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



AMPLIFIED GEOCHEMICAL IMAGING, LLC

210 Executive Drive, Suite 1
Newark, DE 19702-3335 USA
ph: +1-302-266-2428
www.agisurveys.net

PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755235 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM450

Installation Date: 2/3/2015 12:22:00PM

Retrieval Date: 2/17/2015 10:22:00AM

Date Analyzed: 2/25/2015 10:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755236 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM460

Installation Date: 2/3/2015 12:33:00PM

Retrieval Date: 2/17/2015 10:35:00AM

Date Analyzed: 2/25/2015 3:43:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.09 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755236 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM460

Installation Date: 2/3/2015 12:33:00PM

Retrieval Date: 2/17/2015 10:35:00AM

Date Analyzed: 2/25/2015 3:43:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755237 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM470

Installation Date: 2/3/2015 12:45:00PM

Retrieval Date: 2/17/2015 10:47:00AM

Date Analyzed: 2/25/2015 8:21:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.10 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



AMPLIFIED GEOCHEMICAL IMAGING, LLC

210 Executive Drive, Suite 1
Newark, DE 19702-3335 USA
ph: +1-302-266-2428
www.agisurveys.net

PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755237 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM470

Installation Date: 2/3/2015 12:45:00PM

Retrieval Date: 2/17/2015 10:47:00AM

Date Analyzed: 2/25/2015 8:21:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.54 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: February 2015 Sampling

FOR: SWCA Environmental
Houston, Texas 77040
USA

SAMPLER ID: 00755238 TRIP_BLANK

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: TB02

Installation Date: 2/3/2015 9:11:00AM

Retrieval Date: 2/17/2015 12:24:00PM

Date Analyzed: 2/25/2015 11:28:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01315-2

FOR: SWCA Environmental

SITE NAME: Comal & San Marcos Rivers

SITE ADDRESS: February 2015 Sampling

Houston, Texas 77040

USA

SAMPLER ID: 00755238 TRIP_BLANK

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: TB02

Installation Date: 2/3/2015 9:11:00AM

Retrieval Date: 2/17/2015 12:24:00PM

Date Analyzed: 2/25/2015 11:28:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150223-2

Reviewer: Fatima Niazi

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |

AMPLIFIED GEOCHEMICAL IMAGING, LLC
210 EXECUTIVE DRIVE, SUITE 1, NEWARK, DE 19702
SWCA ENVIRONMENTAL, HOUSTON, TX
STANDARD TARGET VOCs/SVOCs
ESTIMATED WATER CONCENTRATIONS
COMAL & SAN MARCOS RIVERS
ORDER # 01315-2

| | | | | | | | | | estimated | | | |
|----------|----------|------------|------------|------------|----|------------|----|----|-----------|------------|--------------|-------------|
| DATAFILE | FIELD | DATE/ TIME | DATE/ TIME | DATE/ TIME | | DATE/ TIME | | | | | | |
| NAME | ID | INSTALLED | RETRIEVED | RECEIVED | | ANALYZED | | DF | TPH, ug/L | MTBE, ug/L | t12DCE, ug/L | 11DCA, ug/L |
| RL = | | | | | | | | | 0.06 | 0.01 | 0.01 | 0.01 |
| 00755224 | HCS410 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | 0.06 | <0.01 | <0.01 | <0.01 |
| 00755225 | HCS420 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | 0.06 | <0.01 | <0.01 | <0.01 |
| 00755226 | HCS430 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | 0.08 | <0.01 | <0.01 | <0.01 |
| 00755227 | HCS440 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | 0.06 | <0.01 | <0.01 | <0.01 |
| 00755228 | FDHCS440 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | 0.06 | <0.01 | <0.01 | <0.01 |
| 00755229 | HCS460 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | <0.05 | <0.01 | <0.01 | <0.01 |
| 00755230 | HSM410 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | 0.07 | <0.01 | <0.01 | <0.01 |
| 00755231 | HSM420 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | <0.06 | <0.01 | <0.01 | <0.01 |
| 00755232 | HSM430 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | 0.08 | <0.01 | <0.01 | <0.01 |
| 00755233 | FDHSM430 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | 0.07 | <0.01 | <0.01 | <0.01 |
| 00755234 | HSM440 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | <0.06 | <0.01 | <0.01 | <0.01 |
| 00755235 | HSM450 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | <0.06 | <0.01 | <0.01 | <0.01 |
| 00755236 | HSM460 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | <0.06 | <0.01 | <0.01 | <0.01 |
| 00755237 | HSM470 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | 0.06 | <0.01 | <0.01 | <0.01 |
| | | | | | | | | | | | | |
| 00755238 | TB02 | 2/3/2015 | 2/17/2015 | 2/18/2015 | ET | 2/25/2015 | ET | 1 | <0.05 | <0.01 | <0.01 | <0.01 |
| | | | | | | | | | | | | |

AMPLIFIED GEOCHEMICAL IMAGING, LLC
210 EXECUTIVE DRIVE, SUITE 1, NEWARK, DE 19702
SWCA ENVIRONMENTAL, HOUSTON, TX
STANDARD TARGET VOCs/SVOCs
ESTIMATED WATER CONCENTRATIONS
COMAL & SAN MARCOS RIVERS
ORDER # 01315-2

| DATAFILE | FIELD | | | | | | | | | |
|----------|----------|--------------|-------------|--------------|-------------|------------|------------|-----------|--------------|-----------|
| NAME | ID | c12DCE, ug/L | CHCl3, ug/L | 111TCA, ug/L | 12DCA, ug/L | BENZ, ug/L | CCl4, ug/L | TCE, ug/L | 112TCA, ug/L | TOL, ug/L |
| RL = | | 0.01 | 0.01 | 0.01 | 0.01 | 0.01 | 0.00 | 0.01 | 0.01 | 0.01 |
| 00755224 | HCS410 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755225 | HCS420 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755226 | HCS430 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755227 | HCS440 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755228 | FDHCS440 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755229 | HCS460 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755230 | HSM410 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755231 | HSM420 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755232 | HSM430 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755233 | FDHSM430 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755234 | HSM440 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755235 | HSM450 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755236 | HSM460 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755237 | HSM470 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |
| 00755238 | TB02 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 |

AMPLIFIED GEOCHEMICAL IMAGING, LLC
210 EXECUTIVE DRIVE, SUITE 1, NEWARK, DE 19702
SWCA ENVIRONMENTAL, HOUSTON, TX
STANDARD TARGET VOCs/SVOCs
ESTIMATED WATER CONCENTRATIONS
COMAL & SAN MARCOS RIVERS
ORDER # 01315-2

| DATAFILE | FIELD | | | | | | | | |
|----------|----------|-----------|-----------|--------------|-----------------|--------------|-------------|------------|-----------------|
| NAME | ID | OCT, ug/L | PCE, ug/L | CIBENZ, ug/L | 1112TetCA, ug/L | ETBENZ, ug/L | mpXYL, ug/L | oXYL, ug/L | 1122TetCA, ug/L |
| RL = | | 0.00 | 0.01 | 0.01 | 0.01 | 0.01 | 0.01 | 0.01 | 0.01 |
| 00755224 | HCS410 | <0.00 | 0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755225 | HCS420 | <0.00 | 0.05 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755226 | HCS430 | <0.00 | 0.06 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755227 | HCS440 | <0.00 | 0.05 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755228 | FDHCS440 | <0.00 | 0.06 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755229 | HCS460 | <0.00 | 0.04 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755230 | HSM410 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755231 | HSM420 | <0.00 | 0.02 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755232 | HSM430 | <0.00 | 0.03 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755233 | FDHSM430 | <0.00 | 0.04 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755234 | HSM440 | <0.00 | 0.03 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755235 | HSM450 | <0.00 | 0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755236 | HSM460 | <0.00 | 0.02 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755237 | HSM470 | <0.00 | 0.02 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00755238 | TB02 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |

AMPLIFIED GEOCHEMICAL IMAGING, LLC
210 EXECUTIVE DRIVE, SUITE 1, NEWARK, DE 19702
SWCA ENVIRONMENTAL, HOUSTON, TX
STANDARD TARGET VOCs/SVOCs
ESTIMATED WATER CONCENTRATIONS
COMAL & SAN MARCOS RIVERS
ORDER # 01315-2

| | | | | | | | | | estimated |
|----------|----------|--------------|--------------|-------------|-------------|-------------|-------------|------------|--------------|
| DATAFILE | FIELD | | | | | | | | |
| NAME | ID | 135TMB, ug/L | 124TMB, ug/L | 13DCB, ug/L | 14DCB, ug/L | 12DCB, ug/L | UNDEC, ug/L | NAPH, ug/L | TRIDEC, ug/L |
| RL = | | 0.00 | 0.01 | 0.01 | 0.01 | 0.01 | 0.02 | 0.02 | 0.02 |
| 00755224 | HCS410 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755225 | HCS420 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755226 | HCS430 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755227 | HCS440 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755228 | FDHCS440 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755229 | HCS460 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755230 | HSM410 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755231 | HSM420 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755232 | HSM430 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755233 | FDHSM430 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755234 | HSM440 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755235 | HSM450 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755236 | HSM460 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00755237 | HSM470 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| | | | | | | | | | |
| 00755238 | TB02 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| | | | | | | | | | |

AMPLIFIED GEOCHEMICAL IMAGING, LLC
210 EXECUTIVE DRIVE, SUITE 1, NEWARK, DE 19702
SWCA ENVIRONMENTAL, HOUSTON, TX
STANDARD TARGET VOCs/SVOCs
ESTIMATED WATER CONCENTRATIONS
COMAL & SAN MARCOS RIVERS
ORDER # 01315-2

| | | | estimated | estimated | estimated | estimated |
|----------|----------|---------------|----------------------|----------------|--------------------|----------------|
| DATAFILE | FIELD | | | | | |
| NAME | ID | 2MeNAPH, ug/L | Acenaphthylene, ug/L | PENTADEC, ug/L | Acenaphthene, ug/L | Fluorene, ug/L |
| RL = | | 0.01 | 0.01 | 0.02 | 0.01 | 0.01 |
| 00755224 | HCS410 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00755225 | HCS420 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00755226 | HCS430 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00755227 | HCS440 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00755228 | FDHCS440 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00755229 | HCS460 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00755230 | HSM410 | <0.01 | <0.02 | <0.02 | <0.02 | <0.02 |
| 00755231 | HSM420 | <0.01 | <0.02 | <0.02 | <0.02 | <0.02 |
| 00755232 | HSM430 | <0.01 | <0.02 | <0.02 | <0.02 | <0.02 |
| 00755233 | FDHSM430 | <0.01 | <0.02 | <0.02 | <0.02 | <0.02 |
| 00755234 | HSM440 | <0.01 | <0.02 | <0.02 | <0.02 | <0.02 |
| 00755235 | HSM450 | <0.01 | <0.02 | <0.02 | <0.02 | <0.02 |
| 00755236 | HSM460 | <0.01 | <0.02 | <0.02 | <0.02 | <0.02 |
| 00755237 | HSM470 | <0.01 | <0.02 | <0.02 | <0.02 | <0.02 |
| | | | | | | |
| 00755238 | TB02 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| | | | | | | |

KEY TO DATA TABLE

UNITS

| | |
|-------------------|--|
| µg | micrograms, relative mass value |
| µg/m ³ | micrograms per cubic meter; estimated soil gas concentration |
| µg/L | micrograms per Liter; calculated water concentration |

DATA QUALIFIERS

| | |
|---|--|
| > | greater than; value exceeds calibration range, estimated value |
| < | less than; compound value is below the LOD and RL |
| J | mass value below LOQ or RL, but above LOD, estimated mass value |
| E | mass value exceeds upper calibration level, estimated mass value |
| Q | one or more quality control parameters failed for the compound |

ABBREVIATIONS

| | |
|--------|--|
| AVG RL | average reporting limit; calculated based on individual field sample RLs |
| LOD | limit of detection |
| LOQ | limit of quantification |
| MDL | method detection limit |
| RL | reporting limit |

| | | | |
|-------------|---|----------|--|
| 1112TetCA | 1,1,1,2-tetrachloroethane | CIBENZ | chlorobenzene |
| 111TCA | 1,1,1-trichloroethane | ct12DCE | cis- & trans-1,2-dichloroethene |
| 1122TetCA | 1,1,2,2-tetrachloroethane | EtBENZ | ethylbenzene |
| 112TCA | 1,1,2-trichloroethane | mpXYL | m-, p-xylene |
| 11DCA | 1,1-dichloroethane | MTBE | methyl t-butyl ether |
| 11DCE | 1,1-dichloroethene | NAPH | naphthalene |
| 124TMB | 1,2,4-trimethylbenzene | OCT | octane |
| 12DCA | 1,2-dichloroethane | oXYL | o-xylene |
| 12DCB | 1,2-dichlorobenzene | PCE | tetrachloroethene |
| 135TMB | 1,3,5-trimethylbenzene | PENTADEC | pentadecane |
| 13DCB | 1,3-dichlorobenzene | PHEN | phenanthrene |
| 14DCB | 1,4-dichlorobenzene | t12DCE | trans-1,2-dichloroethene |
| 2MeNAPH | 2-methyl naphthalene | TCE | trichloroethene |
| BENZ | benzene | TMBs | combined masses of 1,3,5-trimethylbenzene and 1,2,4-trimethylbenzene |
| BTEX | combined masses of benzene, toluene, ethylbenzene, and total xylenes (Gasoline Range Aromatics) | TOL | toluene |
| C11,C13&C15 | combined masses of undecane, tridecane, and pentadecane (C11+C13+C15) (Diesel Range Alkanes) | TPH | total petroleum hydrocarbons |
| c12DCE | cis-1,2-dichloroethene | TRIDEC | tridecane |
| CCl4 | carbon tetrachloride | UNDEC | undecane |
| CHC13 | chloroform | VC | vinyl chloride |

SUMMARY OF SAMPLING RATE CALIBRATION FOR AGI UNIVERSAL SAMPLER IN AQUEOUS PHASES

PURPOSE:

The purpose of this document is to:

1. Summarize the test protocol,
2. Summarize the methodology for analysis of data,
3. Present general results for generating concentration calibration

of the AGI Universal Sampler, part number SPG-0008, in aqueous phase media following AGI's "Standard Practice for Determining the Sampling Rate of Passive Diffusion Samplers in Various Environmental Media": SPG-SOP-0493. The work will be summarized in three parts: Part 1: shallow water, Part 2: deep water, and Part 3: sediment.

Principle of Operation of the AGI Samper

The AGI Universal Sampler is designed with solid adsorbents enclosed inside a tubular microporous PTFE membrane. When placed in water, the pores and hydrophobic nature of the PTFE keep liquid water from entering the membrane until a water head of about 34 feet is reached. The membrane will not keep water vapor from entering but the adsorbents are very hydrophobic and through testing validated to be unaffected by this moisture vapor. In shallow water, <34', volatile and semi-volatile compounds will partition from the dissolved water into the air phase in the PTFE membrane according to Henry's Law. This partitioning is instantaneous and within seconds-minutes, the compound is adsorbed by the adsorbent inside the sealed tube. Because the diffusivity in air is about 10,000 times higher than the diffusivity in water, the sampling rate is controlled by the water contact area of the membrane that allows the Henry's Law effect to occur. This contact area is set by the membrane diameter and length of the sealed tube, which is fixed in AGI's manufacturing process.

Henry's law as well as diffusivity, which are fundamentally incorporated into the sampling rate, are affected by temperature, T , and follow an Arrhenius equation $H_T = H_r \times \exp[-E_a/R](1/T_r - 1/T)$. Because a 5°C temperature change can make a 15% change in sampling rate, the temperature of the sampled water should be known to get the most precise concentration.

The membrane pore size is also small enough that colloidal particles and microbes can not pass through the membrane. This keeps the adsorbent from getting contaminated and eliminates any need to add preservative or chilling during storage or transportation.

When the water pressure exceeds the water entry pressure of the membrane, about 34 feet of water, the water becomes in contact with the solid adsorbent. Under this condition the compounds in the water will partition from the water into the solid. The partitioning coefficient, K_{AW} , can be approximated by the octanol-water coefficient, K_{OW} , but has been measured more precisely in the lab for AGI's specific solid adsorbent. The sampling rate is the product of the sampling rate at <34' of water and the K_{AW} .

In sediment, the sampler measures pore-water concentration, which is generally agreed to be the preferred measurement as it is more indicative of bioavailability. In sediment the volumetric availability of water to the sampler is reduced by the volume fraction solids in the sediment, which typically varies from zero to 35%, but can be as high as 73% in well packed and broad particle size distribution sediments. As a result, sampling rates in sediment are multiplied by the fraction pore water in the sediment to determine concentration.

PART 1: Calibration in shallow water

Part 1 summarizes the work in shallow water generating calibration data, evaluating the physical and chemical factors affecting the sampling rate, and measurement of the actual sampling rates or regression calibration equations needed to determine concentrations.

Sample Generation in water

In this calibration work, solutions of analytes at known concentrations were formulated in clean 4 liter smoked glass jugs by injecting microliter measured amounts of environmental standards using a calibrated syringe into pure or deionized water and stirring for a minimum of 2 hours but generally overnight. Headspace in the jugs was minimized and generally less than 1% by volume during the tests. Jugs were temperature controlled by placing them in a water filled cooler, chilled via a copper tubing loop in the cooler. Temperature was measured with a certified digital temperature gauge and an average value used for each temperature experiment.

AGI samplers were weighted so they won't float and placed in the jugs at time zero. They were removed at various intervals to generate samples along with duplicates that showed mass increasing with exposure time. The sampler exposure time was selected to span minutes to hours and was generally reduced for high concentration tests to maintain uptake with time in roughly the linear dynamic range. Samplers were removed and dried with a paper towel and returned to their original container for analysis. They were analyzed by AGI's 8260C (SPG-WI-318 or SPG-WI-10028) method in duplicate, which is based on EPA SW846 Method 8260C.

Water samples were also taken and measured at an outside accredited lab using EPA SW846 Method 8260B. The concentrations agreed well with the calculated concentrations based on the standard certification, jug volume, and syringe injection. The variability of the outside lab 8260B values were found to be high, so for the sampling rate calculations we used the concentrations based on syringe dosing.

Calibrations were run at five concentrations, nominally at 6, 24, 118, 590, 1420 ug/L and five temperatures nominally at 5, 10, 15, 20, and 25 degrees centigrade. Samples were taken at 4 different exposure times. Samples were run in duplicate. A total of 176 data points were generated using 28 compounds from AGI's standard compounds list. Tridecane and pentadecane were not evaluated due to their very low solubility in water. In addition, another 23 compounds were tested using an 8260 liquid standard at nominal concentrations of 0.5, 1.0, 5.0, 15, 95, and 470 ug/L at a temperature typical of groundwater, 15°C. This is a living calibration and as additional data are generated, they may be qualified and added to this data set to improve the precision of the sampling rate calibration and broaden the compound list.

Key Variable Effects

As expected from theory, at short to moderate exposure times, mass will increase roughly linearly proportional to exposure time, as well as proportional to concentration, and exponentially with temperature following Arrhenius law. Temperature affects the Henry's law as well as diffusivity in water. Sampling rate is generally independent of concentration and time at mass values significantly below saturation. In the following sections we have characterized the sampling rate for each compound as affected by temperature and also developed calibrations using regression which account for the minor impact of time, and mass.

Concentration using Simple Sampling Rate Determination

A simple way to determine concentration is to measure mass on the AGI sampler, divide by exposure time, and divide by sampling rate, SR.

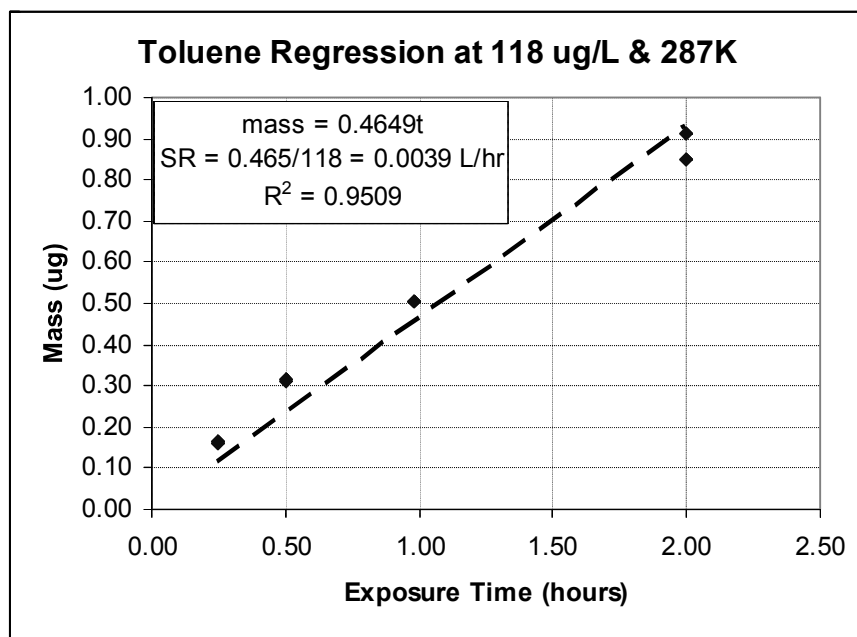
$$\text{Conc [ug/L]} = \text{mass/time/SR} \quad (1)$$

The sampling rate can be determined via measurements of mass versus time at a known concentration and temperature according to the following modification of equation (1).

$$\text{SR} = \text{mass/time/concentration} \quad (2)$$

Sampling rates in L/hr were determined by measuring the trend or regression mass uptake versus time and dividing by the concentration. A measurement like this will use 8 data points (4 times x 2 samples). Such a sampling rate can be measured at any concentration and temperature.

The chart to the right shows a plot of mass versus time for water at 118 ug/L and 287K. This is actual data from a single run. Slope of 0.465 ug/hr divided by the concentration of 118 ug/L yields a sampling rate, SR, of 0.0039 L/hr.



SR's typically range from about 0.004 to 0.007 L/hr at 15°C. Table A shows SR's measured for our standard compound list at 5 temperatures.

Rigorous Concentration using Regression

A preferred method for determining concentration that will yield improved accuracy over a wide range of concentrations, exposure times, and temperatures is to use all data in a regression analysis. This allows adjustments for the minor non-linear influences of mass and time as well as the effects of temperature. This is done by regressing equation (1) or a universal version of equation (1)

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRo} \cdot \exp(-E_a/R/T)] \quad (3)$$

The subtle non-linear effects of mass and time will be evident in the deviation of coefficients b and d from 1.0. This regression generates four constants b, d, SRo, and $-E_a/R$ by regressing $\ln(\text{conc})$ versus $\ln(\text{mass})$, $\ln(\text{time})$, $1/\text{temp}$. These four constants can be used to determine concentration via the equation:

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRo} \times \exp(-E_a/R(1/T))] \quad (4)$$

Where conc is in ug/L, mass is in ug, time in hours, T in degrees Kelvin.

Equation (4) can be also expressed at a reference temperature, T_r , such as 15°C by

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRr} \times \exp(-E_a/R(1/T_r - 1/T))] \quad (5)$$

This allows sampling rates, SRr, at any reference temperature, T_r , and for any analyte to easily be compared. These values of SRr at 293.14K can be found in Table A.

When sampling times are between 0 and 4 hours, using the 4 constant equation (5) is preferred. For concentrations from about 5 to 1500 ug/L one hour exposure times generally give the lowest error, typically with average error of 6-20% and with total error range of 12%-32%. For low concentrations where sampling times are greater than 4 hours, it is preferred to use equation (1) to avoid unrealistic effects from the coefficient d or to set d to 1.0. In such a case SR in equation (1) can be substituted with $[\text{SRr} \times \exp(-E_a/R(1/T_r - 1/T))]$ to use an SR representative of the well temperature, T.

The chart to the right is a plot of the calculated concentration from the 4 constant regression compared to the dosed concentration. Agreement is excellent for the 176 data points.

There does appear to be a slight high bias of 8.6% over the full range of this data, although it is well within acceptable limits of variability.

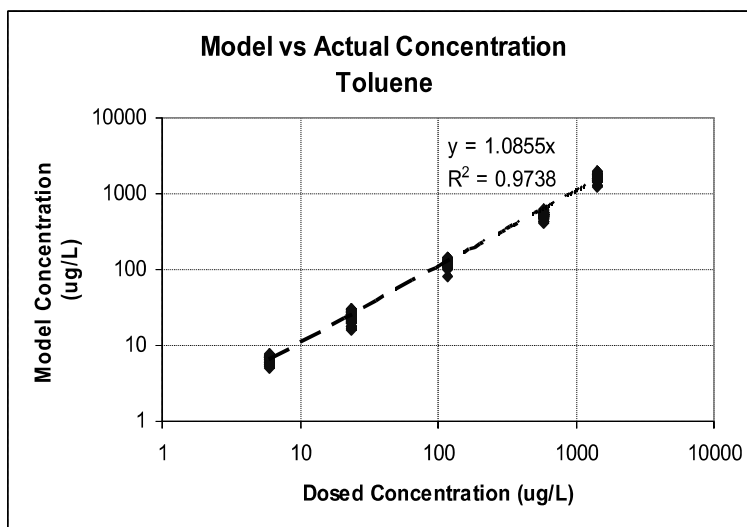


Table B shows the tabulated summary of the 4 constants regression with Rsq values and error estimates for the 4 constants for each analyte. Most regression Rsq values are 0.99 or greater for each analyte. In general, $-E_a/R$ is about 2400 \pm 400, b is about 0.9, d is about -0.75, and SR(15°C) ranges from .004 L/hr to 0.007 L/hr increasing with MW of the compound.

Error Estimates

The error in the water concentration values will depend on both the error in mass from the analytical method as well as the error in the concentration calibration. Table C shows the error in the mass values from the 8260C low sensitivity method.

The standard error of the regression and standard errors of the constants can be found in table B. For each compound we have measured the error between the derived concentration and the actual concentration. The error tends to be lowest at our recommended exposure time of one hour as shown by the example for Toluene to the right.

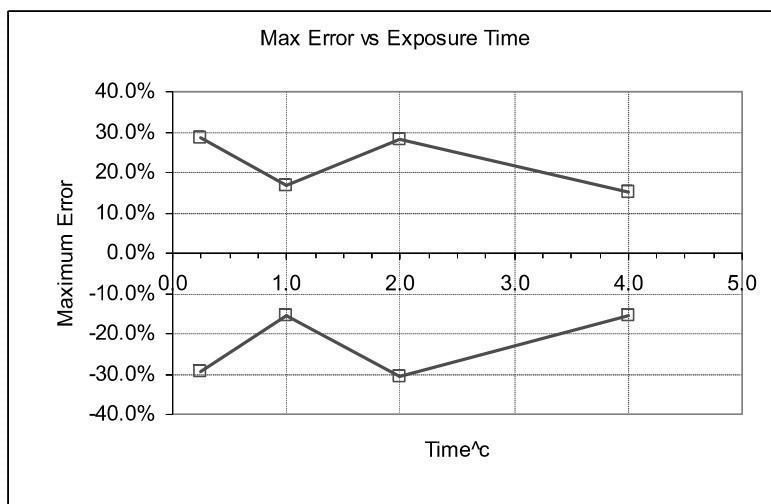
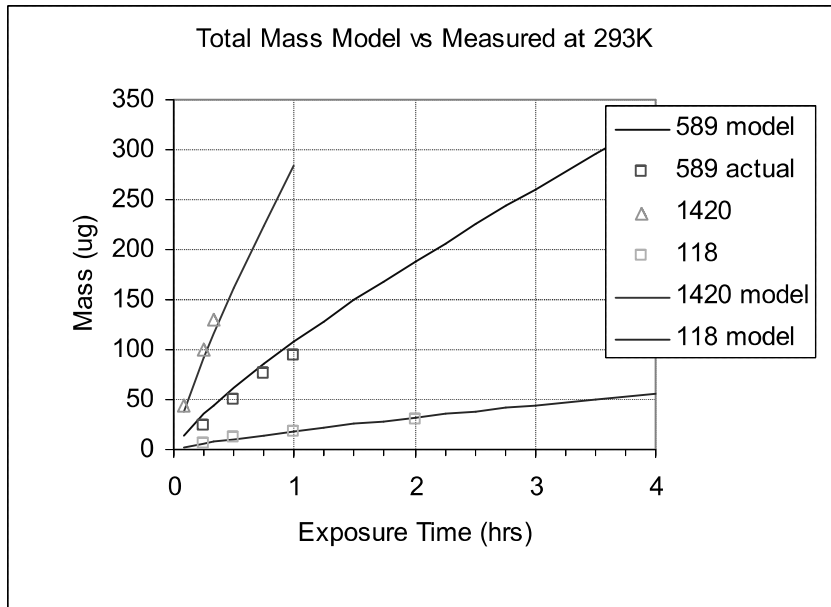


Table D shows the total average error in water concentration by compound as well as the low and high error. The average ranges from about 6% to 20%, which is similar to the analytical method errors. The low and high errors range from 12% to 32% and include contribution from measurement errors in both time and temperature.

Sorbent Saturation

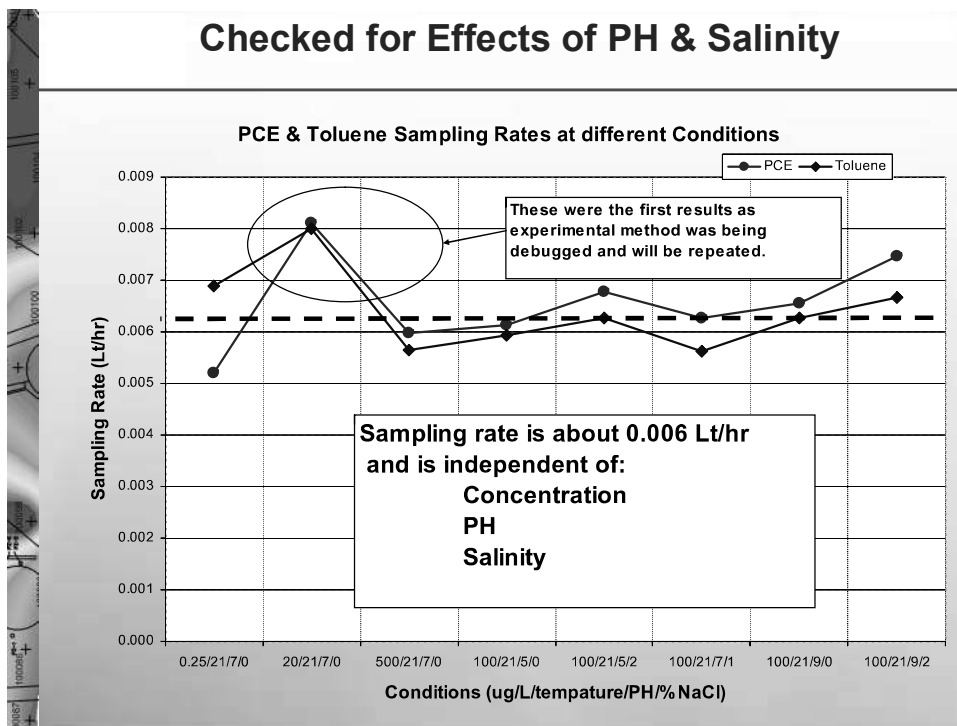
As mass increases on a solid sorbent and approaches saturation, reverse diffusion can occur causing the sampling rate to drop. Eventually the mass level will reach a maximum steady state value at any concentration. A rate of mass uptake with time that deviates significantly from linear, indicates that sorbent saturation could be an issue. When using equation (1), staying in the linear range to avoid the effects of adsorbent saturation is important. We recommend keeping the total mass on the sampler below 50 ug or flagging when this is exceeded.

The 4 constant regression accounts for some of the non linearity allowing good accuracy at higher mass levels. From the experimental data we have found this safe range can be extended to 100 ug or higher as shown in the chart below. This chart compares total mass of all compounds (excluding heavy alkanes, which have solubility issues) versus time in comparison to that predicted from the 4-constant concentration equation.



Effect of PH and Salinity

Because neither PH nor salinity is known to have a significant impact on Henry's law or diffusivity in water, we did not expect them to have a significant impact on sampling rate. To confirm this, experiments were run varying PH from 5 to 9 and NaCl content from 0 to 2%. The chart below shows no significant impact for combinations of PH and NaCl content over this range on the sampling rate of toluene in water at 21°C.

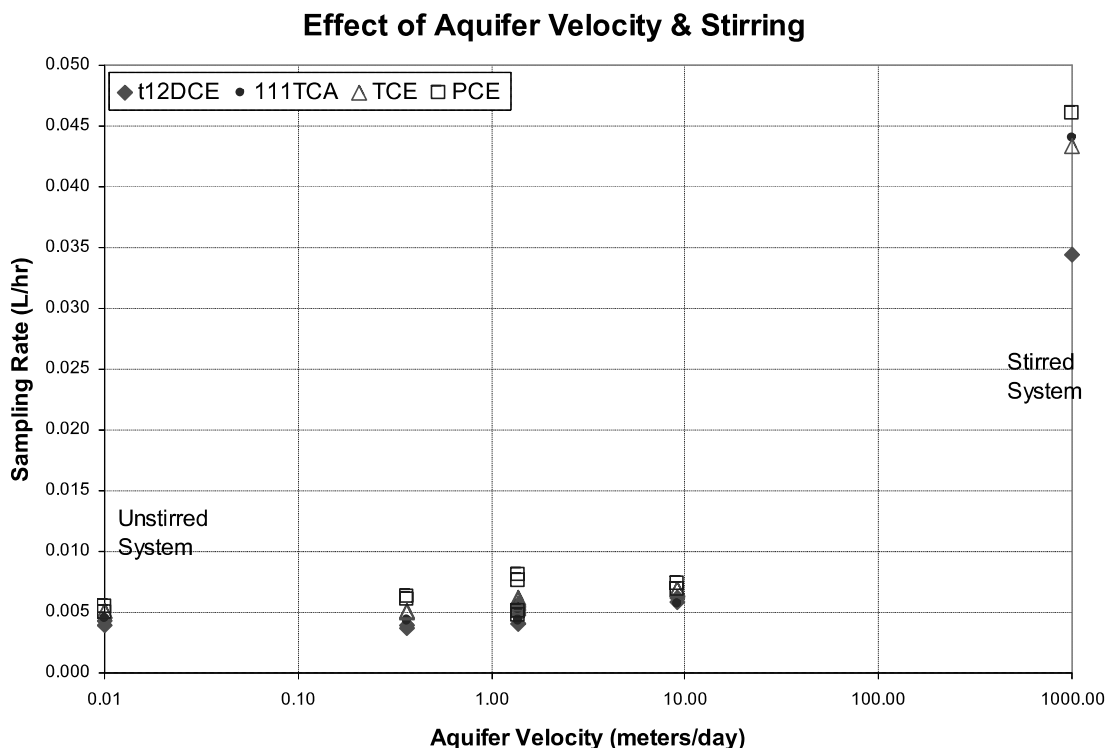


Impact of Aquifer Velocity

The velocity in most aquifers is quite slow, typically a meter/day or less. Occasionally water flow could be much higher such as encountered in karst aquifers, streams or rivers. Mass transfer coefficients are higher in high flow conditions, which will lead to higher sampling rates. We validated that a highly stirred system had sampling rates about 10 times higher than those that were non-stirred. We decided to evaluate the effect of aquifer velocity.

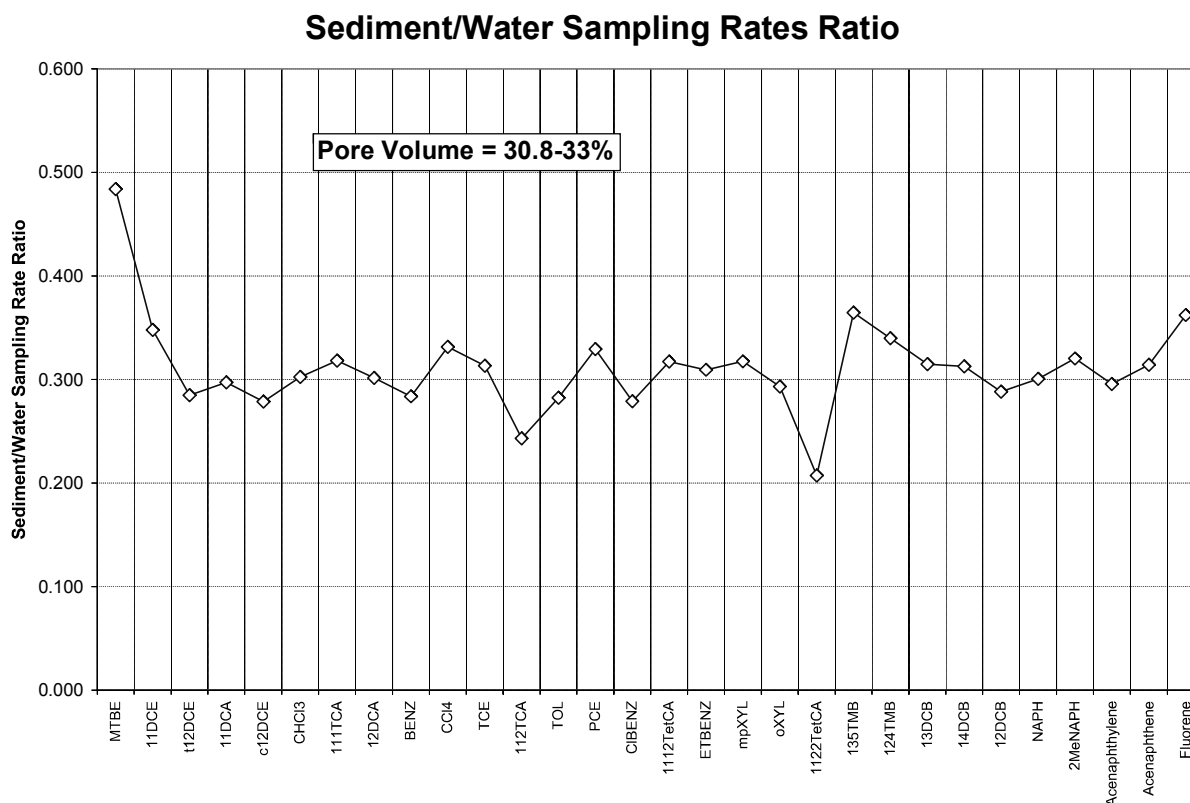
A test apparatus was built comprising a 3" PVC pipe tee filled with clean sand in each of the horizontal straight legs and screened to leave the center open. A test solution was run through this system using a variable flow pump and AGI samplers were placed into the simulated well through the vertical leg of the tee. Tests were run to examine the effect of velocity by varying the pumping rate and hence water velocity.

The chart below shows no significant effect of aquifer velocity up to a speed of about 10 meters/day. At velocities significantly above this, similar to a stirred system, sampling rates are about 10 times higher.



concentration. Pore water fraction can range from 1.0 for water without sediment to as low as 0.25. Typically most sediments have pore fractions of 0.9 to 0.65.

A sampling rate study was done with water and with water added into a well-packed sorted sand. Pore water fraction in this test was measured between 30.8% 33% by volume. Below is a plot of the ratio of sampling rates measured in the sediment to open water. The average ratio is equal to the pore water fraction confirming that sampling rate in sediment is on average equal to the product of pore water fraction times the sampling rate in water.



Summary

The AGI® Sampler can be used to determine the concentration of volatile and semi-volatile compounds in a water phase. This requires knowing the exposure time and water temperature. It also requires knowing if the sample is above or below 34' of water head and if the water has a velocity above 10 meters/day. Regressions of large amounts of data were used to generate a four constant equation to generate concentration values in water. Potential error in the concentration values is excellent typically less than 25%.

TABLE A
WATER SAMPLING RATES STANDARD LIST

| | SRr 293.14 | SR @ 277.54 | SR @ 282.44 | SR @ 287.84 | SR @ 293.24 | SR @ 298.94 |
|-------------------|----------------------|-----------------------|-----------------------|-----------------------|-----------------------|-----------------------|
| MTBE | 0.0025 | 0.0014 | 0.0016 | 0.0018 | 0.0022 | 0.0029 |
| t12DCE | 0.0043 | 0.0028 | 0.0028 | 0.0027 | 0.0037 | 0.0048 |
| 11DCA | 0.0047 | 0.0031 | 0.0033 | 0.0033 | 0.0039 | 0.0052 |
| c12DCE | 0.0046 | 0.0030 | 0.0031 | 0.0031 | 0.0038 | 0.0051 |
| CHCl3 | 0.0046 | 0.0030 | 0.0031 | 0.0031 | 0.0038 | 0.0051 |
| 111TCA | 0.0066 | 0.0043 | 0.0047 | 0.0047 | 0.0056 | 0.0076 |
| 12DCA | 0.0045 | 0.0029 | 0.0029 | 0.0030 | 0.0036 | 0.0050 |
| BENZ | 0.0050 | 0.0031 | 0.0034 | 0.0035 | 0.0042 | 0.0056 |
| CCl4 | 0.0068 | 0.0044 | 0.0048 | 0.0047 | 0.0058 | 0.0080 |
| TCE | 0.0052 | 0.0030 | 0.0034 | 0.0034 | 0.0043 | 0.0058 |
| 112TCA | 0.0043 | 0.0027 | 0.0027 | 0.0028 | 0.0034 | 0.0048 |
| TOL | 0.0056 | 0.0034 | 0.0039 | 0.0039 | 0.0047 | 0.0062 |
| OCT | 0.0064 | 0.0046 | 0.0050 | 0.0040 | 0.0058 | 0.0089 |
| PCE | 0.0061 | 0.0036 | 0.0043 | 0.0043 | 0.0051 | 0.0069 |
| CIBENZ | 0.0054 | 0.0033 | 0.0039 | 0.0040 | 0.0045 | 0.0059 |
| 1112TetCA | 0.0061 | 0.0037 | 0.0042 | 0.0044 | 0.0050 | 0.0065 |
| EtBENZ | 0.0060 | 0.0037 | 0.0045 | 0.0044 | 0.0052 | 0.0069 |
| mpXYL | 0.0064 | 0.0039 | 0.0048 | 0.0046 | 0.0055 | 0.0072 |
| oXYL | 0.0066 | 0.0041 | 0.0050 | 0.0048 | 0.0057 | 0.0074 |
| 1122TetCA | 0.0044 | 0.0027 | 0.0029 | 0.0031 | 0.0036 | 0.0046 |
| 135TMB | 0.0079 | 0.0046 | 0.0059 | 0.0056 | 0.0071 | 0.0093 |
| 124TMB | 0.0078 | 0.0046 | 0.0060 | 0.0055 | 0.0071 | 0.0092 |
| 13DCB | 0.0072 | 0.0041 | 0.0055 | 0.0053 | 0.0063 | 0.0080 |
| 14DCB | 0.0071 | 0.0040 | 0.0054 | 0.0052 | 0.0062 | 0.0079 |
| 12DCB | 0.0070 | 0.0040 | 0.0053 | 0.0051 | 0.0060 | 0.0076 |
| UNDEC | | 0.0026 | 0.0024 | 0.0020 | 0.0031 | 0.0029 |
| NAPH | | 0.0041 | 0.0056 | 0.0054 | 0.0064 | 0.0081 |
| TRIDEC | | | | | | |
| 2MeNAPH | | 0.0043 | 0.0066 | 0.0066 | 0.0080 | 0.0108 |
| PENTADEC | | | | | | |
| Total mass | 0.1177 | 0.0822 | 0.1339 | 0.1334 | 0.1773 | 0.1981 |

Values in L/hr

Total mass does not include UNDEC, TRIDEC, PENTADEC (28 compounds)

TABLE B
4 CONSTANT REGRESSION OUTPUT

| | Adjusted | Standard | | | | | Std | Std | Std | Std |
|------------|----------|----------|---------|-------|-----------|--------|------------------|------------|--------------------|------------|
| | Rsq | Error | ln(SR0) | b | - Ea/R | d | Error ln(SR0) | Error b | Error - Ea/R | Error d |
| MTBE | 0.997 | 0.0960 | -3.217 | 0.981 | 2704 | -0.709 | 0.2881 | 0.0062 | 83 | 0.0082 |
| t12DCE | 0.992 | 0.1659 | -1.877 | 0.905 | 2147 | -0.760 | 0.4971 | 0.0100 | 144 | 0.0138 |
| 11DCA | 0.995 | 0.1272 | -1.346 | 0.916 | 1965 | -0.746 | 0.3809 | 0.0077 | 110 | 0.0106 |
| c12DCE | 0.995 | 0.1299 | -1.905 | 0.911 | 2137 | -0.751 | 0.3892 | 0.0078 | 112 | 0.0109 |
| CHCl3 | 0.996 | 0.1260 | -1.841 | 0.912 | 2118 | -0.748 | 0.3776 | 0.0076 | 109 | 0.0105 |
| 111TCA | 0.995 | 0.1279 | -2.684 | 0.902 | 2259 | -0.761 | 0.3836 | 0.0076 | 111 | 0.0106 |
| 12DCA | 0.995 | 0.1263 | -2.161 | 0.908 | 2218 | -0.746 | 0.3786 | 0.0076 | 109 | 0.0106 |
| BENZ | 0.995 | 0.1323 | -2.207 | 0.920 | 2198 | -0.754 | 0.3965 | 0.0080 | 114 | 0.0110 |
| CCl4 | 0.994 | 0.1405 | -3.121 | 0.889 | 2379 | -0.776 | 0.4220 | 0.0083 | 122 | 0.0116 |
| TCE | 0.992 | 0.1655 | -3.338 | 0.900 | 2522 | -0.772 | 0.4969 | 0.0099 | 144 | 0.0137 |
| 112TCA | 0.995 | 0.1264 | -2.412 | 0.896 | 2302 | -0.724 | 0.3790 | 0.0075 | 109 | 0.0107 |
| TOL | 0.994 | 0.1426 | -2.873 | 0.916 | 2364 | -0.756 | 0.4281 | 0.0087 | 124 | 0.0119 |
| OCT | 0.938 | 0.4698 | -5.984 | 0.822 | 3235 | -0.827 | 1.4231 | 0.0277 | 412 | 0.0388 |
| PCE | 0.991 | 0.1773 | -3.780 | 0.877 | 2601 | -0.775 | 0.5329 | 0.0103 | 154 | 0.0147 |
| CIBENZ | 0.994 | 0.1457 | -2.601 | 0.911 | 2292 | -0.747 | 0.4370 | 0.0088 | 126 | 0.0122 |
| 1112TetCA | 0.996 | 0.1235 | -2.676 | 0.898 | 2281 | -0.725 | 0.3705 | 0.0073 | 107 | 0.0104 |
| EtBENZ | 0.993 | 0.1597 | -2.930 | 0.918 | 2357 | -0.752 | 0.4794 | 0.0097 | 138 | 0.0134 |
| mpXYL | 0.992 | 0.1678 | -3.036 | 0.909 | 2372 | -0.749 | 0.5037 | 0.0101 | 145 | 0.0140 |
| oXYL | 0.993 | 0.1555 | -2.862 | 0.911 | 2312 | -0.740 | 0.4667 | 0.0094 | 135 | 0.0131 |
| 1122TetCA | 0.996 | 0.1118 | -1.971 | 0.913 | 2167 | -0.691 | 0.3351 | 0.0067 | 97 | 0.0096 |
| 135TMB | 0.988 | 0.2024 | -4.435 | 0.897 | 2720 | -0.738 | 0.6093 | 0.0121 | 176 | 0.0170 |
| 124TMB | 0.989 | 0.1997 | -4.126 | 0.890 | 2631 | -0.731 | 0.6009 | 0.0118 | 173 | 0.0169 |
| 13DCB | 0.991 | 0.1832 | -3.422 | 0.888 | 2449 | -0.730 | 0.5503 | 0.0108 | 159 | 0.0155 |
| 14DCB | 0.991 | 0.1802 | -3.263 | 0.892 | 2408 | -0.724 | 0.5413 | 0.0107 | 156 | 0.0153 |
| 12DCB | 0.992 | 0.1697 | -2.970 | 0.894 | 2327 | -0.716 | 0.5092 | 0.0101 | 147 | 0.0144 |
| UNDEC | 0.694 | 0.374 | -1.406 | 0.426 | 1708 | 0.806 | 1.792 | 0.0028 | 517 | 0.053 |
| NAPH | 0.992 | 0.166 | -3.374 | 0.915 | 2430 | 0.671 | 0.497 | 0.010 | 144 | 0.014 |
| TRIDEC | | | | | | | | | | |
| 2MeNAPH | 0.984 | 0.238 | -5.498 | 0.869 | 2990 | 0.689 | 0.72 | 0.014 | 208 | 0.021 |
| PENTADEC | | | | | | | | | | |
| Total mass | 0.993 | 0.1543 | -6.111 | 0.907 | 2419 | -0.732 | 0.4666 | 0.0093 | 134 | 0.0130 |

TABLE C 8260C MASS UNCERTAINTY

AGI 8260C Method for Mass using SPG-0008 Samplers

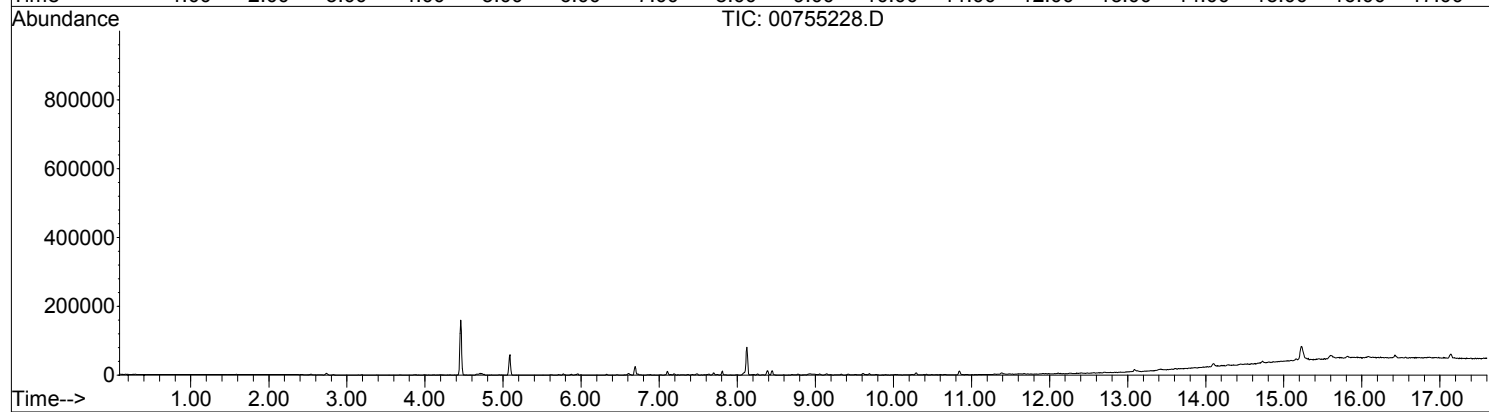
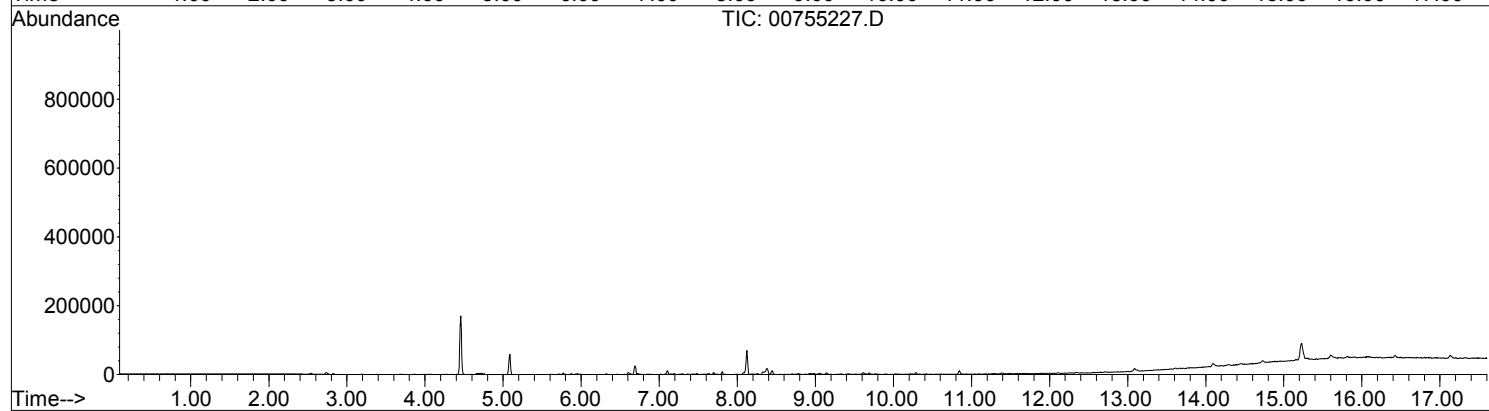
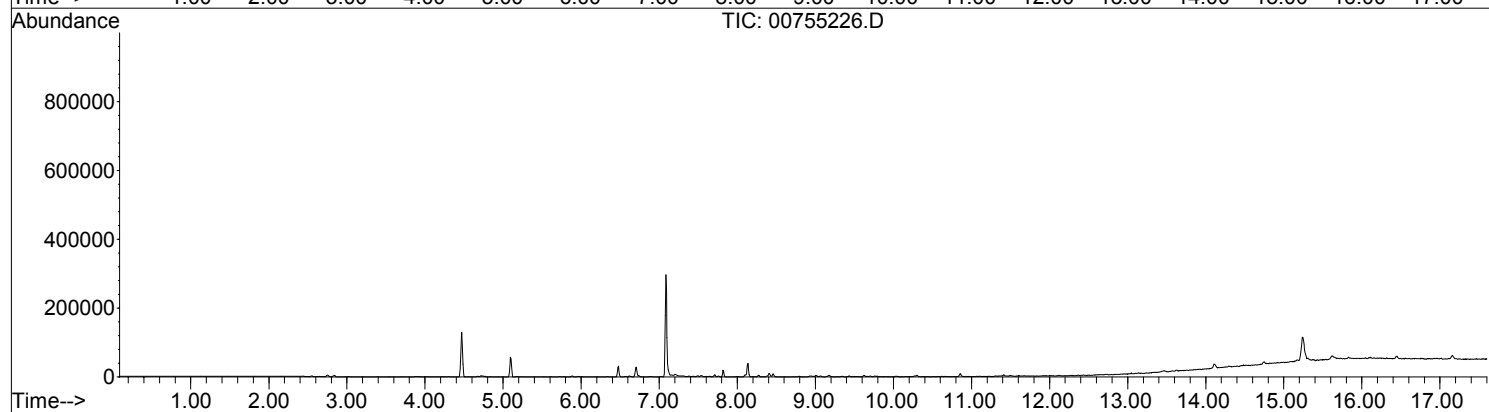
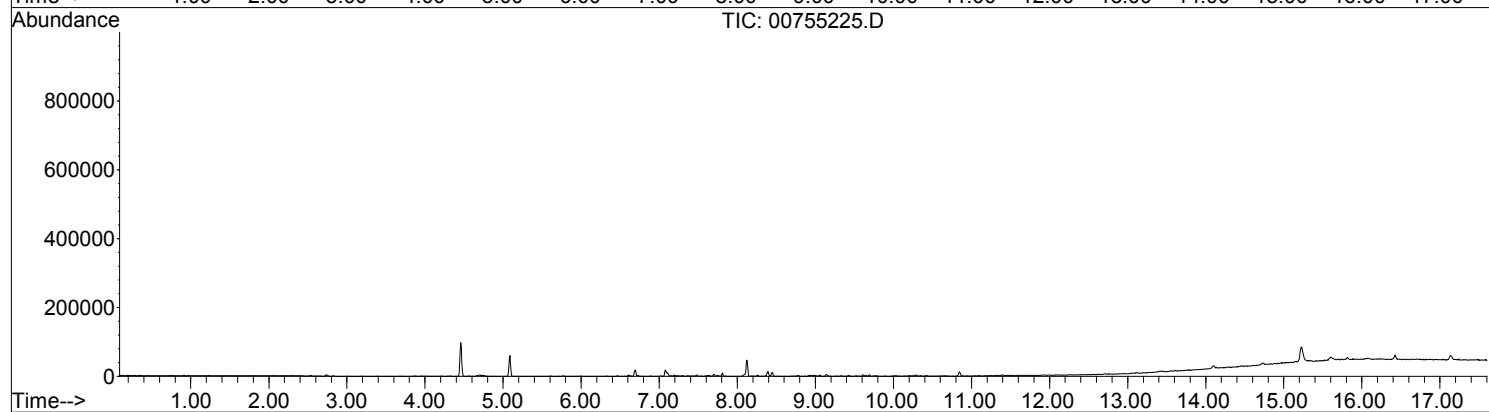
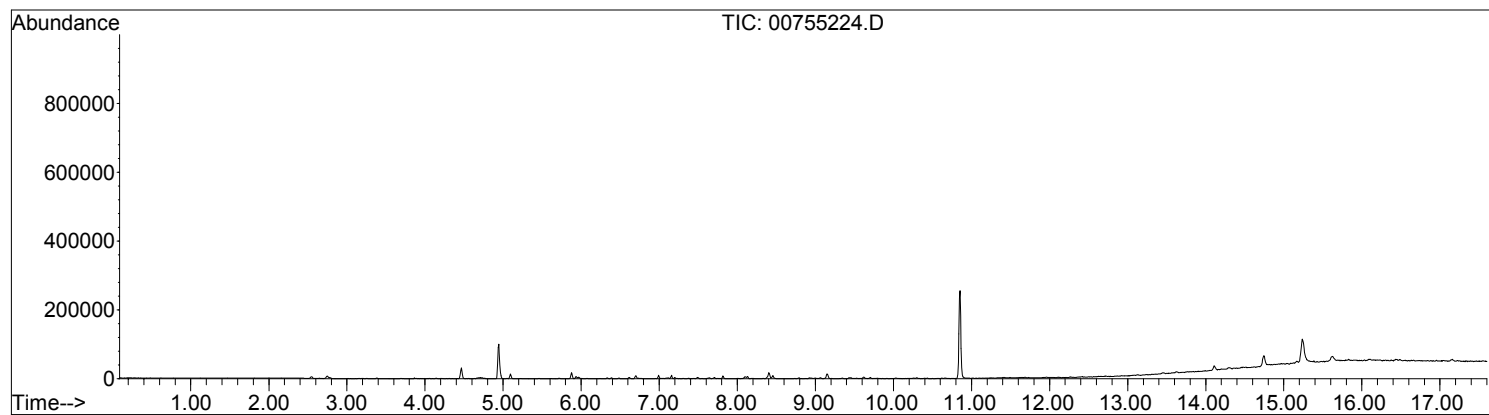
| | 99% Uncertainty Range +/- | 95% Uncertainty Range +/- |
|-----------|---------------------------------|---------------------------------|
| MTBE | 20% | 14% |
| t12DCE | 22% | 15% |
| 11DCA | 18% | 12% |
| c12DCE | 18% | 12% |
| CHCl3 | 16% | 11% |
| 111TCA | 18% | 12% |
| 12DCA | 20% | 13% |
| BENZ | 16% | 10% |
| CCl4 | 19% | 12% |
| TCE | 15% | 10% |
| 112TCA | 18% | 12% |
| TOL | 15% | 10% |
| OCT | 20% | 13% |
| PCE | 16% | 11% |
| CIBENZ | 18% | 12% |
| 1112TetCA | 19% | 13% |
| EtBENZ | 18% | 12% |
| mpXYL | 18% | 12% |
| oXYL | 18% | 12% |
| 1122TetCA | 23% | 15% |
| 135TMB | 21% | 14% |
| 124TMB | 20% | 14% |
| 13DCB | 19% | 13% |
| 14DCB | 19% | 13% |
| 12DCB | 20% | 14% |
| NAPH | 21% | 14% |
| 2MeNAPH | 25% | 17% |

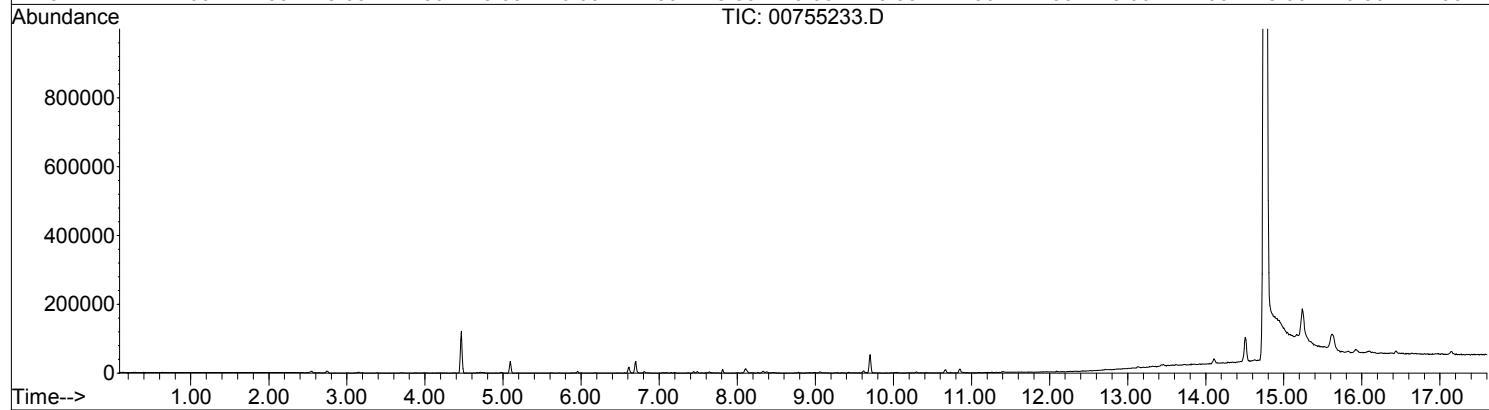
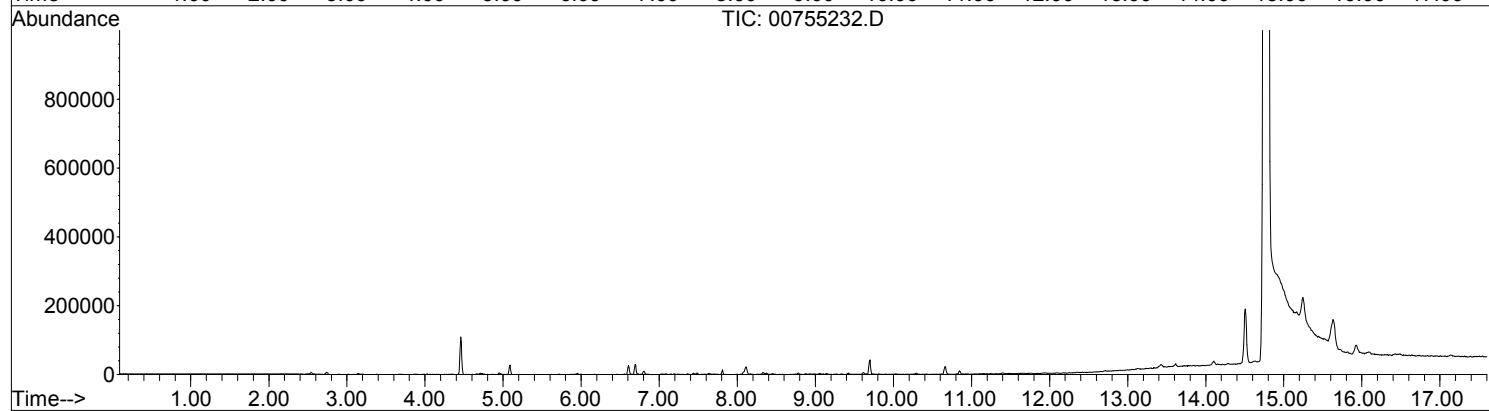
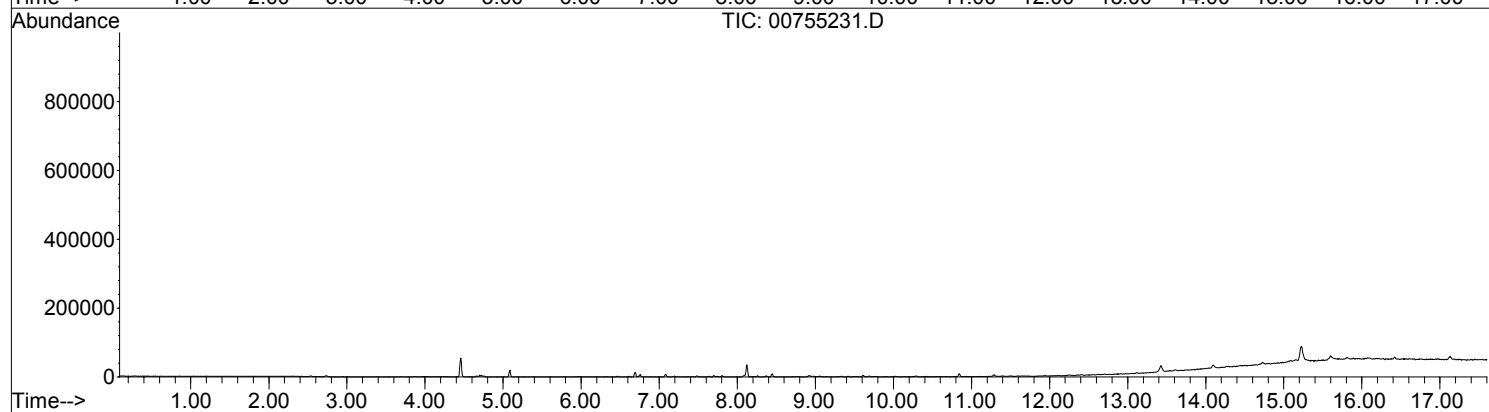
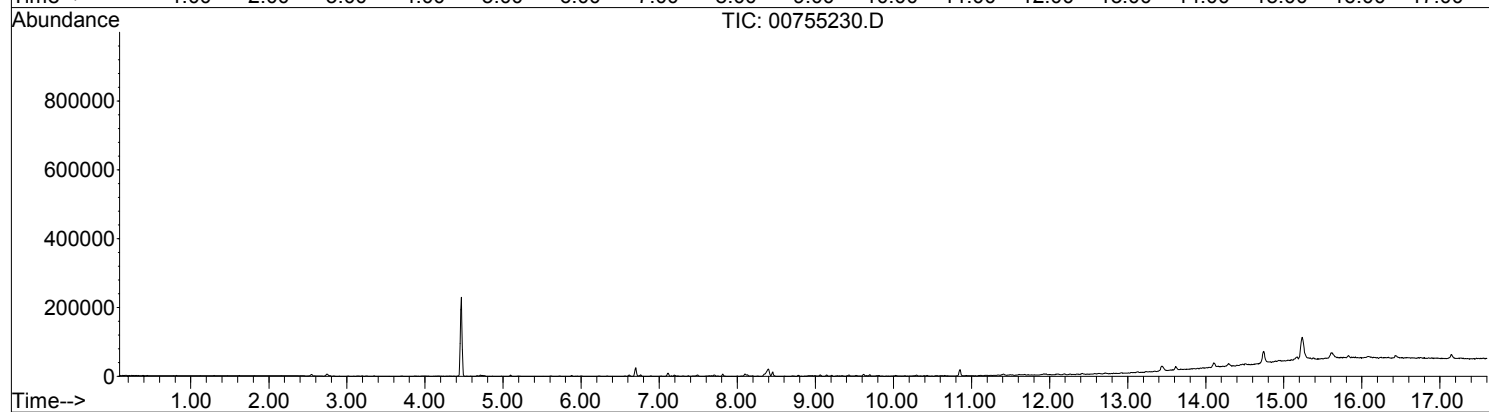
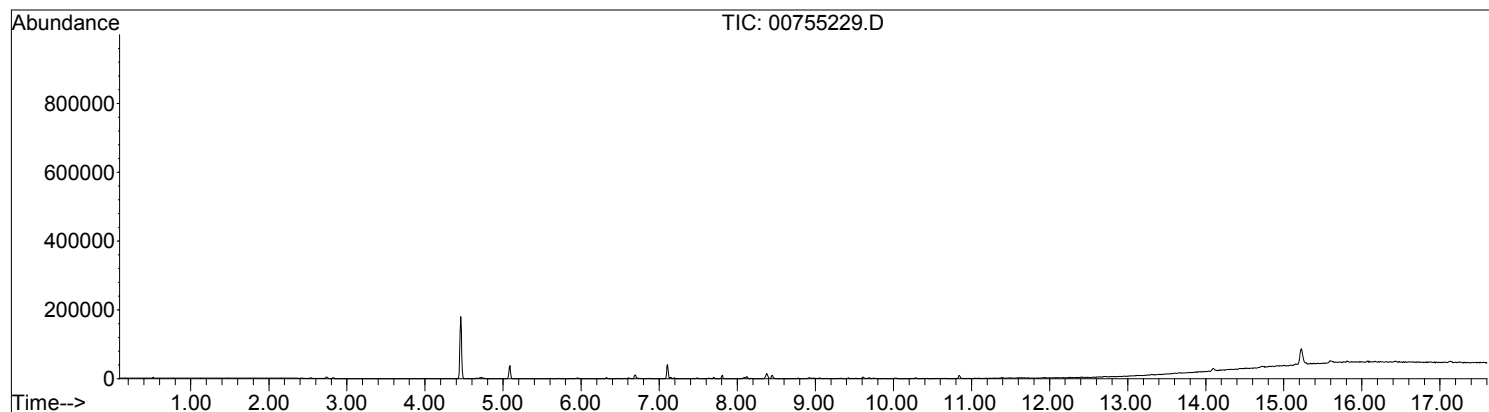
TABLE D
4 CONSTANT WATER CONCENTRATION UNCERTAINTY

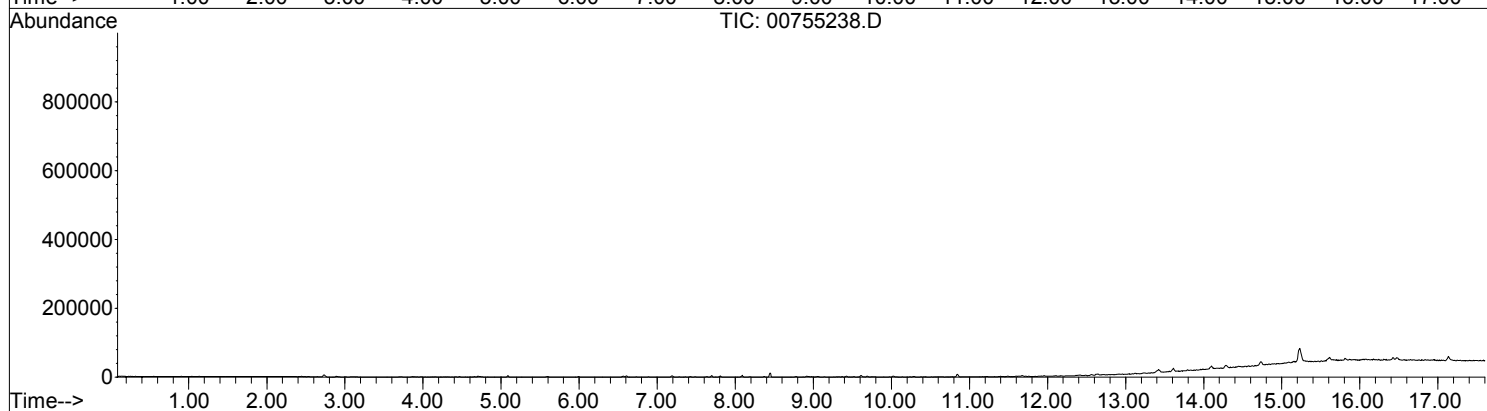
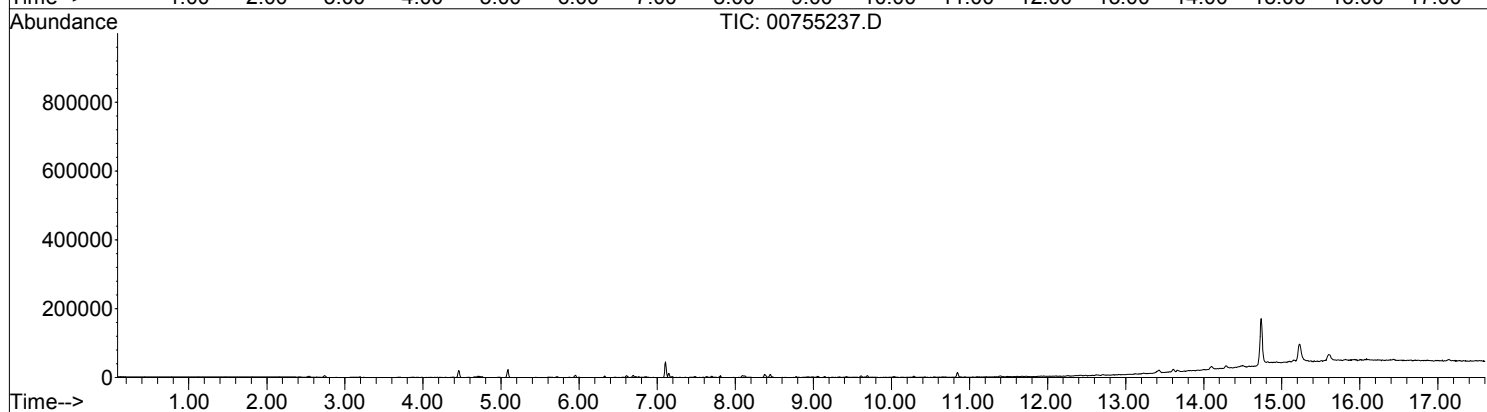
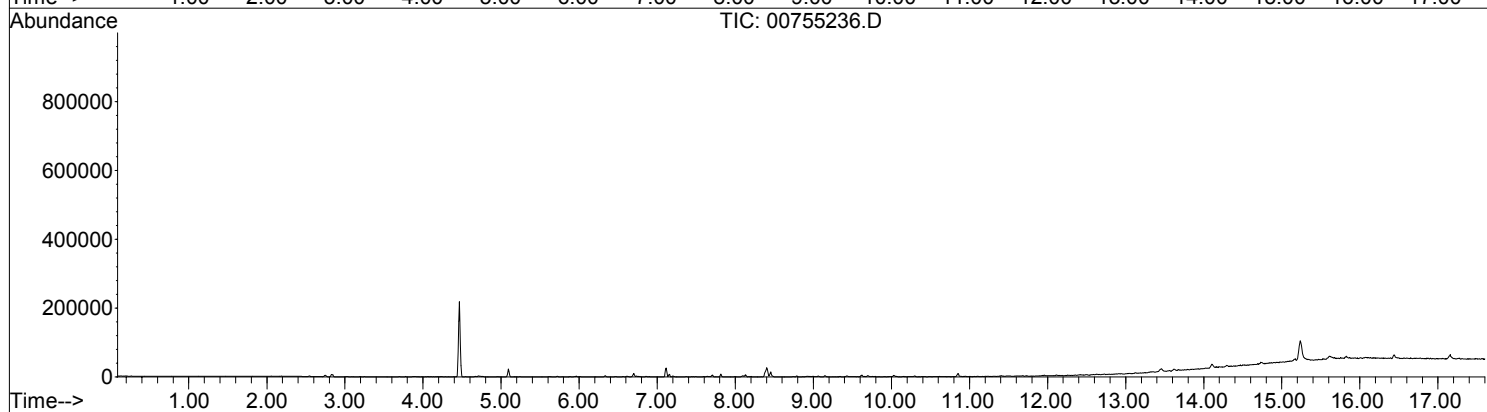
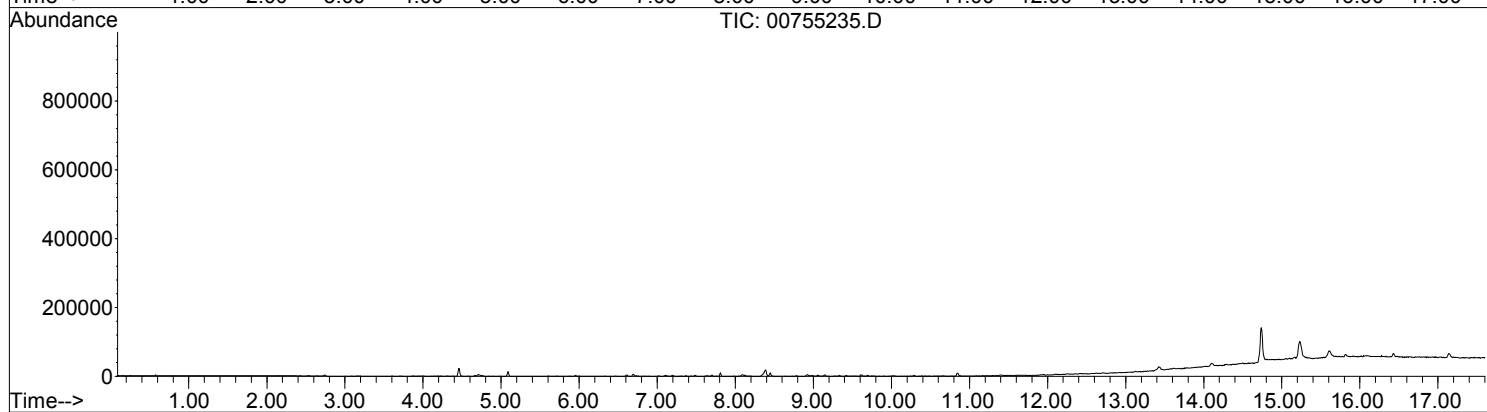
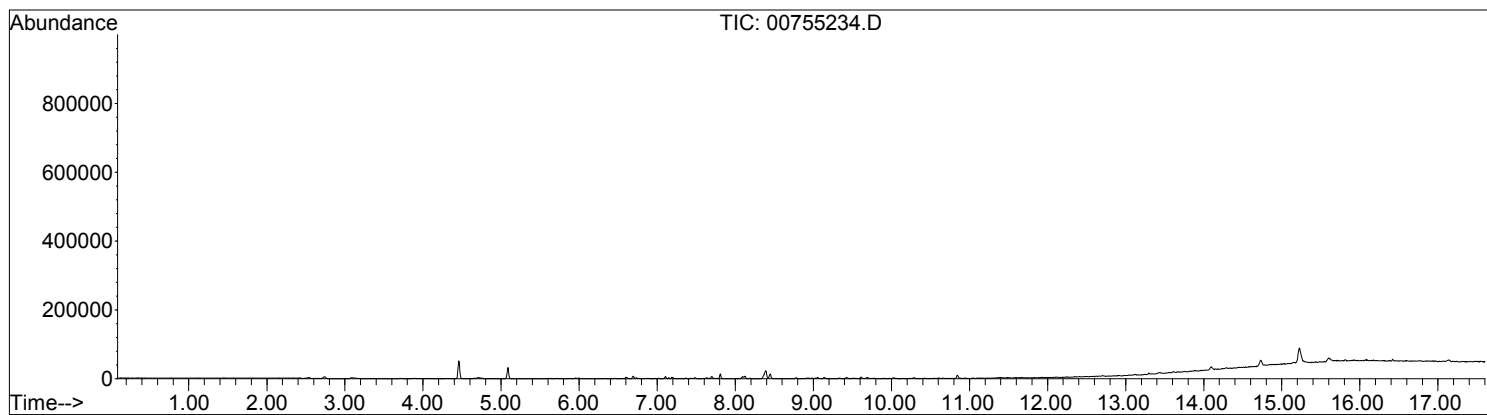
ERROR IN CONCENTRATION REPORTING (1)

| | Average Error | Minimum Error | Maximum Error |
|-----------|--------------------------|--------------------------|--------------------------|
| MTBE | 6% | -12% | 12% |
| t12DCE | 11% | -26% | 21% |
| 11DCA | 8% | -19% | 13% |
| c12DCE | 9% | -19% | 15% |
| CHCl3 | 9% | -20% | 14% |
| 111TCA | 9% | -19% | 23% |
| 12DCA | 10% | -19% | 17% |
| BENZ | 8% | -18% | 13% |
| CCl4 | 10% | -23% | 22% |
| TCE | 10% | -21% | 14% |
| 112TCA | 11% | -21% | 21% |
| TOL | 7% | -17% | 14% |
| OCT | 20% | -41% | 42% |
| PCE | 10% | -24% | 15% |
| CIBENZ | 7% | -16% | 14% |
| 1112TetCA | 8% | -17% | 18% |
| EtBENZ | 6% | -19% | 14% |
| mpXYL | 7% | -22% | 13% |
| oXYL | 7% | -19% | 13% |
| 1122TetCA | 8% | -16% | 17% |
| 135TMB | 9% | -23% | 17% |
| 124TMB | 10% | -28% | 19% |
| 13DCB | 10% | -22% | 17% |
| 14DCB | 10% | -22% | 17% |
| 12DCB | 9% | -23% | 17% |
| NAPH | 10% | -24% | 21% |
| 2MeNAPH | 13% | -32% | 30% |

(1) For 1 hour exposure, includes error related to mass value from AGI analytical method 8260C









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AMPLIFIED
GEOCHEMICAL
IMAGING, LLC

Laboratory Report

Site: Comal and San Marcos Rivers

Prepared for:

SWCA Environmental Consultants
10245 Little York Road Suite 600
San Antonio, TX 77040
UNITED STATES

Prepared on:
May 08, 2015

Project Summary and Objective

Amplified Geochemical Imaging, LLC. (AGI) provided the AGI Environmental Survey used at:

Comal and San Marcos Rivers

The service provided by AGI included delivery of the required quantity of AGI Universal Samplers, analysis by the method described below for the requested organic compounds, reporting of the data, and contour mapping (as needed).

This report includes results for only the samples noted under the Laboratory Sample Report section. If contour maps are part of the project deliverable, the maps will be prepared and issued under a separate report cover, upon receipt of a usable sitemap (electronic) and compound choices for contouring.


Written/submitted by:

Jim E Whetzel

Project Manager

Reviewed/approved by:

Don D'Apolito



Project Manager

Analytical data approved by:

Jasmine Smith

Chemist

Quality Assurance Statement

The AGI Laboratory, at Amplified Geochemical Imaging's facility in Elkton, MD USA, operates under the guidelines of its ISO Standard 17025 DoD ELAP accreditation, and its Quality Assurance Manual, Operating Procedures, and Methods (SPG-SOP-0462).

For this project, the analytical method, results, and observations reported do [] do not [✓] fall within the scope of AGI's ISO 17025 accreditation.

Screening/Concentration Method

The AGI Universal Samplers are analyzed at AGI's fixed laboratory using thermal desorption-gas chromatography/mass spectrometry (TD-GC/MS) instrumentation following U.S. EPA Method 8260 (SPG-WI-0292) which includes the following:

- **BFB Tuning Frequency:** A BFB tune is analyzed at the start of each analytical run and after every 30 samples.
- **Initial Calibration:** A minimum of a five point calibration curve is analyzed prior to the analysis of samples.
- **Linearity of Target Compounds:** If the RSD of any target analyte is less than or equal to 25% then average response factor can be used for quantitation. If the RSD exceeds 25% for a target compound a regression equation can be used for quantitation.
- **Continuing Calibration Verification:** After every 10 samples, and at the end of each analytical batch, and a second-source Reference Standard is analyzed near the mid point of the calibration curve. The acceptance criteria for all target analytes in the reference standards are +/- 50% of the true value.
- **Method Blank:** Analyzed prior to the analysis of field samples and every 30 samples.

Note: Analyte levels reported for the field-deployed AGI Universal Samplers that exceed trip and method blank levels, and/or the reporting limit, are more likely to have originated from on-site sources.

| | |
|----------------------------|-------------------|
| Media Sampled: | WATER |
| Chemist - sample analysis: | Kelly J Stringham |
| Chemist - data processor: | Kelly J Stringham |
| Chemist - data review: | Jasmine Smith |

Method deviations: None.

Please note that data file names ending with R are rerun samples using the second pair of sorbers, in which the original results were not reported. Data file names ending in D are duplicate analysis results for the second set of sorbers from the same sampler, and are reported.

Additional Report Information

- Comments
- Laboratory Sample Report
- Chain of Custody
- Installation and Retrieval Log
- Analytical Results and Key
- Concentration Calculation Method Summary
- Total Ion Chromatograms

Project Specific Comments

None.

Survey period ¹

Samplers were installed on April 7, 2015 and retrieved on April 21, 2015 for an exposure period of 14 days.

Tamper seal intact:

Yes

Date received:

4/22/2015 10:25 AM

By: Darlene Yellowdy

COC returned:

Yes

Comments:

None

¹ - Installation start to end of retrieval, as reported. See installation and retrieval log for individual deployment and retrieval dates and times (i.e., sampler exposure time).

General Comments

Analytical QA/QC

Laboratory instrumentation consists of gas chromatographs equipped with mass selective detectors, coupled with automated thermal desorption units. Sample preparation involves cutting the tip off the bottom of the AGI Universal Sampler, and transferring one or more "sorbents" to a thermal desorption tube for analysis. The insertion/retrieval cord prevents soil, water and other interferences from coming in contact with the adsorbent. No further sample preparation is required. Any replicate sorbents not consumed in the initial analysis will be discarded fifteen (15) days from the date of the laboratory report.

Data are archived and stored in a secure manner as per AGI's Quality Assurance program (SPG-SOP-0462).

Total petroleum hydrocarbons (TPH), gasoline-range petroleum hydrocarbons (GRPH), and/or diesel range petroleum hydrocarbons (DRPH), when reported, are calculated using the area under the peaks observed in m/z 55 and 57 selected ion chromatograms. Quantitation of the mass values was performed using the response factor for a specific alkane (present in the calibration standards). TPH values include the entire chromatogram and provide estimates for aliphatic hydrocarbon ranges of C4 to C20. GRPH and DRPH include only the relevant regions of the chromatograms and provide estimates for C4 to C10 and C10 to C20 aliphatic hydrocarbons, respectively.

Trip blanks were provided to document potential exposures that were not part of the signal of interest (e.g., impact during sampler shipment, installation and/or retrieval, and storage). The trip blanks are identically manufactured and packaged AGI Universal Samplers to those samplers deployed in the field. The trip blanks remain unopened during all phases of the project. Levels reported on the trip blanks may indicate potential impact to the samplers other than the contaminant source of interest.

Unresolved peak envelopes (UPEs) are represented as a series of compound peaks clustered together around a central gas chromatograph elution time in the total ion chromatogram. UPEs may be indicative of complex fluid mixtures. UPEs observed early in the chromatograms are considered to indicate presence of more volatile fluids, while UPEs observed later in the chromatogram may indicate the presence of less volatile fluids. Multiple UPEs may indicate the presence of multiple complex fluids.

Total ion chromatograms (TICs) are included in the Attachments. The eight-digit serial number of each sampler is incorporated in the TIC identification (e.g., 12345678.D represents AGI Universal Sampler 12345678).

General Comments

Soil Gas Sampling

For soil gas sampling, the AGI Environmental Survey reports mass levels migrating through the open pore spaces of the soil and diffusing through the sampler membrane for sorption by the engineered, hydrophobic adsorbents, housed within the membrane tube. During the migration of the soil gas away from the source to the AGI Universal Sampler, the vapors are subject to a variety of attenuation factors. The soil gas masses reported on the samplers compare favorably with the concentrations reported in the soil or groundwater (e.g., where soil gas levels are reported at greater levels to other sampled locations on the site, the matrix data should reveal the same pattern, and vice versa). However, due to a variety of factors, a perfect comparison between matrix data and soil gas levels can rarely be achieved.

Soil gas concentrations ($\mu\text{g}/\text{m}^3$) are calculated following the method described in the Additional Report Information section.

Soil gas signals reported by this method cannot be correlated specifically to soil adsorbed, groundwater, and /or free-phase contamination. The soil gas signal reported from each AGI Universal Sampler can evolve from all of these sources. Differentiation between soil and groundwater contamination can only be achieved with prior knowledge of the site history (i.e., the site is known to have groundwater contamination only).

Air Sampling

For indoor, outdoor, and crawlspace air sampling, the AGI Environmental Survey reports mass levels present in the air and diffusing through the sampler membrane for sorption by the engineered adsorbents housed within the membrane tube.

Air concentrations ($\mu\text{g}/\text{m}^3$) are calculated following the method described in the Additional Report Information section.

Groundwater and Sediment Porewater Sampling

For groundwater and sediment porewater sampling, the AGI Environmental Survey reports the mass levels of compounds present in the water which, when coming in contact with the sampler membrane, partitions out of solution, and diffuses through the sampler membrane for sorption by the engineered adsorbents.

Water concentrations ($\mu\text{g}/\text{L}$) are calculated using the quantified mass, exposure period and the compound specific uptake rate. The rates were measured under controlled experimental conditions. The uptake rates are corrected for water pressure (depth of the AGI Universal Sampler below the water table), water temperature and the aquifer flow rate. For sediment porewater, the uptake rate is corrected for the reduced volume of water in the sediment, by multiplying the uptake rate by the pore water fraction.

LABORATORY SAMPLE REPORT

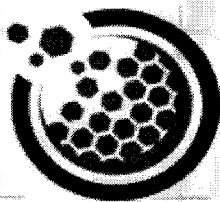
Project: ENV 01347

Site Name: Comal and San Marcos Rivers

Module Type: SPG0008

| Module ID | Sample Type | Field ID | |
|----------------------------|--------------------------|---------------------|-------------------|
| 00757309 | FIELD_SAMPLE | HCS410 | |
| 00757310 | FIELD_SAMPLE | HCS420 | |
| 00757311 | FIELD_SAMPLE | HCS430 | |
| 00757312 | FIELD_SAMPLE | HCS440 | |
| 00757313 | FIELD_SAMPLE | FDHCS440 | |
| 00757314 | LOST | HCS460 | |
| 00757315 | FIELD_SAMPLE | HSM410 | |
| 00757316 | FIELD_SAMPLE | HSM420 | |
| 00757317 | FIELD_SAMPLE | HSM430 | |
| 00757318 | FIELD_SAMPLE | FDHSM430 | |
| 00757319 | FIELD_SAMPLE | HSM440 | |
| 00757320 | FIELD_SAMPLE | HSM450 | |
| 00757321 | FIELD_SAMPLE | HMS460 | |
| 00757322 | FIELD_SAMPLE | HSM470 | |
| 00757323 | TRIP_BLANK | TB05 | |
| 00757324 | FIELD_SAMPLE | Not Provided | |
| Total # "FIELD SAMPLES" | Total # "TRIP BLANKS" | Total # "UNUSED" | Total # "LOST" |
| 14 | 1 | 0 | 1 |

Duplicate samples: 0



AMPLIFIED GEOCHEMICAL IMAGING, LLC

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ph: +1-302-266-2428
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AGI Universal Passive Sampler Chain of Groundwater Sampling

Production Order #: 01347

Customer Name: SWCA Environmental Consultants
Address: 6200 UTSA Boulevard Suite 102

Site Name: Comal and San Marcos Rivers
Site Address:

San Antonio, TX 78249
USA

Project Manager:

Serial # of Samplers Shipped
00757309 - 00757324

| | | | |
|--------------------------------|-----------|------------------|---|
| # of Samplers for Installation | 14.00 | # of Trip Blanks | 2 |
| Total Samplers Shipped | 16.00 | Pieces | |
| Total Samplers Received | <u>16</u> | Pieces | |
| Total Samplers Installed | <u>14</u> | Pieces | |

00757314 lost in field,
Kept vial in case
we find it in the
next day

Serial # of Trip Blanks (Client Decides)

007573243
gr

| | |
|--|--|
| Prepared By: <u>Lisa Bozelle</u> | Is Concurrent water sampling planned this monitoring period? YES <input type="radio"/> NO <input checked="" type="radio"/> |
| Verified By: <u>[Signature]</u> | Scheduled Sampling Date: _____ |
| Installation Performed By: Name: <u>Jennifer Moreland Brittany Rios</u> Company: <u>SWCA</u> | Retrieval Performed By: Name: <u>Jennifer Moreland Brittany Rios</u> Company: <u>SWCA</u> |
| Installation Start Date / Time: <u>4/7/15 9:41</u> | Retrieval Start Date / Time: <u>4/21/15 11:10</u> |
| Installation Complete Date / Time: <u>4/7/15 1302</u> | Retrieval Complete Date / Time: <u>4/21/15 1523</u> |
| Total Samplers Retrieved: <u>13</u> | |
| Total Samplers Lost In Field: <u>1</u> | |
| Total Unused Samplers Returned: <u>1 gr</u> | |
| Relinquished By: <u>Lisa Bozelle</u> Date/Time: <u>1:20 PM 3-24-15</u> Company: <u>AGI</u> | Received By: <u>[Signature]</u> Date/Time: <u>3/26/15 9:00AM</u> Company: <u>SWCA</u> |
| Relinquished By: <u>[Signature]</u> Date/Time: <u>4/21/15 10:00</u> Company: <u>SWCA</u> | Received By: <u>Shelene Galloway</u> Date/Time: <u>4/22/15 10:25</u> Company: <u>AGI</u> |



**AMPLIFIED
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IMAGING, LLC**

210 Executive Drive, Suite 1
Newark, DE USA 19702-3335
ph: 302-266-2428

AGI Project No.

ENV 01077

Site Name:

Comal & San Marco River

Site Location:

AGI Soil Gas Sampling

Installation & Retrieval Log

Company Name:

SWCA Environmental

Location:

Samples collected by:

Jennifer Moreland and Brittany Rios

* Optional or as needed

| | | | | | WATER QUALITY MONITORING | | |
|--------------------|----------|---|--|---|--|--|---|
| SAMPLER SERIAL NO. | WELL ID | SAMPLE TYPE (Field Sample, Trip Blank, Field Blank, etc.) | INSTALLATION DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 13:00 | RETRIEVAL DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 15:00 | INSTALLATION DEPTH FROM TOP OF CASING (feet) | DEPTH TO WATER FROM TOP OF CASING (feet) | SCREENED INTERVAL FROM TOP OF CASING (feet) |
| 00757309 | HCS410 | FIELD_SAMPLE | 4/7/15 10:00 | 4/21/15 11:34 | 1.5 | | |
| 00757310 | HCS420 | FIELD_SAMPLE | 4/7/15 10:11 | 4/21/15 11:50 | 3 | | |
| 00757311 | HCS430 | FIELD_SAMPLE | 4/7/15 9:41 | 4/21/15 11:10 | 7 | | |
| 00757312 | HCS440 | FIELD_SAMPLE | 4/7/15 10:23 | 4/21/15 12:05 | 4 | | |
| 00757313 | FDHCS440 | FIELD_DUPLICATE | 4/7/15 10:23 | 4/21/15 12:05 | 4 | | |
| 00757314 | HCS460 | FIELD_SAMPLE | 4/7/15 10:33 | 4/21/2015 N/A | 5 | | |
| 00757315 | HSM410 | FIELD_SAMPLE | 4/7/15 11:14 | 4/21/15 13:51 | 2 | | |
| 00757316 | HSM420 | FIELD_SAMPLE | 4/7/15 11:26 | 4/21/15 14:02 | 4 | | |
| 00757317 | HSM430 | FIELD_SAMPLE | 4/7/15 11:36 | 4/21/15 14:21 | 0.5 | | |
| 00757318 | FDHSM430 | FIELD_DUPLICATE | 4/7/15 11:36 | 4/21/15 14:21 | 0.5 | | |
| 00757319 | HSM440 | FIELD_SAMPLE | 4/7/15 11:49 | 4/21/15 14:35 | 5 | | |
| 00757320 | HSM450 | FIELD_SAMPLE | 4/7/15 12:31 | 4/21/15 14:52 | 4 | | |
| 00757321 | HMS460 | FIELD_SAMPLE | 4/7/15 12:50 | 4/21/15 15:11 | 3 | | |
| 00757322 | HSM470 | FIELD_SAMPLE | 4/7/15 13:02 | 4/21/15 15:23 | 1.5 | | |
| 00757323 | TB05 | TRIP_BLANK | 4/7/15 9:41 | 4/21/15 15:23 | NA | | |



**AGI Soil Gas Sampling
Installation & Retrieval L**

* Optional or as needed

| MONITORING | | | OBSERVATIONS/COMMENTS* (e.g., observations, location description, missing, pulled from well, etc. - as needed) | YES / NO | |
|--------------------|--|---|--|--|--------|
| SAMPLER SERIAL NO. | WATER TEMPERATURE AT MODULE DEPTH (deg C) | FLOW RATE THROUGH SCREEN: HIGH OR LOW (HIGH = flow > than 10 m/day; LOW = flow < 10 m/day) | | EVIDENCE OF LIQUID PETROLEUM HYDROCARBONS? | ODOR ? |
| 00757309 | 23.3 | High | | No | No |
| 00757310 | 23.3 | High | | No | No |
| 00757311 | 23.3 | High | | No | No |
| 00757312 | 23.3 | High | | No | No |
| 00757313 | 23.3 | High | | No | No |
| 00757314 | 23.3 | High | Sampler was missing upon date of retrieval | No | No |
| 00757315 | 22.0 | High | | No | No |
| 00757316 | 22.0 | High | | No | No |
| 00757317 | 22.0 | High | | No | No |
| 00757318 | 22.0 | High | | No | No |
| 00757319 | 22.0 | High | | No | No |
| 00757320 | 22.0 | High | | No | No |
| 00757321 | 22.0 | High | | No | No |
| 00757322 | 22.0 | High | | No | No |
| 00757323 | | | | No | No |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757309 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS410

Installation Date: 4/7/2015 10:00:00AM

Retrieval Date: 4/21/2015 11:34:00AM

Date Analyzed: 5/6/2015 12:56:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757309 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS410

Installation Date: 4/7/2015 10:00:00AM

Retrieval Date: 4/21/2015 11:34:00AM

Date Analyzed: 5/6/2015 12:56:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|-------------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 1.82 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00757310 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS420

Installation Date: 4/7/2015 10:11:00AM

Retrieval Date: 4/21/2015 11:50:00AM

Date Analyzed: 5/6/2015 2:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.34 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | 0.16 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



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210 Executive Drive, Suite 1
Newark, DE 19702-3335 USA
ph: +1-302-266-2428
www.agisurveys.net

PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757310 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS420

Installation Date: 4/7/2015 10:11:00AM

Retrieval Date: 4/21/2015 11:50:00AM

Date Analyzed: 5/6/2015 2:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|-------------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 1.02 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757311 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS430

Installation Date: 4/7/2015 9:41:00AM

Retrieval Date: 4/21/2015 11:10:00AM

Date Analyzed: 5/6/2015 12:25:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.42 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757311 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS430

Installation Date: 4/7/2015 9:41:00AM

Retrieval Date: 4/21/2015 11:10:00AM

Date Analyzed: 5/6/2015 12:25:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|-------------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 1.26 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757312 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS440

Installation Date: 4/7/2015 10:23:00AM

Retrieval Date: 4/21/2015 12:05:00PM

Date Analyzed: 5/6/2015 2:56:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.41 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757312 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS440

Installation Date: 4/7/2015 10:23:00AM

Retrieval Date: 4/21/2015 12:05:00PM

Date Analyzed: 5/6/2015 2:56:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|-------------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.65 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757313 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: FDHCS440

Installation Date: 4/7/2015 10:23:00AM

Retrieval Date: 4/21/2015 12:05:00PM

Date Analyzed: 5/6/2015 7:56:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.37 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757313 FIELD_SAMPLE
Dilution Factor: 1 Field ID: FDHCS440
Installation Date: 4/7/2015 10:23:00AM
Retrieval Date: 4/21/2015 12:05:00PM
Analyst: Kelly J Stringham
Reviewer: Jasmine Smith

Matrix: WATER Product: SPG0008
Date Analyzed: 5/6/2015 7:56:00PM
Method: SPG-WI-0292
Batch: ENV-150424-1

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|-------------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 1.05 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757315 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM410

Installation Date: 4/7/2015 11:14:00AM

Retrieval Date: 4/21/2015 1:51:00PM

Date Analyzed: 5/6/2015 1:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757315 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM410

Installation Date: 4/7/2015 11:14:00AM

Retrieval Date: 4/21/2015 1:51:00PM

Date Analyzed: 5/6/2015 1:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|-------------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.93 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757316 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM420

Installation Date: 4/7/2015 11:26:00AM

Retrieval Date: 4/21/2015 2:02:00PM

Date Analyzed: 5/6/2015 3:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|--------------------------|-------------|-------------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | 0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.10 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | 0.03 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | 0.09 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757316 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM420

Installation Date: 4/7/2015 11:26:00AM

Retrieval Date: 4/21/2015 2:02:00PM

Date Analyzed: 5/6/2015 3:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|-------------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.54 | 0.50 |
| BTEX | | 0.14 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757317 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM430

Installation Date: 4/7/2015 11:36:00AM

Retrieval Date: 4/21/2015 2:21:00PM

Date Analyzed: 5/6/2015 4:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.80 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757317 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM430

Installation Date: 4/7/2015 11:36:00AM

Retrieval Date: 4/21/2015 2:21:00PM

Date Analyzed: 5/6/2015 4:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|---------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757318 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: FDHSM430

Installation Date: 4/7/2015 11:36:00AM

Retrieval Date: 4/21/2015 2:21:00PM

Date Analyzed: 5/6/2015 6:56:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.72 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757318 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: FDHSM430

Installation Date: 4/7/2015 11:36:00AM

Retrieval Date: 4/21/2015 2:21:00PM

Date Analyzed: 5/6/2015 6:56:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|-------------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.54 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757319 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM440

Installation Date: 4/7/2015 11:49:00AM

Retrieval Date: 4/21/2015 2:35:00PM

Date Analyzed: 5/6/2015 11:55:00AM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.13 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757319 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM440

Installation Date: 4/7/2015 11:49:00AM

Retrieval Date: 4/21/2015 2:35:00PM

Date Analyzed: 5/6/2015 11:55:00AM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|---------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757320 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM450

Installation Date: 4/7/2015 12:31:00PM

Retrieval Date: 4/21/2015 2:52:00PM

Date Analyzed: 5/6/2015 3:56:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.08 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



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Newark, DE 19702-3335 USA
ph: +1-302-266-2428
www.agisurveys.net

PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757320 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM450

Installation Date: 4/7/2015 12:31:00PM

Retrieval Date: 4/21/2015 2:52:00PM

Date Analyzed: 5/6/2015 3:56:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|---------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757321 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HMS460

Installation Date: 4/7/2015 12:50:00PM

Retrieval Date: 4/21/2015 3:11:00PM

Date Analyzed: 5/6/2015 6:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.11 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757321 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HMS460

Installation Date: 4/7/2015 12:50:00PM

Retrieval Date: 4/21/2015 3:11:00PM

Date Analyzed: 5/6/2015 6:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|---------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757322 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM470

Installation Date: 4/7/2015 1:02:00PM

Retrieval Date: 4/21/2015 3:23:00PM

Date Analyzed: 5/6/2015 8:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.09 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757322 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM470

Installation Date: 4/7/2015 1:02:00PM

Retrieval Date: 4/21/2015 3:23:00PM

Date Analyzed: 5/6/2015 8:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|-------------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.68 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757323 TRIP BLANK

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: TB05

Installation Date: 4/7/2015 9:41:00AM

Retrieval Date: 4/21/2015 3:23:00PM

Date Analyzed: 5/6/2015 7:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757323 TRIP BLANK

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: TB05

Installation Date: 4/7/2015 9:41:00AM

Retrieval Date: 4/21/2015 3:23:00PM

Date Analyzed: 5/6/2015 7:26:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150424-1

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|---------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00757324 FIELD_SAMPLE
Dilution Factor: 1

Matrix: WATER

Product: SPG0008

Date Analyzed: 5/6/2015 1:56:00PM

Analyst: Kelly J Stringham
Reviewer: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-150424-1

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01347
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00757324 FIELD_SAMPLE
Dilution Factor: 1

Matrix: WATER

Product: SPG0008

Date Analyzed: 5/6/2015 1:56:00PM

Analyst: Kelly J Stringham
Reviewer: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-150424-1

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------|------------|-------------|---------|
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |

AMPLIFIED GEOCHEMICAL IMAGING, LLC
210 EXECUTIVE DRIVE, SUITE 1, NEWARK, DE 19702
SWCA ENVIRONMENTAL CONSULTANTS, SAN ANTONIO, TX
STANDARD TARGET VOCs/SVOCs
ESTIMATED WATER CONCENTRATIONS
COMAL AND SAN MARCOS RIVERS
ORDER # 01347

| | | | | | | | | | estimated | | |
|---------------|--------------|------------|------------|------------|----|------------|----|----|-----------|------------|--------------|
| DATAFILE | FIELD | DATE/ TIME | DATE/ TIME | DATE/ TIME | | DATE/ TIME | | | | | |
| NAME | ID | INSTALLED | RETRIEVED | RECEIVED | | ANALYZED | | DF | TPH, ug/L | MTBE, ug/L | t12DCE, ug/L |
| RL = | | | | | | | | | 0.05 | 0.01 | 0.01 |
| 00757309 | HCS410 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | 0.09 | <0.01 | <0.01 |
| 00757310 | HCS420 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | 0.07 | <0.01 | <0.01 |
| 00757311 | HCS430 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | 0.08 | <0.01 | <0.01 |
| 00757312 | HCS440 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | 0.06 | <0.01 | <0.01 |
| 00757313 | FDHCS440 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | 0.07 | <0.01 | <0.01 |
| 00757315 | HSM410 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | 0.07 | <0.01 | <0.01 |
| 00757316 | HSM420 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | 0.06 | <0.01 | <0.01 |
| 00757317 | HSM430 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |
| 00757318 | FDHSM430 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | 0.06 | <0.01 | <0.01 |
| 00757319 | HSM440 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |
| 00757320 | HSM450 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |
| 00757321 | HMS460 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |
| 00757322 | HSM470 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | 0.06 | <0.01 | <0.01 |
| 00757323 | TB05 | 4/7/2015 | 4/21/2015 | 4/22/2015 | ET | 5/6/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |
| 00757324 | | | | 4/22/2015 | ET | 5/6/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |
| BLK_ENV1_SPG8 | Method Blank | | | | | 5/6/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |

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ESTIMATED WATER CONCENTRATIONS
COMAL AND SAN MARCOS RIVERS
ORDER # 01347

| DATAFILE | | | | | | | | | |
|---------------|-------------|--------------|-------------|--------------|-------------|------------|------------|-----------|--------------|
| NAME | 11DCA, ug/L | c12DCE, ug/L | CHCl3, ug/L | 111TCA, ug/L | 12DCA, ug/L | BENZ, ug/L | CCl4, ug/L | TCE, ug/L | 112TCA, ug/L |
| RL = | 0.01 | 0.01 | 0.01 | 0.00 | 0.01 | 0.01 | 0.00 | 0.01 | 0.01 |
| 00757309 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757310 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757311 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757312 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757313 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757315 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757316 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757317 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757318 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757319 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757320 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757321 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757322 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757323 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| 00757324 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |
| BLK_ENV1_SPG8 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 |

AMPLIFIED GEOCHEMICAL IMAGING, LLC
 210 EXECUTIVE DRIVE, SUITE 1, NEWARK, DE 19702
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 STANDARD TARGET VOCs/SVOCs
 ESTIMATED WATER CONCENTRATIONS
 COMAL AND SAN MARCOS RIVERS
 ORDER # 01347

| DATAFILE | | | | | | | | | |
|---------------|-----------|-----------|-----------|--------------|-----------------|--------------|-------------|------------|-----------------|
| NAME | TOL, ug/L | OCT, ug/L | PCE, ug/L | CIBENZ, ug/L | 1112TetCA, ug/L | ETBENZ, ug/L | mpXYL, ug/L | oXYL, ug/L | 1122TetCA, ug/L |
| RL = | 0.01 | 0.00 | 0.01 | 0.01 | 0.01 | 0.01 | 0.01 | 0.01 | 0.01 |
| 00757309 | <0.01 | <0.00 | 0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757310 | <0.01 | <0.00 | 0.06 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757311 | <0.01 | <0.00 | 0.07 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757312 | <0.01 | <0.00 | 0.07 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757313 | <0.01 | <0.00 | 0.07 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757315 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757316 | 0.01 | <0.00 | 0.02 | <0.01 | <0.01 | 0.01 | 0.02 | <0.01 | <0.01 |
| 00757317 | <0.01 | <0.00 | 0.14 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757318 | <0.01 | <0.00 | 0.13 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757319 | <0.01 | <0.00 | 0.03 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757320 | <0.01 | <0.00 | 0.02 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757321 | <0.01 | <0.00 | 0.02 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757322 | <0.01 | <0.00 | 0.02 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757323 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00757324 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| BLK_ENV1_SPG8 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |

AMPLIFIED GEOCHEMICAL IMAGING, LLC
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STANDARD TARGET VOCs/SVOCs
ESTIMATED WATER CONCENTRATIONS
COMAL AND SAN MARCOS RIVERS
ORDER # 01347

| | | | | | | | | estimated |
|---------------|--------------|--------------|-------------|-------------|-------------|-------------|------------|--------------|
| DATAFILE | | | | | | | | |
| NAME | 135TMB, ug/L | 124TMB, ug/L | 13DCB, ug/L | 14DCB, ug/L | 12DCB, ug/L | UNDEC, ug/L | NAPH, ug/L | TRIDEC, ug/L |
| RL = | 0.00 | 0.01 | 0.01 | 0.01 | 0.01 | 0.02 | 0.02 | 0.02 |
| 00757309 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757310 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | 0.03 | <0.02 | <0.02 |
| 00757311 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757312 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757313 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757315 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757316 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757317 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757318 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757319 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757320 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757321 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757322 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757323 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| 00757324 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |
| | | | | | | | | |
| BLK_ENV1_SPG8 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.02 | <0.02 | <0.02 |

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 STANDARD TARGET VOCs/SVOCs
 ESTIMATED WATER CONCENTRATIONS
 COMAL AND SAN MARCOS RIVERS
 ORDER # 01347

| | | estimated | estimated | estimated | estimated |
|---------------|---------------|----------------------|----------------|--------------------|----------------|
| DATAFILE | | | | | |
| NAME | 2MeNAPH, ug/L | Acenaphthylene, ug/L | PENTADEC, ug/L | Acenaphthene, ug/L | Fluorene, ug/L |
| RL = | 0.01 | 0.01 | 0.02 | 0.01 | 0.01 |
| 00757309 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757310 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757311 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757312 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757313 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757315 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757316 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757317 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757318 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757319 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757320 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757321 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757322 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757323 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| 00757324 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |
| | | | | | |
| BLK_ENV1_SPG8 | <0.01 | <0.01 | <0.02 | <0.01 | <0.01 |

KEY TO DATA TABLE

UNITS

| | |
|-------------------|--|
| µg | micrograms, relative mass value |
| µg/m ³ | micrograms per cubic meter; estimated soil gas concentration |
| µg/L | micrograms per Liter; calculated water concentration |

DATA QUALIFIERS

| | |
|---|--|
| > | greater than; value exceeds calibration range, estimated value |
| < | less than; compound value is below the LOD and RL |
| J | mass value below LOQ or RL, but above LOD, estimated mass value |
| E | mass value exceeds upper calibration level, estimated mass value |
| Q | one or more quality control parameters failed for the compound |

ABBREVIATIONS

| | | | |
|-------------|---|----------|--|
| AVG RL | average reporting limit; calculated based on individual field sample RLs | | |
| LOD | limit of detection | | |
| LOQ | limit of quantification | | |
| MDL | method detection limit | | |
| RL | reporting limit | | |
| 1112TetCA | 1,1,1,2-tetrachloroethane | CIBENZ | chlorobenzene |
| 111TCA | 1,1,1-trichloroethane | ct12DCE | cis- & trans-1,2-dichloroethene |
| 1122TetCA | 1,1,2,2-tetrachloroethane | EtBENZ | ethylbenzene |
| 112TCA | 1,1,2-trichloroethane | mpXYL | m-, p-xylene |
| 11DCA | 1,1-dichloroethane | MTBE | methyl t-butyl ether |
| 11DCE | 1,1-dichloroethene | NAPH | naphthalene |
| 124TMB | 1,2,4-trimethylbenzene | OCT | octane |
| 12DCA | 1,2-dichloroethane | oXYL | o-xylene |
| 12DCB | 1,2-dichlorobenzene | PCE | tetrachloroethene |
| 135TMB | 1,3,5-trimethylbenzene | PENTADEC | pentadecane |
| 13DCB | 1,3-dichlorobenzene | PHEN | phenanthrene |
| 14DCB | 1,4-dichlorobenzene | t12DCE | trans-1,2-dichloroethene |
| 2MeNAPH | 2-methyl naphthalene | TCE | trichloroethene |
| BENZ | benzene | TMBs | combined masses of 1,3,5-trimethylbenzene and 1,2,4-trimethylbenzene |
| BTEX | combined masses of benzene, toluene, ethylbenzene, and total xylenes (Gasoline Range Aromatics) | TOL | toluene |
| C11,C13&C15 | combined masses of undecane, tridecane, and pentadecane (C11+C13+C15) (Diesel Range Alkanes) | TPH | total petroleum hydrocarbons |
| c12DCE | cis-1,2-dichloroethene | TRIDEC | tridecane |
| CCl4 | carbon tetrachloride | UNDEC | undecane |
| CHC13 | chloroform | VC | vinyl chloride |

SUMMARY OF SAMPLING RATE CALIBRATION FOR AGI UNIVERSAL SAMPLER IN AQUEOUS PHASES

PURPOSE:

The purpose of this document is to:

1. Summarize the test protocol,
2. Summarize the methodology for analysis of data,
3. Present general results for generating concentration calibration

of the AGI Universal Sampler, part number SPG-0008, in aqueous phase media following AGI's "Standard Practice for Determining the Sampling Rate of Passive Diffusion Samplers in Various Environmental Media": SPG-SOP-0493. The work will be summarized in three parts: Part 1: shallow water, Part 2: deep water, and Part 3: sediment.

Principle of Operation of the AGI Samper

The AGI Universal Sampler is designed with solid adsorbents enclosed inside a tubular microporous PTFE membrane. When placed in water, the pores and hydrophobic nature of the PTFE keep liquid water from entering the membrane until a water head of about 34 feet is reached. The membrane will not keep water vapor from entering but the adsorbents are very hydrophobic and through testing validated to be unaffected by this moisture vapor. In shallow water, <34', volatile and semi-volatile compounds will partition from the dissolved water into the air phase in the PTFE membrane according to Henry's Law. This partitioning is instantaneous and within seconds-minutes, the compound is adsorbed by the adsorbent inside the sealed tube. Because the diffusivity in air is about 10,000 times higher than the diffusivity in water, the sampling rate is controlled by the water contact area of the membrane that allows the Henry's Law effect to occur. This contact area is set by the membrane diameter and length of the sealed tube, which is fixed in AGI's manufacturing process.

Henry's law as well as diffusivity, which are fundamentally incorporated into the sampling rate, are affected by temperature, T , and follow an Arrhenius equation $H_T = H_r \times \exp[-E_a/R](1/T_r - 1/T)$. Because a 5°C temperature change can make a 15% change in sampling rate, the temperature of the sampled water should be known to get the most precise concentration.

The membrane pore size is also small enough that colloidal particles and microbes can not pass through the membrane. This keeps the adsorbent from getting contaminated and eliminates any need to add preservative or chilling during storage or transportation.

When the water pressure exceeds the water entry pressure of the membrane, about 34 feet of water, the water becomes in contact with the solid adsorbent. Under this condition the compounds in the water will partition from the water into the solid. The partitioning coefficient, K_{AW} , can be approximated by the octanol-water coefficient, K_{OW} , but has been measured more precisely in the lab for AGI's specific solid adsorbent. The sampling rate is the product of the sampling rate at <34' of water and the K_{AW} .

In sediment, the sampler measures pore-water concentration, which is generally agreed to be the preferred measurement as it is more indicative of bioavailability. In sediment the volumetric availability of water to the sampler is reduced by the volume fraction solids in the sediment, which typically varies from zero to 35%, but can be as high as 73% in well packed and broad particle size distribution sediments. As a result, sampling rates in sediment are multiplied by the fraction pore water in the sediment to determine concentration.

PART 1: Calibration in shallow water

Part 1 summarizes the work in shallow water generating calibration data, evaluating the physical and chemical factors affecting the sampling rate, and measurement of the actual sampling rates or regression calibration equations needed to determine concentrations.

Sample Generation in water

In this calibration work, solutions of analytes at known concentrations were formulated in clean 4 liter smoked glass jugs by injecting microliter measured amounts of environmental standards using a calibrated syringe into pure or deionized water and stirring for a minimum of 2 hours but generally overnight. Headspace in the jugs was minimized and generally less than 1% by volume during the tests. Jugs were temperature controlled by placing them in a water filled cooler, chilled via a copper tubing loop in the cooler. Temperature was measured with a certified digital temperature gauge and an average value used for each temperature experiment.

AGI samplers were weighted so they won't float and placed in the jugs at time zero. They were removed at various intervals to generate samples along with duplicates that showed mass increasing with exposure time. The sampler exposure time was selected to span minutes to hours and was generally reduced for high concentration tests to maintain uptake with time in roughly the linear dynamic range. Samplers were removed and dried with a paper towel and returned to their original container for analysis. They were analyzed by AGI's 8260C (SPG-WI-318 or SPG-WI-10028) method in duplicate, which is based on EPA SW846 Method 8260C.

Water samples were also taken and measured at an outside accredited lab using EPA SW846 Method 8260B. The concentrations agreed well with the calculated concentrations based on the standard certification, jug volume, and syringe injection. The variability of the outside lab 8260B values were found to be high, so for the sampling rate calculations we used the concentrations based on syringe dosing.

Calibrations were run at five concentrations, nominally at 6, 24, 118, 590, 1420 ug/L and five temperatures nominally at 5, 10, 15, 20, and 25 degrees centigrade. Samples were taken at 4 different exposure times. Samples were run in duplicate. A total of 176 data points were generated using 28 compounds from AGI's standard compounds list. Tridecane and pentadecane were not evaluated due to their very low solubility in water. In addition, another 23 compounds were tested using an 8260 liquid standard at nominal concentrations of 0.5, 1.0, 5.0, 15, 95, and 470 ug/L at a temperature typical of groundwater, 15°C. This is a living calibration and as additional data are generated, they may be qualified and added to this data set to improve the precision of the sampling rate calibration and broaden the compound list.

Key Variable Effects

As expected from theory, at short to moderate exposure times, mass will increase roughly linearly proportional to exposure time, as well as proportional to concentration, and exponentially with temperature following Arrhenius law. Temperature affects the Henry's law as well as diffusivity in water. Sampling rate is generally independent of concentration and time at mass values significantly below saturation. In the following sections we have characterized the sampling rate for each compound as affected by temperature and also developed calibrations using regression which account for the minor impact of time, and mass.

Concentration using Simple Sampling Rate Determination

A simple way to determine concentration is to measure mass on the AGI sampler, divide by exposure time, and divide by sampling rate, SR.

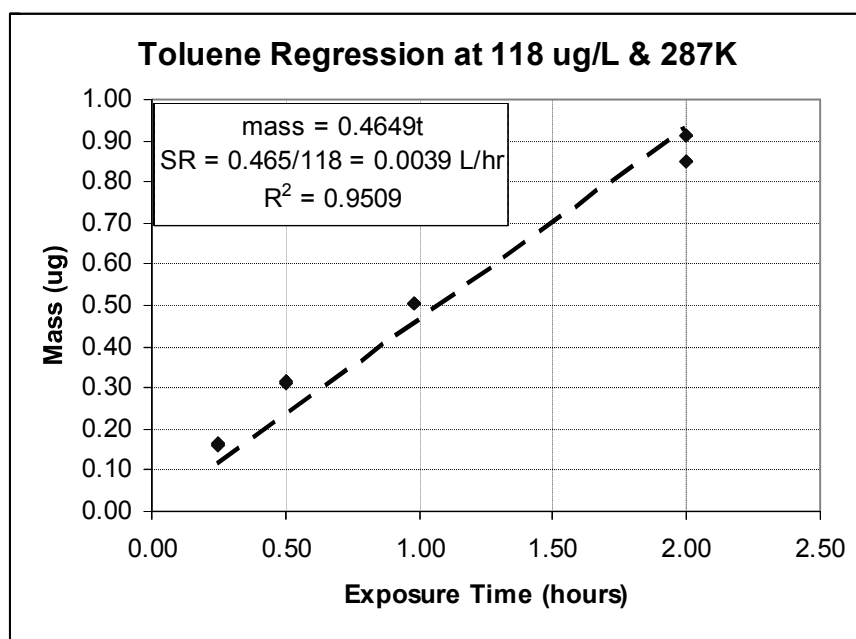
$$\text{Conc [ug/L]} = \text{mass/time/SR} \quad (1)$$

The sampling rate can be determined via measurements of mass versus time at a known concentration and temperature according to the following modification of equation (1).

$$\text{SR} = \text{mass/time/concentration} \quad (2)$$

Sampling rates in L/hr were determined by measuring the trend or regression mass uptake versus time and dividing by the concentration. A measurement like this will use 8 data points (4 times x 2 samples). Such a sampling rate can be measured at any concentration and temperature.

The chart to the right shows a plot of mass versus time for water at 118 ug/L and 287K. This is actual data from a single run. Slope of 0.465 ug/hr divided by the concentration of 118 ug/L yields a sampling rate, SR, of 0.0039 L/hr.



SR's typically range from about 0.004 to 0.007 L/hr at 15°C. Table A shows SR's measured for our standard compound list at 5 temperatures.

Rigorous Concentration using Regression

A preferred method for determining concentration that will yield improved accuracy over a wide range of concentrations, exposure times, and temperatures is to use all data in a regression analysis. This allows adjustments for the minor non-linear influences of mass and time as well as the effects of temperature. This is done by regressing equation (1) or a universal version of equation (1)

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRo} \cdot \exp(-E_a/R/T)] \quad (3)$$

The subtle non-linear effects of mass and time will be evident in the deviation of coefficients b and d from 1.0. This regression generates four constants b, d, SRo, and $-E_a/R$ by regressing $\ln(\text{conc})$ versus $\ln(\text{mass})$, $\ln(\text{time})$, $1/\text{temp}$. These four constants can be used to determine concentration via the equation:

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRo} \times \exp(-E_a/R(1/T))] \quad (4)$$

Where conc is in ug/L, mass is in ug, time in hours, T in degrees Kelvin.

Equation (4) can be also expressed at a reference temperature, T_r , such as 15°C by

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRr} \times \exp(-E_a/R(1/T_r - 1/T))] \quad (5)$$

This allows sampling rates, SRr, at any reference temperature, T_r , and for any analyte to easily be compared. These values of SRr at 293.14K can be found in Table A.

When sampling times are between 0 and 4 hours, using the 4 constant equation (5) is preferred. For concentrations from about 5 to 1500 ug/L one hour exposure times generally give the lowest error, typically with average error of 6-20% and with total error range of 12%-32%. For low concentrations where sampling times are greater than 4 hours, it is preferred to use equation (1) to avoid unrealistic effects from the coefficient d or to set d to 1.0. In such a case SR in equation (1) can be substituted with $[\text{SRr} \times \exp(-E_a/R(1/T_r - 1/T))]$ to use an SR representative of the well temperature, T.

The chart to the right is a plot of the calculated concentration from the 4 constant regression compared to the dosed concentration. Agreement is excellent for the 176 data points.

There does appear to be a slight high bias of 8.6% over the full range of this data, although it is well within acceptable limits of variability.

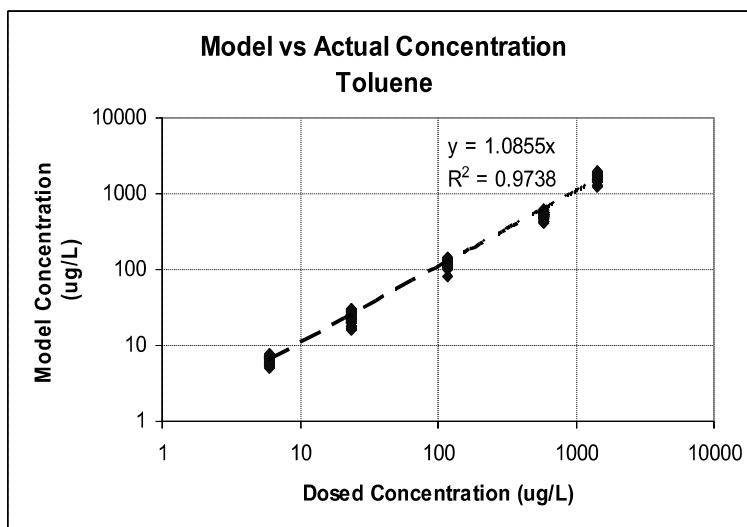


Table B shows the tabulated summary of the 4 constants regression with Rsq values and error estimates for the 4 constants for each analyte. Most regression Rsq values are 0.99 or greater for each analyte. In general, $-E_a/R$ is about 2400 \pm 400, b is about 0.9, d is about -0.75, and SR(15°C) ranges from .004 L/hr to 0.007 L/hr increasing with MW of the compound.

Error Estimates

The error in the water concentration values will depend on both the error in mass from the analytical method as well as the error in the concentration calibration. Table C shows the error in the mass values from the 8260C low sensitivity method.

The standard error of the regression and standard errors of the constants can be found in table B. For each compound we have measured the error between the derived concentration and the actual concentration. The error tends to be lowest at our recommended exposure time of one hour as shown by the example for Toluene to the right.

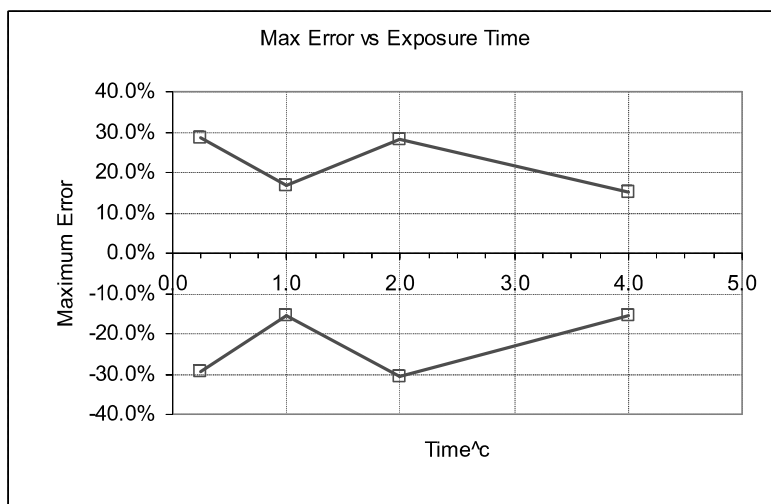
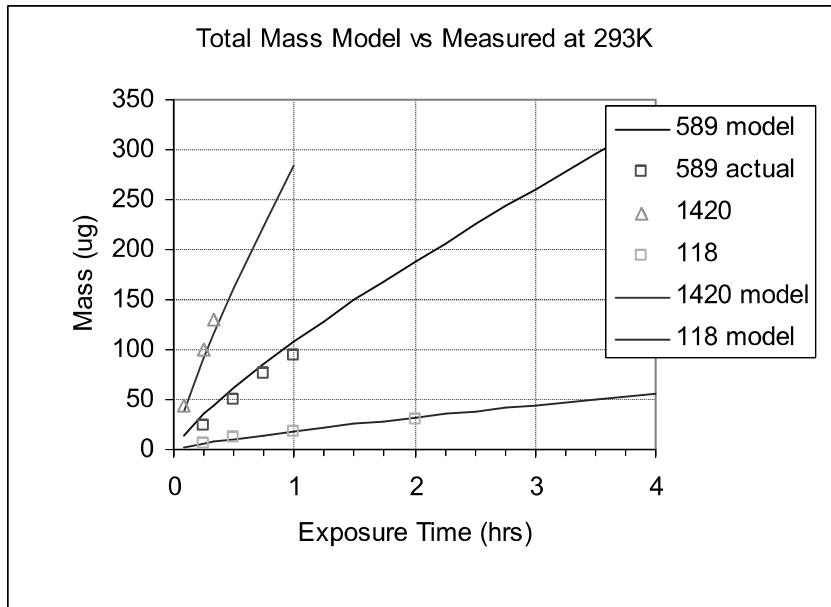


Table D shows the total average error in water concentration by compound as well as the low and high error. The average ranges from about 6% to 20%, which is similar to the analytical method errors. The low and high errors range from 12% to 32% and include contribution from measurement errors in both time and temperature.

Sorbent Saturation

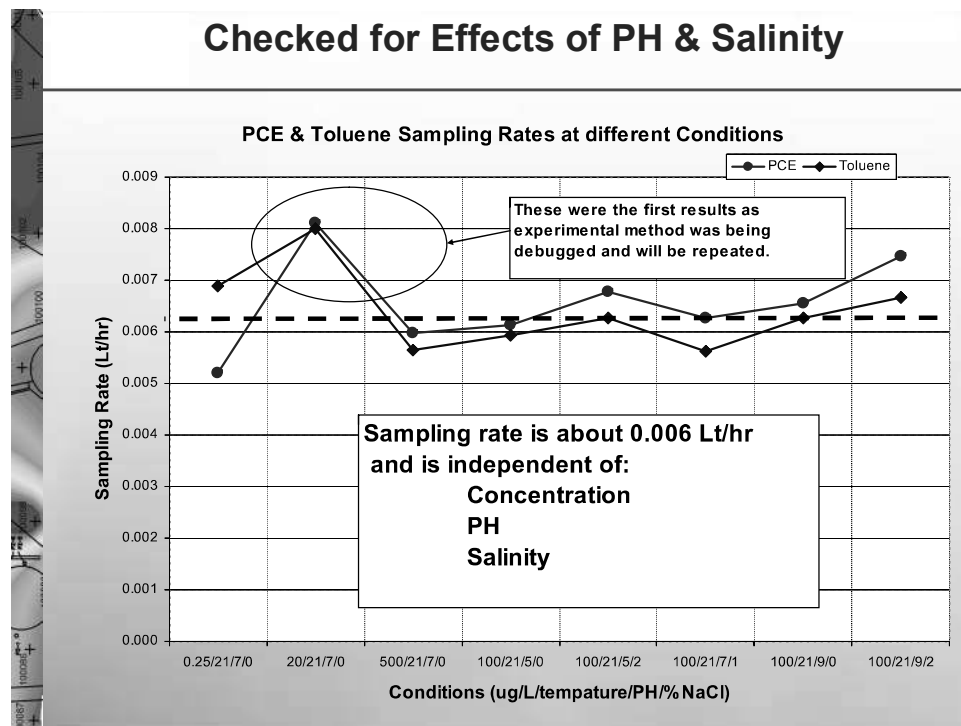
As mass increases on a solid sorbent and approaches saturation, reverse diffusion can occur causing the sampling rate to drop. Eventually the mass level will reach a maximum steady state value at any concentration. A rate of mass uptake with time that deviates significantly from linear, indicates that sorbent saturation could be an issue. When using equation (1), staying in the linear range to avoid the effects of adsorbent saturation is important. We recommend keeping the total mass on the sampler below 50 ug or flagging when this is exceeded.

The 4 constant regression accounts for some of the non linearity allowing good accuracy at higher mass levels. From the experimental data we have found this safe range can be extended to 100 ug or higher as shown in the chart below. This chart compares total mass of all compounds (excluding heavy alkanes, which have solubility issues) versus time in comparison to that predicted from the 4-constant concentration equation.



Effect of PH and Salinity

Because neither PH nor salinity is known to have a significant impact on Henry's law or diffusivity in water, we did not expect them to have a significant impact on sampling rate. To confirm this, experiments were run varying PH from 5 to 9 and NaCl content from 0 to 2%. The chart below shows no significant impact for combinations of PH and NaCl content over this range on the sampling rate of toluene in water at 21°C.

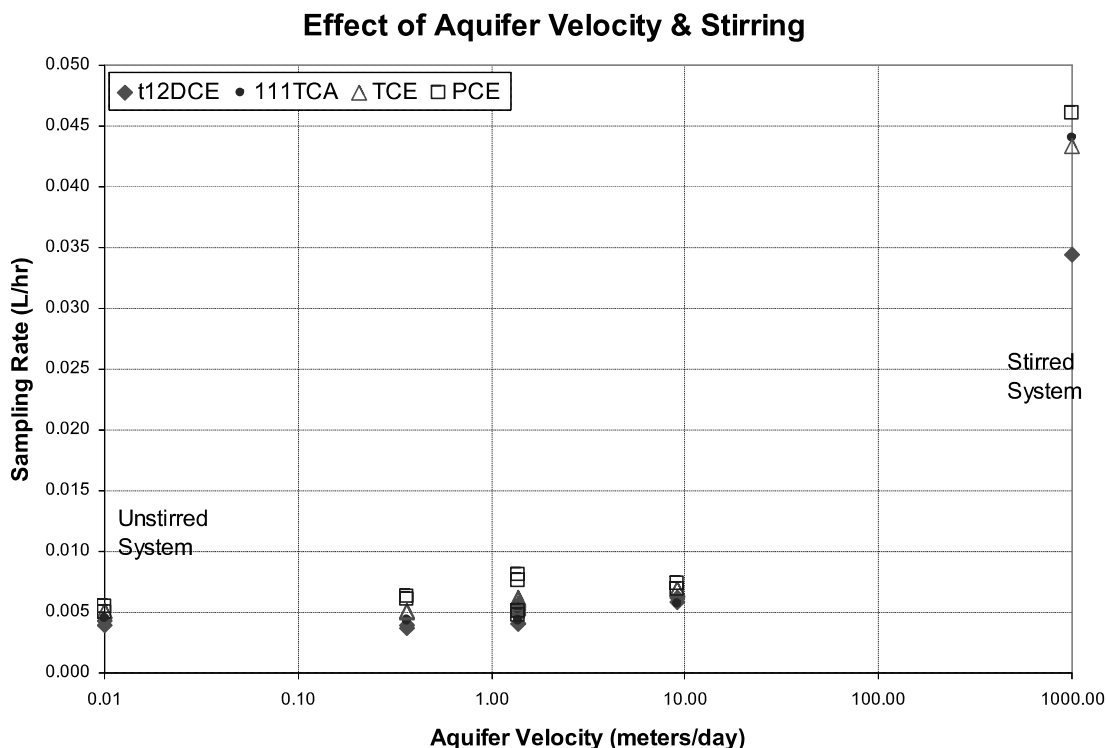


Impact of Aquifer Velocity

The velocity in most aquifers is quite slow, typically a meter/day or less. Occasionally water flow could be much higher such as encountered in karst aquifers, streams or rivers. Mass transfer coefficients are higher in high flow conditions, which will lead to higher sampling rates. We validated that a highly stirred system had sampling rates about 10 times higher than those that were non-stirred. We decided to evaluate the effect of aquifer velocity.

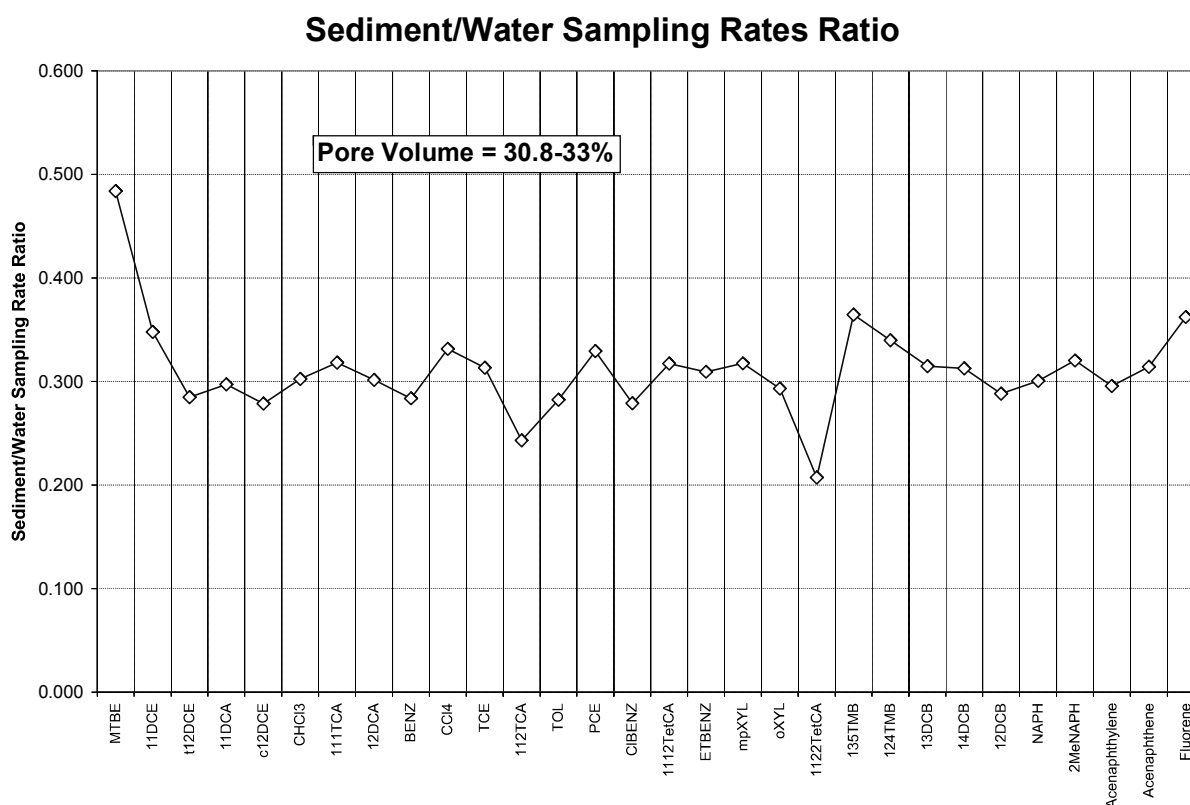
A test apparatus was built comprising a 3" PVC pipe tee filled with clean sand in each of the horizontal straight legs and screened to leave the center open. A test solution was run through this system using a variable flow pump and AGI samplers were placed into the simulated well through the vertical leg of the tee. Tests were run to examine the effect of velocity by varying the pumping rate and hence water velocity.

The chart below shows no significant effect of aquifer velocity up to a speed of about 10 meters/day. At velocities significantly above this, similar to a stirred system, sampling rates are about 10 times higher.



concentration. Pore water fraction can range from 1.0 for water without sediment to as low as 0.25. Typically most sediments have pore fractions of 0.9 to 0.65.

A sampling rate study was done with water and with water added into a well-packed sorted sand. Pore water fraction in this test was measured between 30.8% 33% by volume. Below is a plot of the ratio of sampling rates measured in the sediment to open water. The average ratio is equal to the pore water fraction confirming that sampling rate in sediment is on average equal to the product of pore water fraction times the sampling rate in water.



Summary

The AGI® Sampler can be used to determine the concentration of volatile and semi-volatile compounds in a water phase. This requires knowing the exposure time and water temperature. It also requires knowing if the sample is above or below 34' of water head and if the water has a velocity above 10 meters/day. Regressions of large amounts of data were used to generate a four constant equation to generate concentration values in water. Potential error in the concentration values is excellent typically less than 25%.

TABLE A
WATER SAMPLING RATES STANDARD LIST

| | SRr 293.14 | SR @ 277.54 | SR @ 282.44 | SR @ 287.84 | SR @ 293.24 | SR @ 298.94 |
|-------------------|----------------------|-----------------------|-----------------------|-----------------------|-----------------------|-----------------------|
| MTBE | 0.0025 | 0.0014 | 0.0016 | 0.0018 | 0.0022 | 0.0029 |
| t12DCE | 0.0043 | 0.0028 | 0.0028 | 0.0027 | 0.0037 | 0.0048 |
| 11DCA | 0.0047 | 0.0031 | 0.0033 | 0.0033 | 0.0039 | 0.0052 |
| c12DCE | 0.0046 | 0.0030 | 0.0031 | 0.0031 | 0.0038 | 0.0051 |
| CHCl3 | 0.0046 | 0.0030 | 0.0031 | 0.0031 | 0.0038 | 0.0051 |
| 111TCA | 0.0066 | 0.0043 | 0.0047 | 0.0047 | 0.0056 | 0.0076 |
| 12DCA | 0.0045 | 0.0029 | 0.0029 | 0.0030 | 0.0036 | 0.0050 |
| BENZ | 0.0050 | 0.0031 | 0.0034 | 0.0035 | 0.0042 | 0.0056 |
| CCl4 | 0.0068 | 0.0044 | 0.0048 | 0.0047 | 0.0058 | 0.0080 |
| TCE | 0.0052 | 0.0030 | 0.0034 | 0.0034 | 0.0043 | 0.0058 |
| 112TCA | 0.0043 | 0.0027 | 0.0027 | 0.0028 | 0.0034 | 0.0048 |
| TOL | 0.0056 | 0.0034 | 0.0039 | 0.0039 | 0.0047 | 0.0062 |
| OCT | 0.0064 | 0.0046 | 0.0050 | 0.0040 | 0.0058 | 0.0089 |
| PCE | 0.0061 | 0.0036 | 0.0043 | 0.0043 | 0.0051 | 0.0069 |
| CIBENZ | 0.0054 | 0.0033 | 0.0039 | 0.0040 | 0.0045 | 0.0059 |
| 1112TetCA | 0.0061 | 0.0037 | 0.0042 | 0.0044 | 0.0050 | 0.0065 |
| EtBENZ | 0.0060 | 0.0037 | 0.0045 | 0.0044 | 0.0052 | 0.0069 |
| mpXYL | 0.0064 | 0.0039 | 0.0048 | 0.0046 | 0.0055 | 0.0072 |
| oXYL | 0.0066 | 0.0041 | 0.0050 | 0.0048 | 0.0057 | 0.0074 |
| 1122TetCA | 0.0044 | 0.0027 | 0.0029 | 0.0031 | 0.0036 | 0.0046 |
| 135TMB | 0.0079 | 0.0046 | 0.0059 | 0.0056 | 0.0071 | 0.0093 |
| 124TMB | 0.0078 | 0.0046 | 0.0060 | 0.0055 | 0.0071 | 0.0092 |
| 13DCB | 0.0072 | 0.0041 | 0.0055 | 0.0053 | 0.0063 | 0.0080 |
| 14DCB | 0.0071 | 0.0040 | 0.0054 | 0.0052 | 0.0062 | 0.0079 |
| 12DCB | 0.0070 | 0.0040 | 0.0053 | 0.0051 | 0.0060 | 0.0076 |
| UNDEC | | 0.0026 | 0.0024 | 0.0020 | 0.0031 | 0.0029 |
| NAPH | | 0.0041 | 0.0056 | 0.0054 | 0.0064 | 0.0081 |
| TRIDEC | | | | | | |
| 2MeNAPH | | 0.0043 | 0.0066 | 0.0066 | 0.0080 | 0.0108 |
| PENTADEC | | | | | | |
| Total mass | 0.1177 | 0.0822 | 0.1339 | 0.1334 | 0.1773 | 0.1981 |

Values in L/hr

Total mass does not include UNDEC, TRIDEC, PENTADEC (28 compounds)

TABLE B
4 CONSTANT REGRESSION OUTPUT

| | Adjusted Rsq | Standard Error | | | | | Std Error ln(SR0) | Std Error b | Std Error - Ea/R | Std Error d |
|------------|-----------------|-------------------|--------|-------|------|--------|-------------------------|-------------------|---------------------------|-------------------|
| MTBE | 0.997 | 0.0960 | -3.217 | 0.981 | 2704 | -0.709 | 0.2881 | 0.0062 | 83 | 0.0082 |
| t12DCE | 0.992 | 0.1659 | -1.877 | 0.905 | 2147 | -0.760 | 0.4971 | 0.0100 | 144 | 0.0138 |
| 11DCA | 0.995 | 0.1272 | -1.346 | 0.916 | 1965 | -0.746 | 0.3809 | 0.0077 | 110 | 0.0106 |
| c12DCE | 0.995 | 0.1299 | -1.905 | 0.911 | 2137 | -0.751 | 0.3892 | 0.0078 | 112 | 0.0109 |
| CHCl3 | 0.996 | 0.1260 | -1.841 | 0.912 | 2118 | -0.748 | 0.3776 | 0.0076 | 109 | 0.0105 |
| 111TCA | 0.995 | 0.1279 | -2.684 | 0.902 | 2259 | -0.761 | 0.3836 | 0.0076 | 111 | 0.0106 |
| 12DCA | 0.995 | 0.1263 | -2.161 | 0.908 | 2218 | -0.746 | 0.3786 | 0.0076 | 109 | 0.0106 |
| BENZ | 0.995 | 0.1323 | -2.207 | 0.920 | 2198 | -0.754 | 0.3965 | 0.0080 | 114 | 0.0110 |
| CCl4 | 0.994 | 0.1405 | -3.121 | 0.889 | 2379 | -0.776 | 0.4220 | 0.0083 | 122 | 0.0116 |
| TCE | 0.992 | 0.1655 | -3.338 | 0.900 | 2522 | -0.772 | 0.4969 | 0.0099 | 144 | 0.0137 |
| 112TCA | 0.995 | 0.1264 | -2.412 | 0.896 | 2302 | -0.724 | 0.3790 | 0.0075 | 109 | 0.0107 |
| TOL | 0.994 | 0.1426 | -2.873 | 0.916 | 2364 | -0.756 | 0.4281 | 0.0087 | 124 | 0.0119 |
| OCT | 0.938 | 0.4698 | -5.984 | 0.822 | 3235 | -0.827 | 1.4231 | 0.0277 | 412 | 0.0388 |
| PCE | 0.991 | 0.1773 | -3.780 | 0.877 | 2601 | -0.775 | 0.5329 | 0.0103 | 154 | 0.0147 |
| CIBENZ | 0.994 | 0.1457 | -2.601 | 0.911 | 2292 | -0.747 | 0.4370 | 0.0088 | 126 | 0.0122 |
| 1112TetCA | 0.996 | 0.1235 | -2.676 | 0.898 | 2281 | -0.725 | 0.3705 | 0.0073 | 107 | 0.0104 |
| EtBENZ | 0.993 | 0.1597 | -2.930 | 0.918 | 2357 | -0.752 | 0.4794 | 0.0097 | 138 | 0.0134 |
| mpXYL | 0.992 | 0.1678 | -3.036 | 0.909 | 2372 | -0.749 | 0.5037 | 0.0101 | 145 | 0.0140 |
| oXYL | 0.993 | 0.1555 | -2.862 | 0.911 | 2312 | -0.740 | 0.4667 | 0.0094 | 135 | 0.0131 |
| 1122TetCA | 0.996 | 0.1118 | -1.971 | 0.913 | 2167 | -0.691 | 0.3351 | 0.0067 | 97 | 0.0096 |
| 135TMB | 0.988 | 0.2024 | -4.435 | 0.897 | 2720 | -0.738 | 0.6093 | 0.0121 | 176 | 0.0170 |
| 124TMB | 0.989 | 0.1997 | -4.126 | 0.890 | 2631 | -0.731 | 0.6009 | 0.0118 | 173 | 0.0169 |
| 13DCB | 0.991 | 0.1832 | -3.422 | 0.888 | 2449 | -0.730 | 0.5503 | 0.0108 | 159 | 0.0155 |
| 14DCB | 0.991 | 0.1802 | -3.263 | 0.892 | 2408 | -0.724 | 0.5413 | 0.0107 | 156 | 0.0153 |
| 12DCB | 0.992 | 0.1697 | -2.970 | 0.894 | 2327 | -0.716 | 0.5092 | 0.0101 | 147 | 0.0144 |
| UNDEC | 0.694 | 0.374 | -1.406 | 0.426 | 1708 | 0.806 | 1.792 | 0.0028 | 517 | 0.053 |
| NAPH | 0.992 | 0.166 | -3.374 | 0.915 | 2430 | 0.671 | 0.497 | 0.010 | 144 | 0.014 |
| TRIDEC | | | | | | | | | | |
| 2MeNAPH | 0.984 | 0.238 | -5.498 | 0.869 | 2990 | 0.689 | 0.72 | 0.014 | 208 | 0.021 |
| PENTADEC | | | | | | | | | | |
| Total mass | 0.993 | 0.1543 | -6.111 | 0.907 | 2419 | -0.732 | 0.4666 | 0.0093 | 134 | 0.0130 |

TABLE C 8260C MASS UNCERTAINTY

AGI 8260C Method for Mass using SPG-0008 Samplers

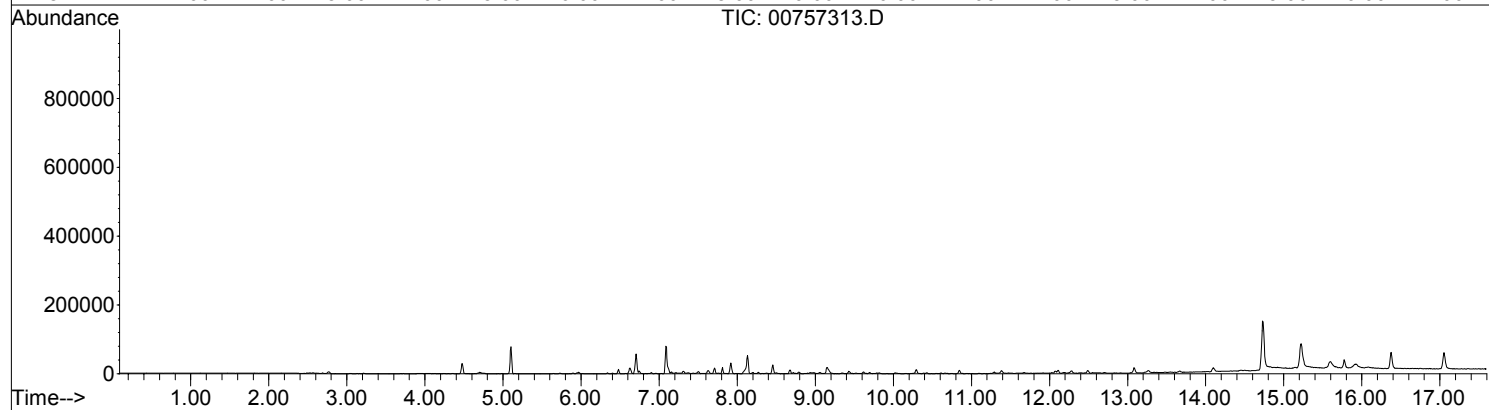
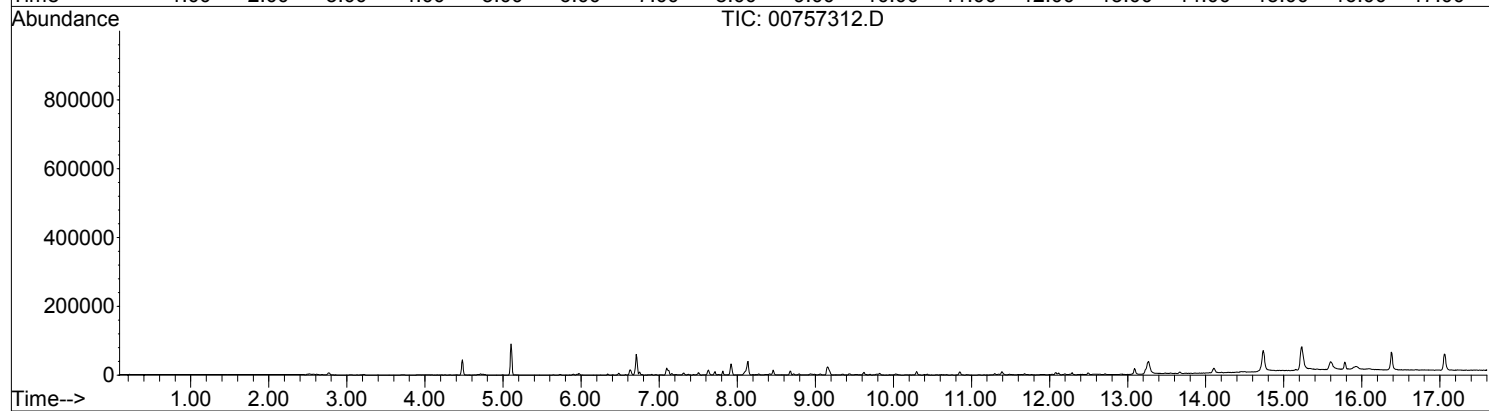
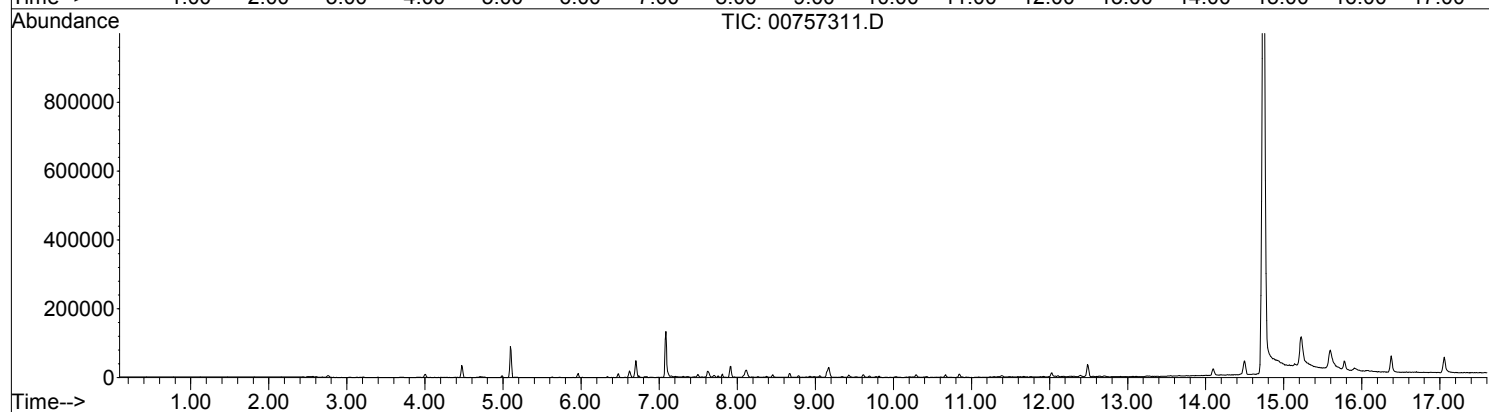
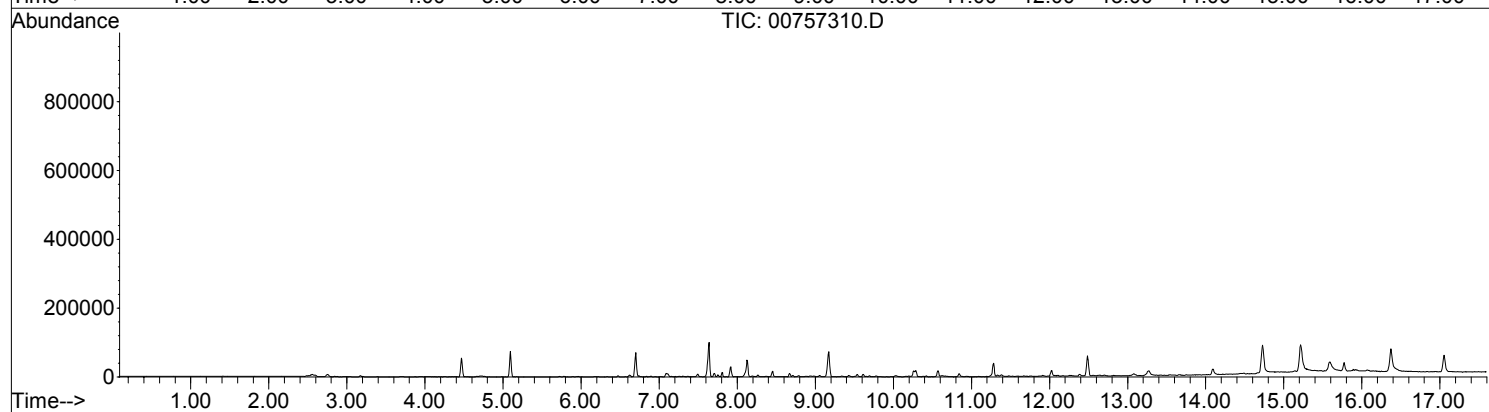
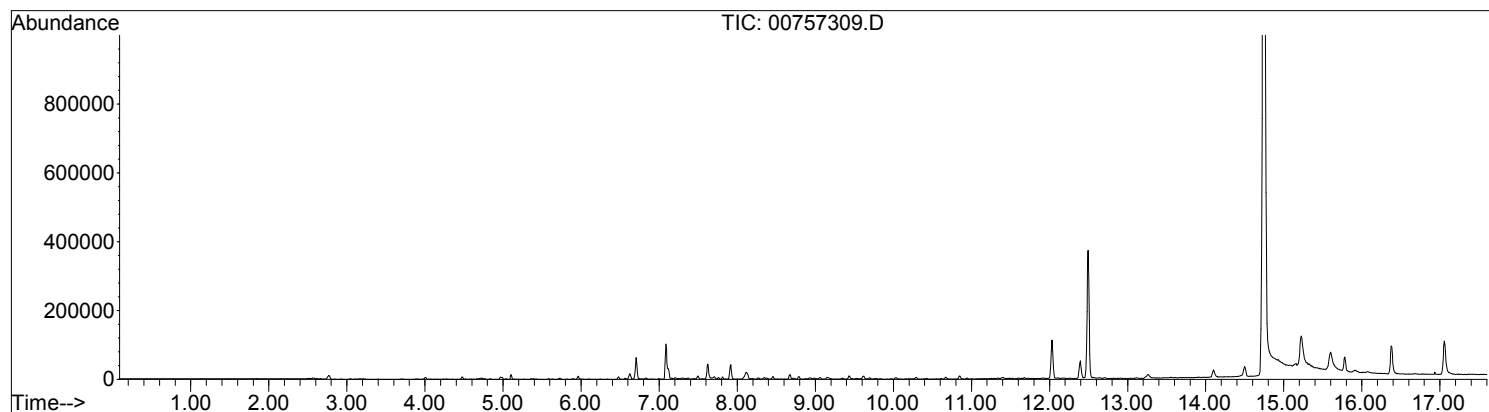
| | 99% Uncertainty Range +/- | 95% Uncertainty Range +/- |
|-----------|---------------------------------|---------------------------------|
| MTBE | 20% | 14% |
| t12DCE | 22% | 15% |
| 11DCA | 18% | 12% |
| c12DCE | 18% | 12% |
| CHCl3 | 16% | 11% |
| 111TCA | 18% | 12% |
| 12DCA | 20% | 13% |
| BENZ | 16% | 10% |
| CCl4 | 19% | 12% |
| TCE | 15% | 10% |
| 112TCA | 18% | 12% |
| TOL | 15% | 10% |
| OCT | 20% | 13% |
| PCE | 16% | 11% |
| CIBENZ | 18% | 12% |
| 1112TetCA | 19% | 13% |
| EtBENZ | 18% | 12% |
| mpXYL | 18% | 12% |
| oXYL | 18% | 12% |
| 1122TetCA | 23% | 15% |
| 135TMB | 21% | 14% |
| 124TMB | 20% | 14% |
| 13DCB | 19% | 13% |
| 14DCB | 19% | 13% |
| 12DCB | 20% | 14% |
| NAPH | 21% | 14% |
| 2MeNAPH | 25% | 17% |

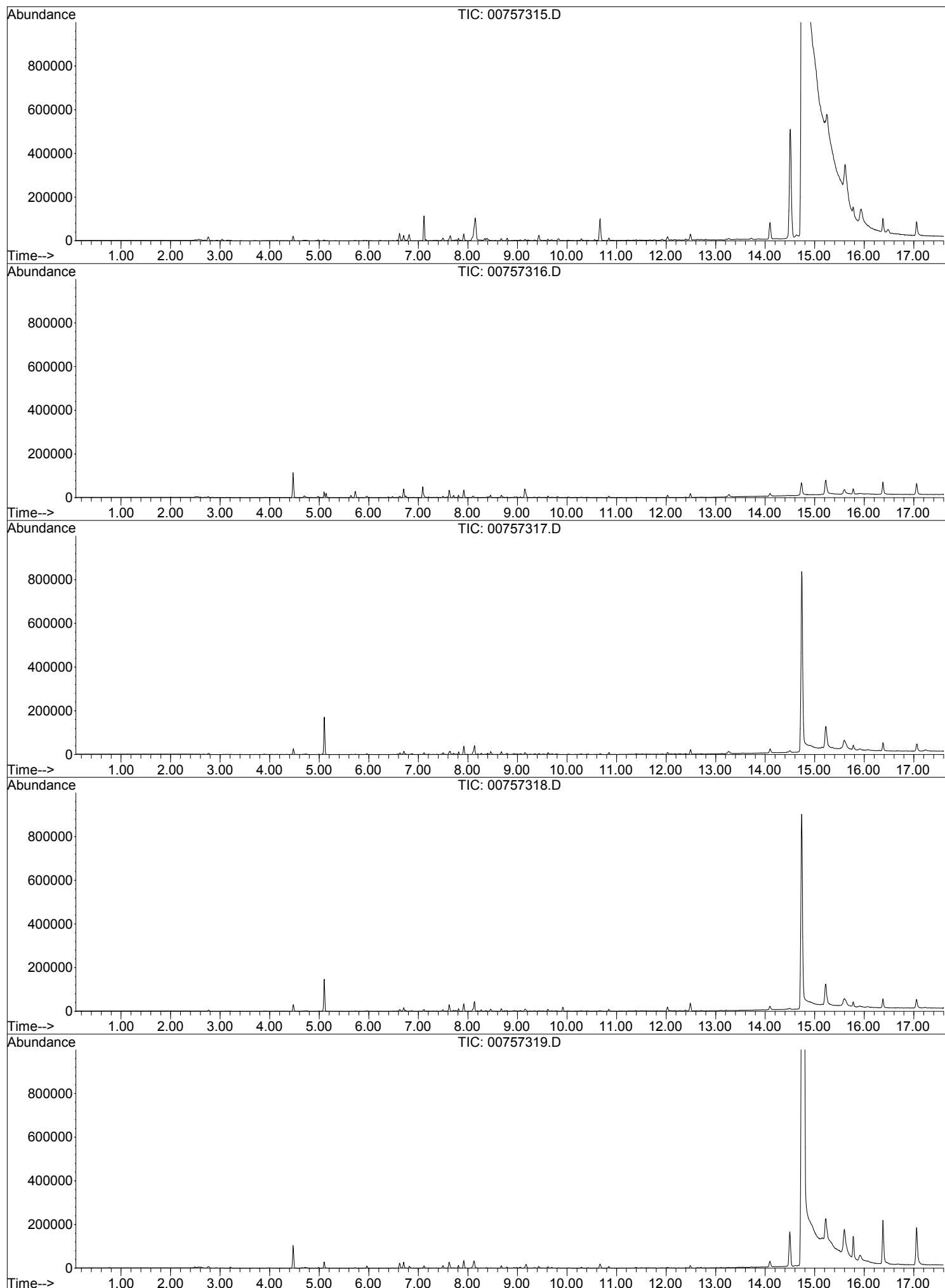
TABLE D
4 CONSTANT WATER CONCENTRATION UNCERTAINTY

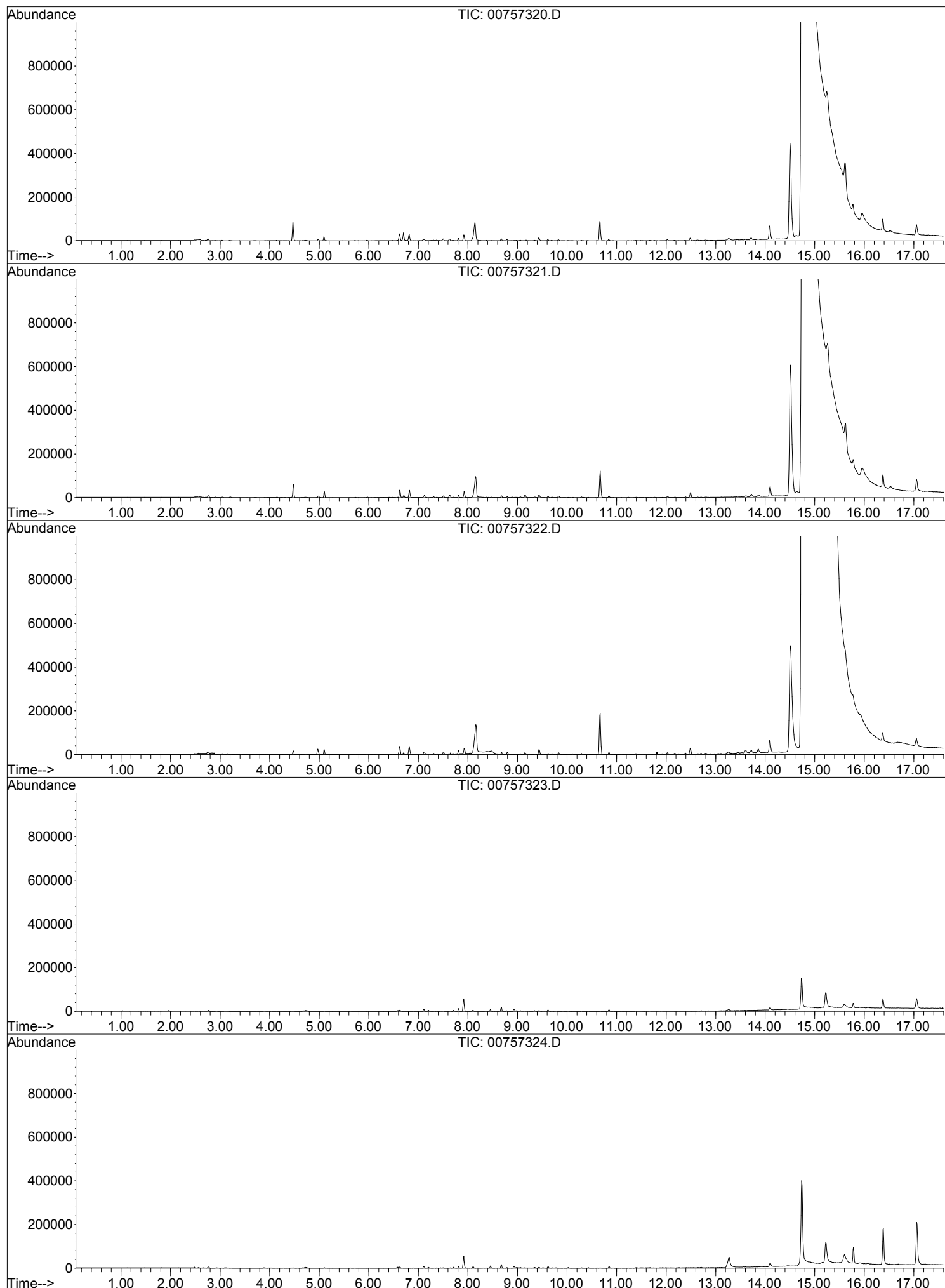
ERROR IN CONCENTRATION REPORTING (1)

| | Average Error | Minimum Error | Maximum Error |
|-----------|--------------------------|--------------------------|--------------------------|
| MTBE | 6% | -12% | 12% |
| t12DCE | 11% | -26% | 21% |
| 11DCA | 8% | -19% | 13% |
| c12DCE | 9% | -19% | 15% |
| CHCl3 | 9% | -20% | 14% |
| 111TCA | 9% | -19% | 23% |
| 12DCA | 10% | -19% | 17% |
| BENZ | 8% | -18% | 13% |
| CCl4 | 10% | -23% | 22% |
| TCE | 10% | -21% | 14% |
| 112TCA | 11% | -21% | 21% |
| TOL | 7% | -17% | 14% |
| OCT | 20% | -41% | 42% |
| PCE | 10% | -24% | 15% |
| CIBENZ | 7% | -16% | 14% |
| 1112TetCA | 8% | -17% | 18% |
| EtBENZ | 6% | -19% | 14% |
| mpXYL | 7% | -22% | 13% |
| oXYL | 7% | -19% | 13% |
| 1122TetCA | 8% | -16% | 17% |
| 135TMB | 9% | -23% | 17% |
| 124TMB | 10% | -28% | 19% |
| 13DCB | 10% | -22% | 17% |
| 14DCB | 10% | -22% | 17% |
| 12DCB | 9% | -23% | 17% |
| NAPH | 10% | -24% | 21% |
| 2MeNAPH | 13% | -32% | 30% |

(1) For 1 hour exposure, includes error related to mass value from AGI analytical method 8260C









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AMPLIFIED
GEOCHEMICAL
IMAGING, LLC

Laboratory Report

Site: Comal and San Marcos Rivers

Prepared for:

SWCA Environmental Consultants
6200 UTSA Boulevard
Suite 102
San Antonio, TX
UNITED STATES

Prepared on:
July 02, 2015

Project Summary and Objective

Amplified Geochemical Imaging, LLC. (AGI) provided the AGI Environmental Survey used at:

Comal and San Marcos Rivers

The service provided by AGI included delivery of the required quantity of AGI Universal Samplers, analysis by the method described below for the requested organic compounds, reporting of the data, and contour mapping (as needed).

This report includes results for only the samples noted under the Laboratory Sample Report section. If contour maps are part of the project deliverable, the maps will be prepared and issued under a separate report cover, upon receipt of a usable sitemap (electronic) and compound choices for contouring.

Written/submitted by:

Jim E Whetzel

Project Manager

Reviewed/approved by:

Don D'Apolito

Project Manager

Analytical data approved by:

Jasmine Smith

Chemist

Quality Assurance Statement

The AGI Laboratory, at Amplified Geochemical Imaging's facility in Elkton, MD USA, operates under the guidelines of its ISO Standard 17025 DoD ELAP accreditation, and its Quality Assurance Manual, Operating Procedures, and Methods (SPG-SOP-0462).

For this project, the analytical method, results, and observations reported do [] do not [✓] fall within the scope of AGI's ISO 17025 accreditation.

Screening/Concentration Method

The AGI Universal Samplers are analyzed at AGI's fixed laboratory using thermal desorption-gas chromatography/mass spectrometry (TD-GC/MS) instrumentation following U.S. EPA Method 8260 (SPG-WI-0292) which includes the following:

- **BFB Tuning Frequency:** A BFB tune is analyzed at the start of each analytical run and after every 30 samples.
- **Initial Calibration:** A minimum of a five point calibration curve is analyzed prior to the analysis of samples.
- **Linearity of Target Compounds:** If the RSD of any target analyte is less than or equal to 25% then average response factor can be used for quantitation. If the RSD exceeds 25% for a target compound a regression equation can be used for quantitation.
- **Continuing Calibration Verification:** After every 10 samples, and at the end of each analytical batch, and a second-source Reference Standard is analyzed near the mid point of the calibration curve. The acceptance criteria for all target analytes in the reference standards are +/- 50% of the true value.
- **Method Blank:** Analyzed prior to the analysis of field samples and every 30 samples.

Note: Analyte levels reported for the field-deployed AGI Universal Samplers that exceed trip and method blank levels, and/or the reporting limit, are more likely to have originated from on-site sources.

| | |
|----------------------------|-------------------|
| Media Sampled: | WATER |
| Chemist - sample analysis: | Kelly J Stringham |
| Chemist - data processor: | Fatima Niazi |
| Chemist - data review: | Jasmine Smith |

Method deviations: None.

Please note that data file names ending with R are rerun samples using the second pair of sorbers, in which the original results were not reported. Data file names ending in D are duplicate analysis results for the second set of sorbers from the same sampler, and are reported.

Additional Report Information

- Comments
- Laboratory Sample Report
- Chain of Custody
- Installation and Retrieval Log
- Analytical Results and Key
- Concentration Calculation Method Summary (as applicable)
- Total Ion Chromatograms

Project Specific Comments

In addition to AGI's standard VOC/ SVOC target compounds pesticides and additional PAH compounds were included in the target compound list. These additional compounds were quantified using a single level calibration standard. Analytical results for these additional compounds were reported only in relative mass values.

Survey period ¹

Samplers were installed on June 1, 2015 and retrieved on June 15, 2015 for an exposure period of 14 days.

Tamper seal intact:

Yes

Date received:

6/18/2015 10:30 AM

By: Clarence W Whigham

COC returned:

Yes

Comments:

Upon retrieval several samples were placed in incorrect sample vials. New labels were placed on vials to indicate actual sample.

Sample 759950 was returned loose in the box. Sample was noted as a trip blank on the installation log, but not on chain of custody. Samples was not analyzed.

¹ - Installation start to end of retrieval, as reported. See installation and retrieval log for individual deployment and retrieval dates and times (i.e., sampler exposure time).

General Comments

Analytical QA/QC

Laboratory instrumentation consists of gas chromatographs equipped with mass selective detectors, coupled with automated thermal desorption units. Sample preparation involves cutting the tip off the bottom of the AGI Universal Sampler, and transferring one or more "sorbents" to a thermal desorption tube for analysis. The insertion/retrieval cord prevents soil, water and other interferences from coming in contact with the adsorbent. No further sample preparation is required. Any replicate sorbents not consumed in the initial analysis will be discarded fifteen (15) days from the date of the laboratory report.

Data are archived and stored in a secure manner as per AGI's Quality Assurance program (SPG-SOP-0462).

Total petroleum hydrocarbons (TPH), gasoline-range petroleum hydrocarbons (GRPH), and/or diesel range petroleum hydrocarbons (DRPH), when reported, are calculated using the area under the peaks observed in m/z 55 and 57 selected ion chromatograms. Quantitation of the mass values was performed using the response factor for a specific alkane (present in the calibration standards). TPH values include the entire chromatogram and provide estimates for aliphatic hydrocarbon ranges of C4 to C20. GRPH and DRPH include only the relevant regions of the chromatograms and provide estimates for C4 to C10 and C10 to C20 aliphatic hydrocarbons, respectively.

Trip blanks were provided to document potential exposures that were not part of the signal of interest (e.g., impact during sampler shipment, installation and/or retrieval, and storage). The trip blanks are identically manufactured and packaged AGI Universal Samplers to those samplers deployed in the field. The trip blanks remain unopened during all phases of the project. Levels reported on the trip blanks may indicate potential impact to the samplers other than the contaminant source of interest.

Unresolved peak envelopes (UPEs) are represented as a series of compound peaks clustered together around a central gas chromatograph elution time in the total ion chromatogram. UPEs may be indicative of complex fluid mixtures. UPEs observed early in the chromatograms are considered to indicate presence of more volatile fluids, while UPEs observed later in the chromatogram may indicate the presence of less volatile fluids. Multiple UPEs may indicate the presence of multiple complex fluids.

Total ion chromatograms (TICs) are included in the Attachments. The eight-digit serial number of each sampler is incorporated in the TIC identification (e.g., 12345678.D represents AGI Universal Sampler 12345678).

Soil Gas Sampling

For soil gas sampling, the AGI Environmental Survey reports mass levels migrating through the open pore spaces of the soil and diffusing through the sampler membrane for sorption by the engineered, hydrophobic adsorbents, housed within the membrane tube. During the migration of the soil gas away from the source to the AGI Universal Sampler, the vapors are subject to a variety of attenuation factors. The soil gas masses reported on the samplers compare favorably with the concentrations reported in the soil or groundwater (e.g., where soil gas levels are reported at greater levels to other sampled locations on the site, the matrix data should reveal the same pattern, and vice versa). However, due to a variety of factors, a perfect comparison between matrix data and soil gas levels can rarely be achieved.

Soil gas concentrations ($\mu\text{g}/\text{m}^3$) are calculated following the method described in the Additional Report Information section.

Soil gas signals reported by this method cannot be correlated specifically to soil adsorbed, groundwater, and /or free-phase contamination. The soil gas signal reported from each AGI Universal Sampler can evolve from all of these sources. Differentiation between soil and groundwater contamination can only be achieved with prior knowledge of the site history (i.e., the site is known to have groundwater contamination only).

Air Sampling

For indoor, outdoor, and crawlspace air sampling, the AGI Environmental Survey reports mass levels present in the air and diffusing through the sampler membrane for sorption by the engineered adsorbents housed within the membrane tube.

Air concentrations ($\mu\text{g}/\text{m}^3$) are calculated following the method described in the Additional Report Information section.

Groundwater and Sediment Porewater Sampling

For groundwater and sediment porewater sampling, the AGI Environmental Survey reports the mass levels of compounds present in the water which, when coming in contact with the sampler membrane, partitions out of solution, and diffuses through the sampler membrane for sorption by the engineered adsorbents.

Water concentrations ($\mu\text{g}/\text{L}$) are calculated using the quantified mass, exposure period and the compound specific uptake rate. The rates were measured under controlled experimental conditions. The uptake rates are corrected for water pressure (depth of the AGI Universal Sampler below the water table), water temperature and the aquifer flow rate. For sediment porewater, the uptake rate is corrected for the reduced volume of water in the sediment, by multiplying the uptake rate by the pore water fraction.

LABORATORY SAMPLE REPORT

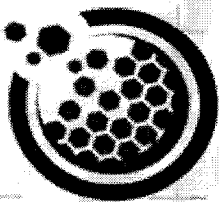
Project: ENV 01387

Site Name: Comal and San Marcos Rivers

Module Type: SPG0008

| Module ID | Sample Type | Field ID | |
|----------------------------|--------------------------|---------------------|-------------------|
| 00759935 | FIELD_SAMPLE | HCS430 | |
| 00759936 | FIELD_SAMPLE | HCS410 | |
| 00759937 | FIELD_SAMPLE | HCS420 | |
| 00759938 | FIELD_SAMPLE | HCS440 | |
| 00759939 | FIELD_SAMPLE | FDHCS440 | |
| 00759940 | FIELD_SAMPLE | HCS460 | |
| 00759941 | FIELD_SAMPLE | HSM410 | |
| 00759942 | FIELD_SAMPLE | HSM420 | |
| 00759943 | FIELD_SAMPLE | HSM430 | |
| 00759944 | FIELD_SAMPLE | FDHSM430 | |
| 00759945 | FIELD_SAMPLE | HSM440 | |
| 00759946 | FIELD_SAMPLE | HSM450 | |
| 00759947 | FIELD_SAMPLE | HSM460 | |
| 00759948 | FIELD_SAMPLE | HSM470 | |
| 00759949 | TRIP_BLANK | TB09 | |
| 00759950 | UNUSED | Not Provided | |
| Total # "FIELD SAMPLES" | Total # "TRIP BLANKS" | Total # "UNUSED" | Total # "LOST" |
| 14 | 1 | 1 | 0 |

Duplicate samples: 0



AMPLIFIED GEOCHEMICAL IMAGING, LLC

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AGI Universal Passive Sampler Chain of Groundwater Sampling

Production Order #: **01387**

Customer Name: SWCA Environmental Consultants
Address: 6200 UTSA Boulevard
Suite 102

San Antonio, TX 78249
USA

Site Name: Comal and San Marcos Rivers
Site Address:

Project Manager:

Serial # of Samplers Shipped
00759935 - 00759950

| | | | |
|--------------------------------|-----------|------------------|---|
| # of Samplers for Installation | 14.00 | # of Trip Blanks | 2 |
| Total Samplers Shipped | 16.00 | Pieces | |
| Total Samplers Received | <u>16</u> | Pieces | |
| Total Samplers Installed | <u>14</u> | Pieces | |

Vial for 00759947
lost in field. Used vial
from extra sampler #00759950
00759950 enclosed in box with loose,
not in a vial

Serial # of Trip Blanks (Client Decides)

| | | |
|----------|--|--|
| 00759949 | | |
|----------|--|--|

| | |
|--|--|
| Prepared By: <u>[Signature]</u> | Is Concurrent water sampling planned this monitoring period? YES <input type="radio"/> NO <input checked="" type="radio"/> |
| Verified By: <u>Lisa Bozette</u> | Scheduled Sampling Date: _____ |
| Installation Performed By: Name: <u>Jennifer Moreland/Brittany Rios</u> Company: <u>SWCA</u> | Retrieval Performed By: Name: <u>Jennifer Moreland/Brittany Rios</u> Company: <u>SWCA</u> |
| Installation Start Date / Time: <u>6/1/15 10:36</u> | Retrieval Start Date / Time: <u>6/15/15 10:26</u> |
| Installation Complete Date / Time: <u>6/1/15 15:59</u> | Retrieval Complete Date / Time: <u>6/15/15 14:01</u> |
| Total Samplers Retrieved: <u>14</u> | |
| Total Samplers Lost In Field: <u>0</u> | |
| Total Unused Samplers Returned: <u>1</u> | |
| Relinquished By: <u>[Signature]</u> Date/Time: <u>5-14-15</u> Company: <u>AGI</u> 9:00 AM | Received By: <u>[Signature]</u> Date/Time: <u>5/18/15</u> Company: <u>SWCA</u> 14:00 |
| Relinquished By: <u>[Signature]</u> Date/Time: <u>6/15/15</u> Company: <u>SWCA</u> 16:15 | Received By: <u>Clarence Whigham</u> Date/Time: <u>6/18/15</u> Company: <u>AGI</u> 10:30 |

of vials labeling of vials with site numbers reconfused in field. Seals w/ site numbers added to clarify incase of any issues. Numbers on installation log correct ~ [Signature]



210 Executive Drive, Suite 1
Newark, DE USA 19702-3335
ph: 302-266-2428

AGI Project No. ENV 01387
Site Name: Comal and San Marcos Rovers
Site Location:

AGI Soil Gas Sampling
Installation & Retrieval Log

Company Name: SWCA
Location: New Braunfels and San Marcos, TX
Samples collected by: Jennifer Moreland, Brittany Rios

* Optional or as needed

| SAMPLER SERIAL NO. | WELL ID | SAMPLE TYPE (Field Sample, Trip Blank, Field Blank, etc.) | INSTALLATION DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 13:00 | RETRIEVAL DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 15:00 | WATER QUALITY MONITORING | | | | | OBSERVATIONS/COMMENTS* (e.g., observations, location description, missing, pulled from well, etc. - as needed) |
|--------------------|----------|---|--|---|--|--|---|---|--|--|
| | | | | | INSTALLATION DEPTH FROM TOP OF CASING (feet) | DEPTH TO WATER FROM TOP OF CASING (feet) | SCREENED INTERVAL FROM TOP OF CASING (feet) | WATER TEMPERATURE AT MODULE DEPTH (deg C) | FLOW RATE THROUGH SCREEN: HIGH OR LOW (HIGH = flow > than 10 m/day; LOW = flow < 10 m/day) | |
| 00759935 | HCS430 | FIELD SAMPLE | 6/1/15 10:36 | 6/15/15 10:26 | 3.25 | | | 24 | HIGH | Recent heavy rain events, both river syst |
| 00759936 | HCS410 | FIELD SAMPLE | 6/1/15 11:03 | 6/15/15 10:39 | 1.1 | | | 24 | HIGH | Depths are from water surface to top of s |
| 00759937 | HCS420 | FIELD SAMPLE | 6/1/15 11:30 | 6/15/15 10:57 | 2.1 | | | 24 | HIGH | |
| 00759938 | HCS440 | FIELD SAMPLE | 6/1/15 11:44 | 6/15/15 11:07 | 1.9 | | | 24 | HIGH | |
| 00759939 | FDHCS440 | FIELD DUPLICATE | 6/1/15 11:44 | 6/15/15 11:07 | 1.9 | | | 24 | HIGH | |
| 00759940 | HCS460 | FIELD SAMPLE | 6/1/15 11:58 | 6/15/15 11:17 | 2.65 | | | 24 | HIGH | |
| 00759941 | HSM410 | FIELD SAMPLE | 6/1/15 13:47 | 6/15/15 12:14 | 1.5 | | | 23 | HIGH | |
| 00759942 | HSM420 | FIELD SAMPLE | 6/1/15 13:58 | 6/15/15 12:26 | 2.9 | | | 23 | HIGH | |
| 00759943 | HSM430 | FIELD SAMPLE | 6/1/15 14:08 | 6/15/15 12:34 | 0.55 | | | 23 | HIGH | |
| 00759944 | FDHSM430 | FIELD DUPLICATE | 6/1/15 14:08 | 6/15/15 12:34 | 0.55 | | | 23 | HIGH | |
| 00759945 | HSM440 | FIELD SAMPLE | 6/1/15 14:23 | 6/15/15 12:47 | 3.75 | | | 23 | HIGH | |
| 00759946 | HSM450 | FIELD SAMPLE | 6/1/15 14:46 | 6/15/15 13:30 | 2.1 | | | 23 | HIGH | |
| 00759947 | HSM460 | FIELD SAMPLE | 6/1/15 15:34 | 6/15/15 13:43 | 0.8 | | | 23 | HIGH | |
| 00759948 | HSM470 | FIELD SAMPLE | 6/1/15 15:59 | 6/15/15 14:01 | 1.8 | | | 23 | HIGH | |
| 00759949 | TB09 | TRIP BLANK | 6/1/15 10:36 | 6/15/15 14:01 | NA | | | | | |
| 00759950 | | TRIP BLANK | | | | | | | | |



**AGI Soil Gas Sampling
Installation & Retrieval L**

* Optional or as needed

| SAMPLER SERIAL NO. | YES / NO | |
|--------------------|--|--------|
| | EVIDENCE OF LIQUID PETROLEUM HYDROCARBONS? | ODOR ? |
| 00759935 | ems up | |
| 00759936 | ampler | |
| 00759937 | | |
| 00759938 | | |
| 00759939 | | |
| 00759940 | | |
| 00759941 | | |
| 00759942 | | |
| 00759943 | | |
| 00759944 | | |
| 00759945 | | |
| 00759946 | | |
| 00759947 | | |
| 00759948 | | |
| 00759949 | | |
| 00759950 | | |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00759935 FIELD SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS430

Installation Date: 6/1/2015 10:36:00AM

Retrieval Date: 6/15/2015 10:26:00AM

Date Analyzed: 6/24/2015 4:48:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.60 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759935 FIELD_SAMPLE

Dilution Factor: 1 Field ID: HCS430

Installation Date: 6/1/2015 10:36:00AM

Retrieval Date: 6/15/2015 10:26:00AM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Reviewer: Jasmine Smith

Matrix: WATER

Product: SPG0008

Date Analyzed: 6/24/2015 4:48:00PM

Batch: ENV-150623-3

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.54 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759936 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS410

Installation Date: 6/1/2015 11:03:00AM

Retrieval Date: 6/15/2015 10:39:00AM

Date Analyzed: 6/24/2015 2:47:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | 0.03 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.14 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759936 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS410

Installation Date: 6/1/2015 11:03:00AM

Retrieval Date: 6/15/2015 10:39:00AM

Date Analyzed: 6/24/2015 2:47:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.66 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759937 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS420

Installation Date: 6/1/2015 11:30:00AM

Retrieval Date: 6/15/2015 10:57:00AM

Date Analyzed: 6/24/2015 9:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.20 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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210 Executive Drive, Suite 1
Newark, DE 19702-3335 USA
ph: +1-302-266-2428
www.agisurveys.net

PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759937 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS420

Installation Date: 6/1/2015 11:30:00AM

Retrieval Date: 6/15/2015 10:57:00AM

Date Analyzed: 6/24/2015 9:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00759938 FIELD_SAMPLE

Dilution Factor: 1 Field ID: HCS440

Installation Date: 6/1/2015 11:44:00AM

Retrieval Date: 6/15/2015 11:07:00AM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Reviewer: Jasmine Smith

Matrix: WATER

Product: SPG0008

Date Analyzed: 6/24/2015 4:18:00PM

Batch: ENV-150623-3

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.40 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00759938 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS440

Installation Date: 6/1/2015 11:44:00AM

Retrieval Date: 6/15/2015 11:07:00AM

Date Analyzed: 6/24/2015 4:18:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.72 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759939 FIELD_SAMPLE
Dilution Factor: 1 Field ID: FDHCS440
Installation Date: 6/1/2015 11:44:00AM
Retrieval Date: 6/15/2015 11:07:00AM
Analyst: Kelly J Stringham
Reviewer: Jasmine Smith

Matrix: WATER Product: SPG0008

Date Analyzed: 6/24/2015 10:49:00PM
Batch: ENV-150623-3
Method: SPG-WI-0292

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.42 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759939 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: FDHCS440

Installation Date: 6/1/2015 11:44:00AM

Retrieval Date: 6/15/2015 11:07:00AM

Date Analyzed: 6/24/2015 10:49:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759940 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS460

Installation Date: 6/1/2015 11:58:00AM

Retrieval Date: 6/15/2015 11:17:00AM

Date Analyzed: 6/24/2015 10:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|--------------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.26 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | 0.03 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00759940 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HCS460

Installation Date: 6/1/2015 11:58:00AM

Retrieval Date: 6/15/2015 11:17:00AM

Date Analyzed: 6/24/2015 10:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 1.90 | 0.50 |
| BTEX | | 0.03 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759941 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM410

Installation Date: 6/1/2015 1:47:00PM

Retrieval Date: 6/15/2015 12:14:00PM

Date Analyzed: 6/24/2015 3:18:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759941 FIELD_SAMPLE
Dilution Factor: 1 Field ID: HSM410
Installation Date: 6/1/2015 1:47:00PM
Retrieval Date: 6/15/2015 12:14:00PM
Analyst: Kelly J Stringham
Reviewer: Jasmine Smith

Matrix: WATER Product: SPG0008

Date Analyzed: 6/24/2015 3:18:00PM
Batch: ENV-150623-3

Method: SPG-WI-0292

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.60 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759942 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM420

Installation Date: 6/1/2015 1:58:00PM

Retrieval Date: 6/15/2015 12:26:00PM

Date Analyzed: 6/24/2015 8:49:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.14 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759942 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM420

Installation Date: 6/1/2015 1:58:00PM

Retrieval Date: 6/15/2015 12:26:00PM

Date Analyzed: 6/24/2015 8:49:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759943 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM430

Installation Date: 6/1/2015 2:08:00PM

Retrieval Date: 6/15/2015 12:34:00PM

Date Analyzed: 6/24/2015 3:48:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 1.98 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759943 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM430

Installation Date: 6/1/2015 2:08:00PM

Retrieval Date: 6/15/2015 12:34:00PM

Date Analyzed: 6/24/2015 3:48:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.96 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759944 FIELD_SAMPLE
Dilution Factor: 1 Field ID: FDHSM430
Installation Date: 6/1/2015 2:08:00PM
Retrieval Date: 6/15/2015 12:34:00PM
Analyst: Kelly J Stringham
Reviewer: Jasmine Smith

Matrix: WATER Product: SPG0008

Date Analyzed: 6/24/2015 7:18:00PM
Batch: ENV-150623-3

Method: SPG-WI-0292

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 2.35 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387

SITE NAME: Comal and San Marcos Rivers

SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**

**San Antonio, TX 78249
USA**

SAMPLER ID: 00759944 FIELD_SAMPLE

Dilution Factor: 1 Field ID: FDHSM430

Installation Date: 6/1/2015 2:08:00PM

Retrieval Date: 6/15/2015 12:34:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Reviewer: Jasmine Smith

Matrix: WATER

Product: SPG0008

Date Analyzed: 6/24/2015 7:18:00PM

Batch: ENV-150623-3

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.54 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759945 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM440

Installation Date: 6/1/2015 2:23:00PM

Retrieval Date: 6/15/2015 12:47:00PM

Date Analyzed: 6/24/2015 8:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.19 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387

SITE NAME: Comal and San Marcos Rivers

SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**

**San Antonio, TX 78249
USA**

SAMPLER ID: 00759945 FIELD_SAMPLE

Dilution Factor: 1 Field ID: HSM440

Installation Date: 6/1/2015 2:23:00PM

Retrieval Date: 6/15/2015 12:47:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Reviewer: Jasmine Smith

Matrix: WATER

Product: SPG0008

Date Analyzed: 6/24/2015 8:19:00PM

Batch: ENV-150623-3

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.68 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759946 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM450

Installation Date: 6/1/2015 2:46:00PM

Retrieval Date: 6/15/2015 1:30:00PM

Date Analyzed: 6/24/2015 6:48:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.09 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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www.agisurveys.net

PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00759946 **FIELD_SAMPLE**

Dilution Factor: 1 Field ID: HSM450

Installation Date: 6/1/2015 2:46:00PM

Retrieval Date: 6/15/2015 1:30:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Reviewer: Jasmine Smith

Matrix: WATER

Product: SPG0008

Date Analyzed: 6/24/2015 6:48:00PM

Batch: ENV-150623-3

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759947 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM460

Installation Date: 6/1/2015 3:34:00PM

Retrieval Date: 6/15/2015 1:43:00PM

Date Analyzed: 6/24/2015 11:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.15 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759947 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM460

Installation Date: 6/1/2015 3:34:00PM

Retrieval Date: 6/15/2015 1:43:00PM

Date Analyzed: 6/24/2015 11:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.67 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00759948 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: HSM470

Installation Date: 6/1/2015 3:59:00PM

Retrieval Date: 6/15/2015 2:01:00PM

Date Analyzed: 6/24/2015 9:49:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|--------------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.18 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | 0.03 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759948 FIELD_SAMPLE
Dilution Factor: 1 Field ID: HSM470
Installation Date: 6/1/2015 3:59:00PM
Retrieval Date: 6/15/2015 2:01:00PM
Analyst: Kelly J Stringham
Reviewer: Jasmine Smith

Matrix: WATER Product: SPG0008

Date Analyzed: 6/24/2015 9:49:00PM
Batch: ENV-150623-3

Method: SPG-WI-0292

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.60 | 0.50 |
| BTEX | | 0.03 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants**
San Antonio, TX 78249
USA

SAMPLER ID: 00759949 TRIP BLANK

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: TB09

Installation Date: 6/1/2015 10:36:00AM

Retrieval Date: 6/15/2015 2:01:00PM

Date Analyzed: 6/24/2015 7:48:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01387
SITE NAME: Comal and San Marcos Rivers
SITE ADDRESS:

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00759949 TRIP BLANK

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: TB09

Installation Date: 6/1/2015 10:36:00AM

Retrieval Date: 6/15/2015 2:01:00PM

Date Analyzed: 6/24/2015 7:48:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150623-3

Reviewer: Jasmine Smith

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |

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ESTIMATED WATER CONCENTRATIONS
COMAL AND SAN MARCOS RIVERS
ORDER # 01387

| | | | | | | | | estimated | | | | | | | | | |
|----------------|--------------|-------------|--------------|--------------|--------------|----|-----------|------------|--------------|-------------|--------------|-------------|--------------|-------------|------------|------------|--|
| DATAFILE | FIELD | DATE/ TIME | DATE/ TIME | DATE/ TIME | DATE/ TIME | | | | | | | | | | | | |
| NAME | ID | INSTALLED | RETRIEVED | RECEIVED | ANALYZED | DF | TPH, ug/L | MTBE, ug/L | t12DCE, ug/L | 11DCA, ug/L | c12DCE, ug/L | CHCl3, ug/L | 111TCA, ug/L | 12DCA, ug/L | BENZ, ug/L | CCl4, ug/L | |
| RL = | | | | | | | 0.053 | 0.013 | 0.007 | 0.007 | 0.007 | 0.007 | 0.005 | 0.008 | 0.006 | 0.005 | |
| 00759935 | HCS430 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | 0.054 | <0.013 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.004 | |
| 00759936 | HCS410 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | 0.059 | <0.013 | <0.007 | <0.007 | <0.007 | 0.012 | <0.005 | <0.008 | <0.006 | <0.004 | |
| 00759937 | HCS420 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | <0.053 | <0.013 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.004 | |
| 00759938 | HCS440 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | 0.062 | <0.013 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.004 | |
| 00759939 | FDHCS440 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | <0.053 | <0.013 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.004 | |
| 00759940 | HCS460 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | 0.093 | <0.013 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.004 | |
| 00759941 | HSM410 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | 0.058 | <0.013 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | |
| 00759942 | HSM420 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | <0.054 | <0.013 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | |
| 00759943 | HSM430 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | 0.071 | <0.013 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | |
| 00759944 | FDHSM430 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | 0.056 | <0.013 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | |
| 00759945 | HSM440 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | 0.061 | <0.013 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | |
| 00759946 | HSM450 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | <0.054 | <0.013 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | |
| 00759947 | HSM460 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | 0.061 | <0.013 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | |
| 00759948 | HSM470 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | 0.058 | <0.013 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | |
| 00759949 | TB09 | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | <0.053 | <0.013 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | |
| BLK_ENV-261689 | Method Blnak | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/24/2015:ET | 1 | <0.053 | <0.013 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | |
| BLK_ENV-261869 | Method Blnak | 6/1/2015:CT | 6/15/2015:CT | 6/18/2015:ET | 6/25/2015:ET | 1 | <0.053 | <0.013 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | |

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ESTIMATED WATER CONCENTRATIONS
COMAL AND SAN MARCOS RIVERS
ORDER # 01387

| DATAFILE | FIELD | | | | | | | | | | | | | | | |
|----------------|--------------|-----------|--------------|-----------|-----------|-----------|--------------|-----------------|--------------|-------------|------------|-----------------|--------------|--------------|-------------|--|
| NAME | ID | TCE, ug/L | 112TCA, ug/L | TOL, ug/L | OCT, ug/L | PCE, ug/L | CIBENZ, ug/L | 1122TetCA, ug/L | ETBENZ, ug/L | mpXYL, ug/L | oXYL, ug/L | 1122TetCA, ug/L | 135TMB, ug/L | 124TMB, ug/L | 13DCB, ug/L | |
| RL = | | 0.006 | 0.009 | 0.006 | 0.005 | 0.005 | 0.006 | 0.007 | 0.005 | 0.005 | 0.005 | 0.010 | 0.005 | 0.005 | 0.006 | |
| 00759935 | HCS430 | <0.006 | <0.009 | <0.005 | <0.004 | 0.101 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.010 | <0.005 | <0.005 | <0.006 | |
| 00759936 | HCS410 | <0.006 | <0.009 | <0.005 | <0.004 | 0.028 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.010 | <0.005 | <0.005 | <0.006 | |
| 00759937 | HCS420 | <0.006 | <0.009 | <0.005 | <0.004 | 0.038 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.010 | <0.005 | <0.005 | <0.006 | |
| 00759938 | HCS440 | <0.006 | <0.009 | <0.005 | <0.004 | 0.071 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.010 | <0.005 | <0.005 | <0.006 | |
| 00759939 | FDHCS440 | <0.006 | <0.009 | <0.005 | <0.004 | 0.075 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.010 | <0.005 | <0.005 | <0.006 | |
| 00759940 | HCS460 | <0.006 | <0.009 | <0.005 | <0.004 | 0.048 | <0.006 | <0.007 | <0.005 | 0.007 | <0.005 | <0.010 | <0.005 | <0.005 | <0.006 | |
| 00759941 | HSM410 | <0.006 | <0.010 | <0.006 | <0.005 | <0.005 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 | |
| 00759942 | HSM420 | <0.006 | <0.010 | <0.006 | <0.005 | 0.030 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 | |
| 00759943 | HSM430 | <0.006 | <0.010 | <0.006 | <0.005 | 0.299 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 | |
| 00759944 | FDHSM430 | <0.006 | <0.010 | <0.006 | <0.005 | 0.349 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 | |
| 00759945 | HSM440 | <0.006 | <0.010 | <0.006 | <0.005 | 0.039 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 | |
| 00759946 | HSM450 | <0.006 | <0.010 | <0.006 | <0.005 | 0.020 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 | |
| 00759947 | HSM460 | <0.006 | <0.010 | <0.006 | <0.005 | 0.031 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 | |
| 00759948 | HSM470 | <0.006 | <0.010 | <0.006 | <0.005 | 0.037 | <0.006 | <0.007 | <0.005 | 0.007 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 | |
| 00759949 | TB09 | <0.006 | <0.009 | <0.006 | <0.004 | <0.005 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.010 | <0.005 | <0.005 | <0.006 | |
| BLK_ENV-261689 | Method Blnak | <0.006 | <0.009 | <0.006 | <0.005 | <0.005 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.010 | <0.005 | <0.005 | <0.006 | |
| BLK_ENV-261869 | Method Blnak | <0.006 | <0.009 | <0.006 | <0.005 | <0.005 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.010 | <0.005 | <0.005 | <0.006 | |

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ESTIMATED WATER CONCENTRATIONS
COMAL AND SAN MARCOS RIVERS
ORDER # 01387

| DATAFILE | FIELD | | | | | estimated: | | estimated: | estimated: | estimated: | estimated |
|----------------|--------------|-------------|-------------|-------------|------------|--------------|---------------|----------------------|----------------|--------------------|----------------|
| NAME | ID | 14DCB, ug/L | 12DCB, ug/L | UNDEC, ug/L | NAPH, ug/L | TRIDEC, ug/L | 2MeNAPH, ug/L | Acenaphthylene, ug/L | PENTADEC, ug/L | Acenaphthene, ug/L | Fluorene, ug/L |
| RL = | | 0.006 | 0.006 | 0.020 | 0.016 | 0.020 | 0.013 | 0.014 | 0.020 | 0.014 | 0.014 |
| 00759935 | HCS430 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759936 | HCS410 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759937 | HCS420 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759938 | HCS440 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759939 | FDHCS440 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759940 | HCS460 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759941 | HSM410 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759942 | HSM420 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759943 | HSM430 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759944 | FDHSM430 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759945 | HSM440 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759946 | HSM450 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759947 | HSM460 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759948 | HSM470 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00759949 | TB09 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| BLK_ENV-261869 | Method Blnak | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| BLK_ENV-261869 | Method Blnak | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |

KEY TO DATA TABLE

UNITS

| | |
|-------------------|--|
| µg | micrograms, relative mass value |
| µg/m ³ | micrograms per cubic meter; estimated soil gas concentration |
| µg/L | micrograms per Liter; calculated water concentration |

DATA QUALIFIERS

| | |
|---|--|
| > | greater than; value exceeds calibration range, estimated value |
| < | less than; compound value is below the LOD and RL |
| J | mass value below LOQ or RL, but above LOD, estimated mass value |
| E | mass value exceeds upper calibration level, estimated mass value |
| Q | one or more quality control parameters failed for the compound |

ABBREVIATIONS

| | |
|--------|--|
| AVG RL | average reporting limit; calculated based on individual field sample RLs |
| LOD | limit of detection |
| LOQ | limit of quantification |
| MDL | method detection limit |
| RL | reporting limit |

| | | | |
|-------------|---|----------|--|
| 1112TetCA | 1,1,1,2-tetrachloroethane | CIBENZ | chlorobenzene |
| 111TCA | 1,1,1-trichloroethane | ct12DCE | cis- & trans-1,2-dichloroethene |
| 1122TetCA | 1,1,2,2-tetrachloroethane | EtBENZ | ethylbenzene |
| 112TCA | 1,1,2-trichloroethane | mpXYL | m-, p-xylene |
| 11DCA | 1,1-dichloroethane | MTBE | methyl t-butyl ether |
| 11DCE | 1,1-dichloroethene | NAPH | naphthalene |
| 124TMB | 1,2,4-trimethylbenzene | OCT | octane |
| 12DCA | 1,2-dichloroethane | oXYL | o-xylene |
| 12DCB | 1,2-dichlorobenzene | PCE | tetrachloroethene |
| 135TMB | 1,3,5-trimethylbenzene | PENTADEC | pentadecane |
| 13DCB | 1,3-dichlorobenzene | PHEN | phenanthrene |
| 14DCB | 1,4-dichlorobenzene | t12DCE | trans-1,2-dichloroethene |
| 2MeNAPH | 2-methyl naphthalene | TCE | trichloroethene |
| BENZ | benzene | TMBs | combined masses of 1,3,5-trimethylbenzene and 1,2,4-trimethylbenzene |
| BTEX | combined masses of benzene, toluene, ethylbenzene, and total xylenes (Gasoline Range Aromatics) | TOL | toluene |
| C11,C13&C15 | combined masses of undecane, tridecane, and pentadecane (C11+C13+C15) (Diesel Range Alkanes) | TPH | total petroleum hydrocarbons |
| c12DCE | cis-1,2-dichloroethene | TRIDEC | tridecane |
| CCl4 | carbon tetrachloride | UNDEC | undecane |
| CHC13 | chloroform | VC | vinyl chloride |

SUMMARY OF SAMPLING RATE CALIBRATION FOR AGI UNIVERSAL SAMPLER IN AQUEOUS PHASES

INTRODUCTION:

The Amplified Geochemical Imaging, LLC (AGI) passive vapor sampler is designed to be used for soil gas, water, sediment pore water, and air sampling. This document describes the process used to calibrate the sampler's compound specific sampling or uptake rates in aqueous phases.

Sampling rates are measured following AGI's "Standard Practice for Determining the Sampling Rate of Passive Diffusion Samplers in Various Environmental Media": SPG-SOP-0493. Rates are used to calculate dissolved phase concentrations of volatile and semi-volatile contaminants in water. The calibration process is summarized in three parts: Part 1: shallow water, Part 2: deep water, and Part 3: sediment.

PURPOSE:

The purpose of this document is to:

1. Summarize the test protocol,
2. Summarize the methodology for analysis of data,
3. Present general results for generating concentration calibration of the AGI Universal Sampler

Principle of Operation of the AGI Sampler

The AGI Universal Sampler is designed with solid adsorbents enclosed inside a tubular microporous PTFE membrane. When placed in water, the pores and hydrophobic nature of the PTFE keep liquid water from entering the membrane until a water head of about 34 feet is reached. The membrane will not keep water vapor from entering but the adsorbents are very hydrophobic and through testing validated to be unaffected by this moisture vapor. In shallow water, <34', volatile and semi-volatile compounds will partition from the dissolved water into the air phase in the PTFE membrane according to Henry's Law. This partitioning is instantaneous and within seconds-minutes, the compound is adsorbed by the adsorbent inside the sealed tube. Because the diffusivity in air is about 10,000 times higher than the diffusivity in water, the sampling rate is controlled by the water contact area of the membrane that allows the Henry's Law effect to occur. This contact area is set by the membrane diameter and length of the sealed tube, which is fixed in AGI's manufacturing process.

Henry's law as well as diffusivity, which are fundamentally incorporated into the sampling rate, are affected by temperature, T , and follow an Arrhenius equation $H_T = H_r \times \exp\left(\frac{-E_a/R}{1/T_r - 1/T}\right)$. Because a 5°C temperature change can make a 15% change in sampling rate, the temperature of the sampled water should be known to get the most precise concentration.

The membrane pore size is also small enough that colloidal particles and microbes cannot pass through the membrane. This keeps the adsorbent from getting contaminated and eliminates any need to add preservative or chilling during storage or transportation.

When the water pressure exceeds the water entry pressure of the membrane, about 34 feet of water, the water becomes in contact with the solid adsorbent. Under this condition the compounds in the water will partition from the water into the solid. The partitioning coefficient, K_{AW} , can be approximated by the octanol-water coefficient, K_{OW} , but has been measured more precisely in the lab for AGI's specific solid adsorbent. The sampling rate is the product of the sampling rate at <34' of water and the K_{AW} .

In sediment, the sampler measures pore-water concentration, which is generally agreed to be the preferred measurement as it is more indicative of bioavailability. In sediment the volumetric

availability of water to the sampler is reduced by the volume fraction solids in the sediment, which typically varies from zero to 35%, but can be as high as 73% in well packed and broad particle size distribution sediments. As a result, sampling rates in sediment are multiplied by the fraction pore water in the sediment to determine concentration.

PART 1: Calibration in shallow water

Part 1 summarizes the work in shallow water generating calibration data, evaluating the physical and chemical factors affecting the sampling rate, and measurement of the actual sampling rates or regression calibration equations needed to determine concentrations.

Sample Generation in water

In this calibration work, solutions of analytes at known concentrations were formulated in clean 4 liter smoked glass jugs by injecting microliter measured amounts of environmental standards using a calibrated syringe into pure or deionized water and stirring for a minimum of 2 hours but generally overnight. Headspace in the jugs was minimized and generally less than 1% by volume during the tests. Jugs were temperature controlled by placing them in a water filled cooler, chilled via a copper tubing loop in the cooler. Temperature was measured with a certified digital temperature gauge and an average value used for each temperature experiment.

AGI samplers were weighted so they won't float and placed in the jugs at time zero. They were removed at various intervals to generate samples along with duplicates that showed mass increasing with exposure time. The sampler exposure time was selected to span minutes to hours and was generally reduced for high concentration tests to maintain uptake with time in roughly the linear dynamic range. Samplers were removed and dried with a paper towel and returned to their original container for analysis. They were analyzed by AGI's 8260C (SPG-WI-318 or SPG-WI-10028) method in duplicate, which is based on EPA SW846 Method 8260C.

Water samples were also taken and measured at an outside accredited lab using EPA SW846 Method 8260B. The concentrations agreed well with the calculated concentrations based on the standard certification, jug volume, and syringe injection. The variability of the outside lab 8260B values were found to be high, so for the sampling rate calculations we used the concentrations based on syringe dosing.

Calibrations were run at five concentrations, nominally at 6, 24, 118, 590, 1420 ug/L and five temperatures nominally at 5, 10, 15, 20, and 25 degrees centigrade. Samples were taken at 4 different exposure times. Samples were run in duplicate. A total of 176 data points were generated using 28 compounds from AGI's standard compounds list. Tridecane and pentadecane were not evaluated due to their very low solubility in water. In addition, another 23 compounds were tested using an 8260 liquid standard at nominal concentrations of 0.5, 1.0, 5.0, 15, 95, and 470 ug/L at a temperature typical of groundwater, 15°C. This is a living calibration and as additional data are generated, they may be qualified and added to this data set to improve the precision of the sampling rate calibration and broaden the compound list.

Key Variable Effects

As expected from theory, at short to moderate exposure times, mass will increase roughly linearly proportional to exposure time, as well as proportional to concentration, and exponentially with temperature following Arrhenius law. Temperature affects the Henry's law as well as diffusivity in water. Sampling rate is generally independent of concentration and time at mass values significantly below saturation. In the following sections we have characterized the sampling rate for each compound as affected by temperature and also developed calibrations using regression which account for the minor impact of time, and mass.

Concentration using Simple Sampling Rate Determination

A simple way to determine concentration is to measure mass on the AGI sampler, divide by exposure time, and divide by sampling rate, SR.

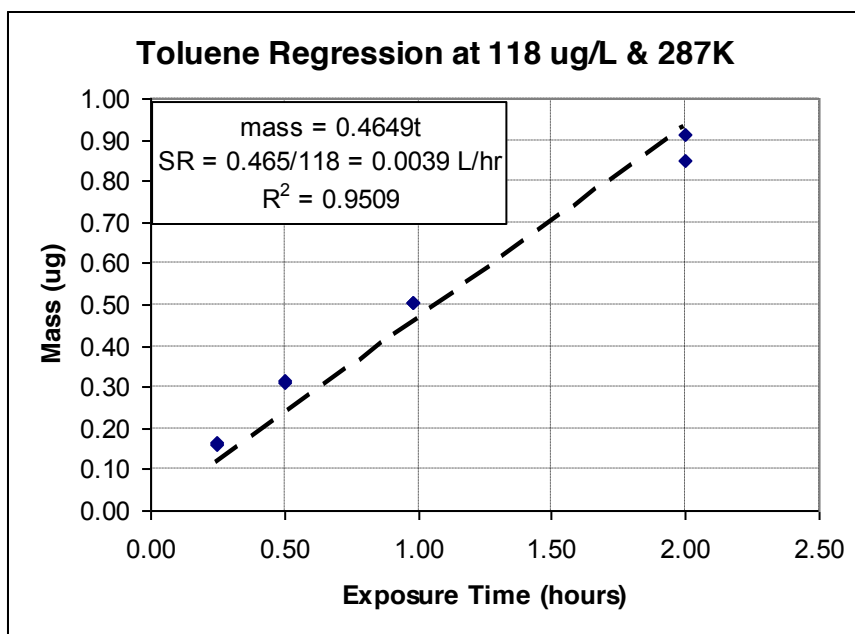
$$\text{Conc [ug/L]} = \text{mass/time/SR} \quad (1)$$

The sampling rate can be determined via measurements of mass versus time at a known concentration and temperature according to the following modification of equation (1).

$$\text{SR} = \text{mass/time/concentration} \quad (2)$$

Sampling rates in L/hr were determined by measuring the trend or regression mass uptake versus time and dividing by the concentration. A measurement like this will use 8 data points (4 times x 2 samples). Such a sampling rate can be measured at any concentration and temperature.

The chart to the right shows a plot of mass versus time for water at 118 ug/L and 287K (actual data from a single run). Slope of 0.465 ug/hr divided by the concentration of 118 ug/L yields a sampling rate, SR, of 0.0039 L/hr.



SR's typically range from about 0.004 to 0.007 L/hr at 15°C. Table A shows SR's measured for our standard compound list at 5 temperatures.

Rigorous Concentration using Regression

A preferred method for determining concentration that will yield improved accuracy over a wide range of concentrations, exposure times, and temperatures is to use all data in a regression analysis, which allows adjustments for the minor non-linear influences of mass and time as well as the effects of temperature. This step is done by regressing equation (1) or a universal version of equation (1):

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRo} * \exp(-\text{Ea}/\text{R}/\text{T})] \quad (3)$$

The subtle non-linear effects of mass and time will be evident in the deviation of coefficients b and d from 1.0. This regression generates four constants b, d, SRo, and $-\text{Ea}/\text{R}$ by regressing $\ln(\text{conc})$ versus $\ln(\text{mass})$, $\ln(\text{time})$, $1/\text{temp}$. These four constants can be used to determine concentration via the equation:

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRo} * \exp(-\text{Ea}/\text{R}(1/\text{T}))] \quad (4)$$

Where conc is in ug/L, mass is in ug, time in hours, T in degrees Kelvin.

Equation (4) can be also expressed at a reference temperature, Tr , such as 15°C by

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRr} * \exp(-\text{Ea}/\text{R}(1/\text{Tr}-1/\text{T}))] \quad (5)$$

This step allows sampling rates, SRr, at any reference temperature, Tr , and for any analyte to easily be compared. The values of SRr at 293.14K can be found in Table A.

When sampling times are between 0 and 4 hours, using the 4 constant equation (5) is preferred. For concentrations from about 5 to 1500 ug/L one hour exposure times generally give the lowest error, typically with average error of 6-20% and with total error range of 12%-32%. For low concentrations where sampling times are greater than 4 hours, it is preferred to use equation (1) to avoid unrealistic effects from the coefficient d or to set d to 1.0. In such a case SR in equation (1) can be substituted with $[\text{SRr} * \exp(-\text{Ea}/\text{R}(1/\text{Tr}-1/\text{T}))]$ to use an SR representative of the well temperature, T.

The chart to the right is a plot of the calculated concentration from the 4 constant regression compared to the dosed concentration. Agreement is excellent for the 176 data points.

However, there does appear to be a slight high bias of 8.6% over the full range of this data, although it is well within acceptable limits of variability.

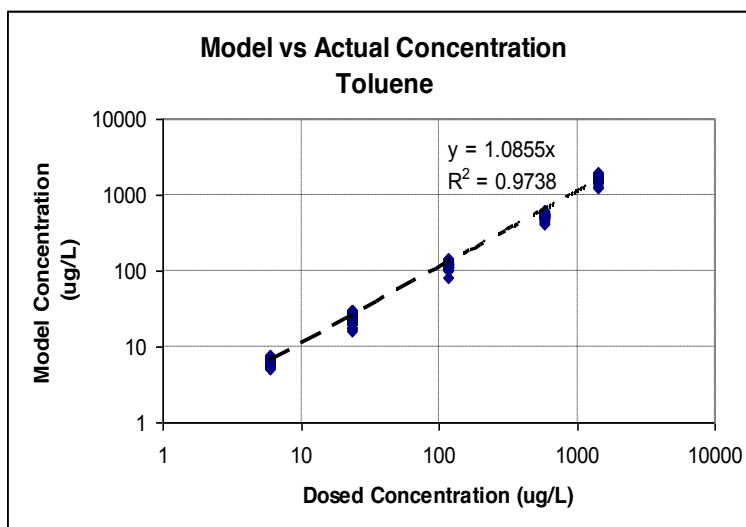


Table B shows the tabulated summary of the 4 constants regression with Rsq values and error estimates for the 4 constants for each analyte. Most regression Rsq values are 0.99 or greater for each analyte. In general, $-E_a/R$ is about 2400 \pm 400, b is about 0.9, d is about -0.75, and SR(15°C) ranges from .004 L/hr to 0.007 L/hr increasing with MW of the compound.

Error Estimates

The error in the water concentration values will depend on both the error in mass from the analytical method as well as the error in the concentration calibration. Table C shows the error in the mass values from the 8260C low sensitivity method.

The standard error of the regression and standard errors of the constants can be found in table B. For each compound we have measured the error between the derived concentration and the actual concentration. The error tends to be lowest at our recommended exposure time of one hour as shown by the example for Toluene to the right.

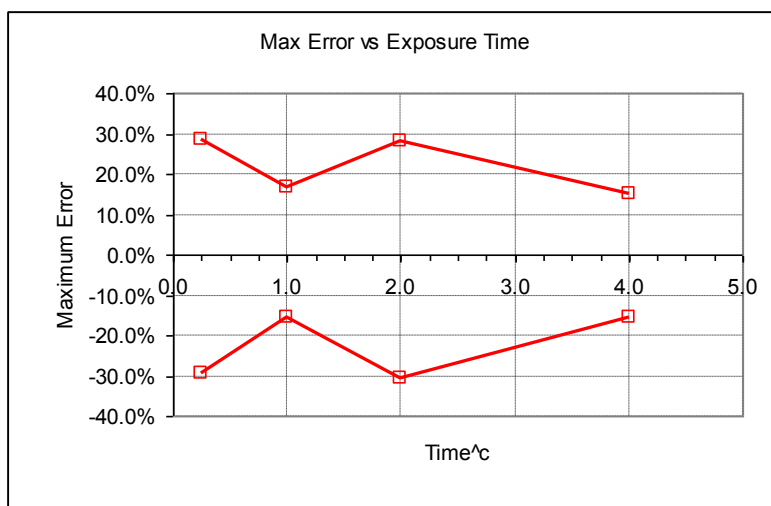
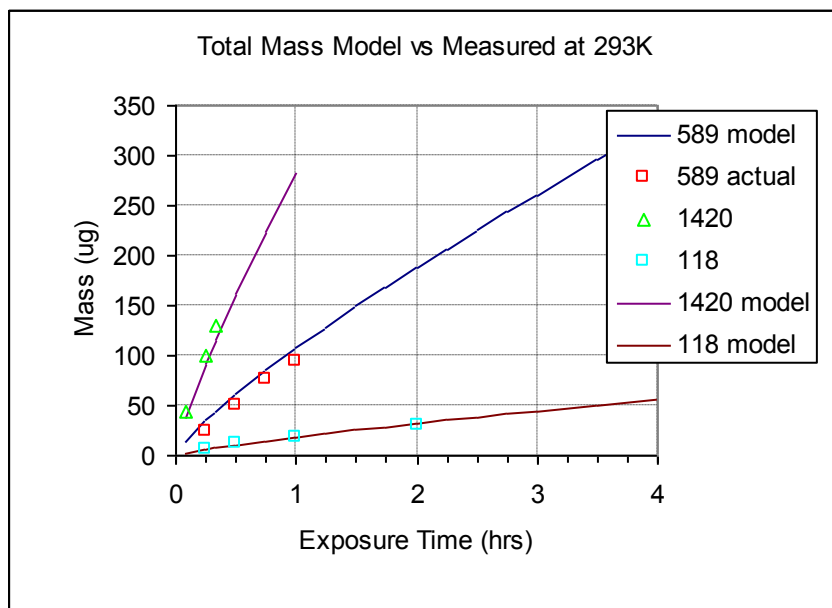


Table D shows the total average error in water concentration by compound as well as the low and high error. The average ranges from about 6% to 20%, which is similar to the analytical method errors. The low and high errors range from 12% to 32% and include contribution from measurement errors in both time and temperature.

Sorbent Saturation

As mass increases on a solid sorbent and approaches saturation, reverse diffusion can occur causing the sampling rate to drop. Eventually the mass level will reach a maximum steady state value at any concentration. A rate of mass uptake with time that deviates significantly from linear, indicates that sorbent saturation could be an issue. When using equation (1), staying in the linear range to avoid the effects of adsorbent saturation is important. We recommend keeping the total mass on the sampler below 50 ug or flagging when this is exceeded.

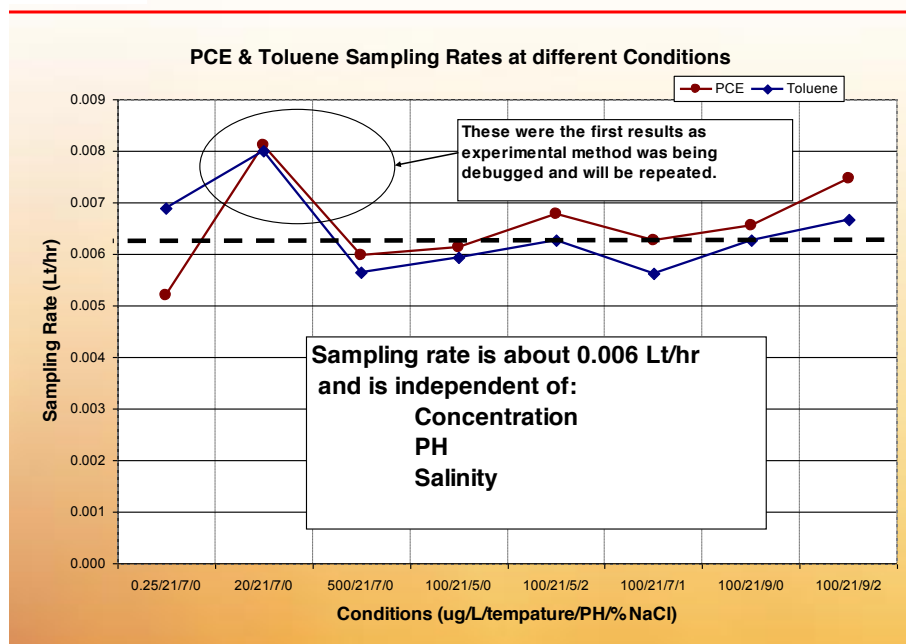
The 4 constant regression accounts for some of the non linearity allowing good accuracy at higher mass levels. From the experimental data we have found this safe range can be extended to 100 ug or higher as shown in the chart below. This chart compares total mass of all compounds (excluding heavy alkanes, which have solubility issues) versus time in comparison to that predicted from the 4-constant concentration equation.



Effect of PH and Salinity

Because neither PH nor salinity is known to have a significant impact on Henry's law or diffusivity in water, we did not expect them to have a significant impact on sampling rate. To confirm this, experiments were run varying PH from 5 to 9 and NaCl content from 0 to 2%. The chart below shows no significant impact for combinations of PH and NaCl content over this range on the sampling rate of toluene in water at 21°C.

Checked for Effects of PH & Salinity

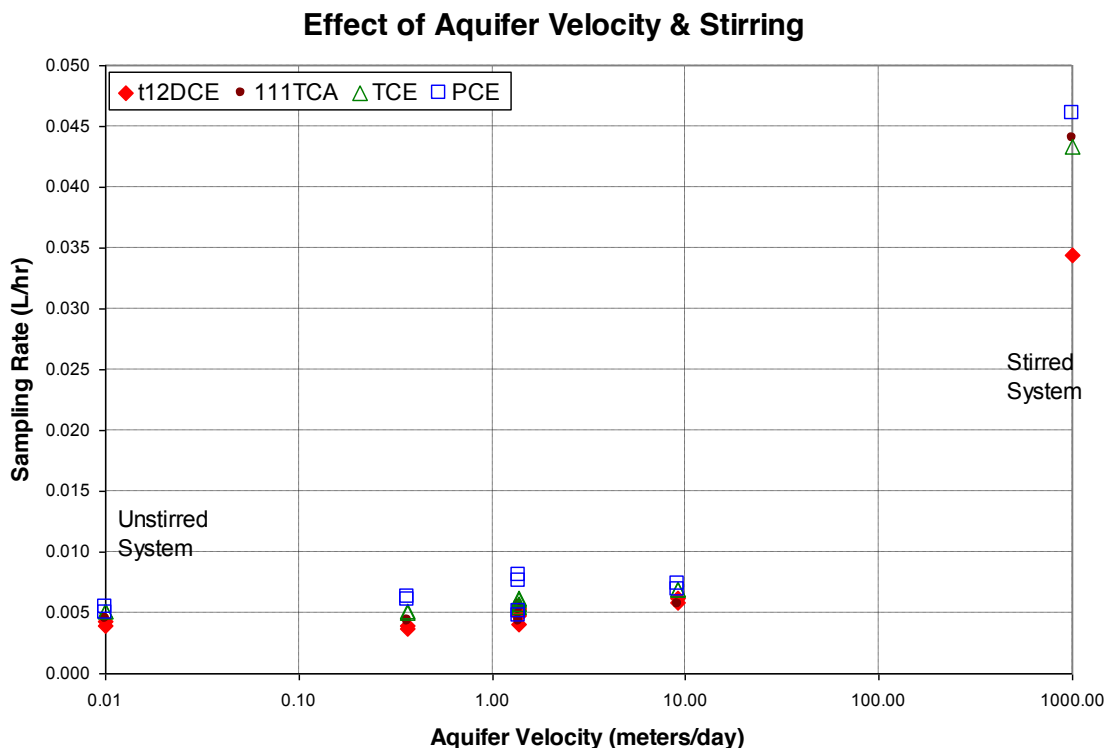


Impact of Aquifer Velocity

The velocity in most aquifers is quite slow, typically a meter/day or less. Occasionally water flow could be much higher such as encountered in karst aquifers, streams or rivers. Mass transfer coefficients are higher in high flow conditions, which will lead to higher sampling rates. We validated that a highly stirred system had sampling rates about 10 times higher than those that were non-stirred. We decided to evaluate the effect of aquifer velocity.

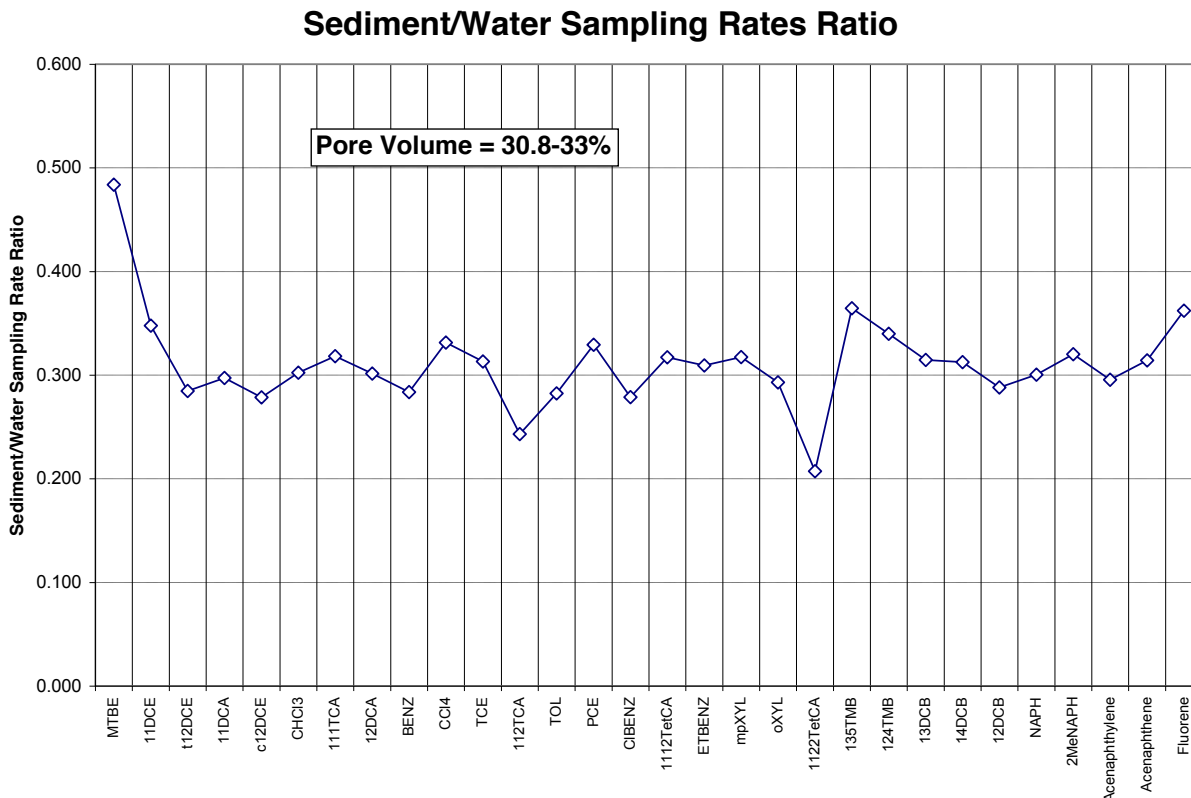
A test apparatus was built comprising a 3" PVC pipe tee filled with clean sand in each of the horizontal straight legs and screened to leave the center open. A test solution was run through this system using a variable flow pump and AGI samplers were placed into the simulated well through the vertical leg of the tee. Tests were run to examine the effect of velocity by varying the pumping rate and hence water velocity.

The chart below shows no significant effect of aquifer velocity up to a speed of about 10 meters/day. At velocities significantly above this, similar to a stirred system, sampling rates are about 10 times higher.



concentration. Pore water fraction can range from 1.0 for water without sediment to as low as 0.25. Typically most sediments have pore fractions of 0.9 to 0.65.

A sampling rate study was done with water and with water added into a well-packed sorted sand. Pore water fraction in this test was measured between 30.8% and 33% by volume. Below is a plot of the ratio of sampling rates measured in the sediment to open water. The average ratio is equal to the pore water fraction confirming that sampling rate in sediment is on average equal to the product of pore water fraction times the sampling rate in water.



Summary

The AGI Sampler can be used to determine the concentration of volatile and semi-volatile compounds in a water phase. This requires knowing the exposure time and water temperature. It also requires knowing if the sample is above or below 34' of water head and if the water has a velocity above 10 meters/day. Regressions of large amounts of data were used to generate a four constant equation to generate concentration values in water. Potential error in the concentration values is excellent typically less than 25%.

TABLE A
WATER SAMPLING RATES STANDARD LIST

| | SRr 293.14 | SR @ 277.54 | SR @ 282.44 | SR @ 287.84 | SR @ 293.24 | SR @ 298.94 |
|-------------------|----------------------|-----------------------|-----------------------|-----------------------|-----------------------|-----------------------|
| MTBE | 0.0025 | 0.0014 | 0.0016 | 0.0018 | 0.0022 | 0.0029 |
| t12DCE | 0.0043 | 0.0028 | 0.0028 | 0.0027 | 0.0037 | 0.0048 |
| 11DCA | 0.0047 | 0.0031 | 0.0033 | 0.0033 | 0.0039 | 0.0052 |
| c12DCE | 0.0046 | 0.0030 | 0.0031 | 0.0031 | 0.0038 | 0.0051 |
| CHCl3 | 0.0046 | 0.0030 | 0.0031 | 0.0031 | 0.0038 | 0.0051 |
| 111TCA | 0.0066 | 0.0043 | 0.0047 | 0.0047 | 0.0056 | 0.0076 |
| 12DCA | 0.0045 | 0.0029 | 0.0029 | 0.0030 | 0.0036 | 0.0050 |
| BENZ | 0.0050 | 0.0031 | 0.0034 | 0.0035 | 0.0042 | 0.0056 |
| CCl4 | 0.0068 | 0.0044 | 0.0048 | 0.0047 | 0.0058 | 0.0080 |
| TCE | 0.0052 | 0.0030 | 0.0034 | 0.0034 | 0.0043 | 0.0058 |
| 112TCA | 0.0043 | 0.0027 | 0.0027 | 0.0028 | 0.0034 | 0.0048 |
| TOL | 0.0056 | 0.0034 | 0.0039 | 0.0039 | 0.0047 | 0.0062 |
| OCT | 0.0064 | 0.0046 | 0.0050 | 0.0040 | 0.0058 | 0.0089 |
| PCE | 0.0061 | 0.0036 | 0.0043 | 0.0043 | 0.0051 | 0.0069 |
| CIBENZ | 0.0054 | 0.0033 | 0.0039 | 0.0040 | 0.0045 | 0.0059 |
| 1112TetCA | 0.0061 | 0.0037 | 0.0042 | 0.0044 | 0.0050 | 0.0065 |
| EtBENZ | 0.0060 | 0.0037 | 0.0045 | 0.0044 | 0.0052 | 0.0069 |
| mpXYL | 0.0064 | 0.0039 | 0.0048 | 0.0046 | 0.0055 | 0.0072 |
| oXYL | 0.0066 | 0.0041 | 0.0050 | 0.0048 | 0.0057 | 0.0074 |
| 1122TetCA | 0.0044 | 0.0027 | 0.0029 | 0.0031 | 0.0036 | 0.0046 |
| 135TMB | 0.0079 | 0.0046 | 0.0059 | 0.0056 | 0.0071 | 0.0093 |
| 124TMB | 0.0078 | 0.0046 | 0.0060 | 0.0055 | 0.0071 | 0.0092 |
| 13DCB | 0.0072 | 0.0041 | 0.0055 | 0.0053 | 0.0063 | 0.0080 |
| 14DCB | 0.0071 | 0.0040 | 0.0054 | 0.0052 | 0.0062 | 0.0079 |
| 12DCB | 0.0070 | 0.0040 | 0.0053 | 0.0051 | 0.0060 | 0.0076 |
| UNDEC | | 0.0026 | 0.0024 | 0.0020 | 0.0031 | 0.0029 |
| NAPH | | 0.0041 | 0.0056 | 0.0054 | 0.0064 | 0.0081 |
| TRIDEC | | | | | | |
| 2MeNAPH | | 0.0043 | 0.0066 | 0.0066 | 0.0080 | 0.0108 |
| PENTADEC | | | | | | |
| Total mass | 0.1177 | 0.0822 | 0.1339 | 0.1334 | 0.1773 | 0.1981 |

Notes:

Values in L/hr

Total mass does not include UNDEC, TRIDEC, PENTADEC (28 compounds)

TABLE B
4 CONSTANT REGRESSION OUTPUT

| | Adjusted Rsq | Standard Error | ln(SR0) | b | -Ea/R | d | Std Error ln(SR0) | Std Error b | Std Error - Ea/R | Std Error d |
|------------|-------------------------|---------------------------|----------------|----------|--------------|----------|----------------------------------|----------------------------|-------------------------------------|----------------------------|
| MTBE | 0.997 | 0.0960 | -3.217 | 0.981 | 2704 | -0.709 | 0.2881 | 0.0062 | 83 | 0.0082 |
| t12DCE | 0.992 | 0.1659 | -1.877 | 0.905 | 2147 | -0.760 | 0.4971 | 0.0100 | 144 | 0.0138 |
| 11DCA | 0.995 | 0.1272 | -1.346 | 0.916 | 1965 | -0.746 | 0.3809 | 0.0077 | 110 | 0.0106 |
| c12DCE | 0.995 | 0.1299 | -1.905 | 0.911 | 2137 | -0.751 | 0.3892 | 0.0078 | 112 | 0.0109 |
| CHCl3 | 0.996 | 0.1260 | -1.841 | 0.912 | 2118 | -0.748 | 0.3776 | 0.0076 | 109 | 0.0105 |
| 111TCA | 0.995 | 0.1279 | -2.684 | 0.902 | 2259 | -0.761 | 0.3836 | 0.0076 | 111 | 0.0106 |
| 12DCA | 0.995 | 0.1263 | -2.161 | 0.908 | 2218 | -0.746 | 0.3786 | 0.0076 | 109 | 0.0106 |
| BENZ | 0.995 | 0.1323 | -2.207 | 0.920 | 2198 | -0.754 | 0.3965 | 0.0080 | 114 | 0.0110 |
| CCl4 | 0.994 | 0.1405 | -3.121 | 0.889 | 2379 | -0.776 | 0.4220 | 0.0083 | 122 | 0.0116 |
| TCE | 0.992 | 0.1655 | -3.338 | 0.900 | 2522 | -0.772 | 0.4969 | 0.0099 | 144 | 0.0137 |
| 112TCA | 0.995 | 0.1264 | -2.412 | 0.896 | 2302 | -0.724 | 0.3790 | 0.0075 | 109 | 0.0107 |
| TOL | 0.994 | 0.1426 | -2.873 | 0.916 | 2364 | -0.756 | 0.4281 | 0.0087 | 124 | 0.0119 |
| OCT | 0.938 | 0.4698 | -5.984 | 0.822 | 3235 | -0.827 | 1.4231 | 0.0277 | 412 | 0.0388 |
| PCE | 0.991 | 0.1773 | -3.780 | 0.877 | 2601 | -0.775 | 0.5329 | 0.0103 | 154 | 0.0147 |
| CIBENZ | 0.994 | 0.1457 | -2.601 | 0.911 | 2292 | -0.747 | 0.4370 | 0.0088 | 126 | 0.0122 |
| 1112TetCA | 0.996 | 0.1235 | -2.676 | 0.898 | 2281 | -0.725 | 0.3705 | 0.0073 | 107 | 0.0104 |
| EtBENZ | 0.993 | 0.1597 | -2.930 | 0.918 | 2357 | -0.752 | 0.4794 | 0.0097 | 138 | 0.0134 |
| mpXYL | 0.992 | 0.1678 | -3.036 | 0.909 | 2372 | -0.749 | 0.5037 | 0.0101 | 145 | 0.0140 |
| oXYL | 0.993 | 0.1555 | -2.862 | 0.911 | 2312 | -0.740 | 0.4667 | 0.0094 | 135 | 0.0131 |
| 1122TetCA | 0.996 | 0.1118 | -1.971 | 0.913 | 2167 | -0.691 | 0.3351 | 0.0067 | 97 | 0.0096 |
| 135TMB | 0.988 | 0.2024 | -4.435 | 0.897 | 2720 | -0.738 | 0.6093 | 0.0121 | 176 | 0.0170 |
| 124TMB | 0.989 | 0.1997 | -4.126 | 0.890 | 2631 | -0.731 | 0.6009 | 0.0118 | 173 | 0.0169 |
| 13DCB | 0.991 | 0.1832 | -3.422 | 0.888 | 2449 | -0.730 | 0.5503 | 0.0108 | 159 | 0.0155 |
| 14DCB | 0.991 | 0.1802 | -3.263 | 0.892 | 2408 | -0.724 | 0.5413 | 0.0107 | 156 | 0.0153 |
| 12DCB | 0.992 | 0.1697 | -2.970 | 0.894 | 2327 | -0.716 | 0.5092 | 0.0101 | 147 | 0.0144 |
| UNDEC | 0.694 | 0.374 | -1.406 | 0.426 | 1708 | -0.806 | 1.792 | 0.028 | 517 | 0.053 |
| NAPH | 0.992 | 0.166 | -3.374 | 0.915 | 2430 | -0.671 | 0.497 | 0.010 | 144 | 0.014 |
| TRIDEC | | | | | | | | | | |
| 2MeNAPH | 0.984 | 0.238 | -5.498 | 0.869 | 2990 | -0.689 | 0.72 | 0.014 | 208 | 0.021 |
| PENTADEC | | | | | | | | | | |
| Total mass | 0.993 | 0.1543 | -6.111 | 0.907 | 2419 | -0.732 | 0.4666 | 0.0093 | 134 | 0.0130 |

TABLE C
8260C MASS UNCERTAINTY

**AGI 8260C Method for Mass using SPG-0008
Samplers**

| | 99% Uncertainty Range +/- | 95% Uncertainty Range +/- |
|-----------|---------------------------------|---------------------------------|
| MTBE | 20% | 14% |
| t12DCE | 22% | 15% |
| 11DCA | 18% | 12% |
| c12DCE | 18% | 12% |
| CHCl3 | 16% | 11% |
| 111TCA | 18% | 12% |
| 12DCA | 20% | 13% |
| BENZ | 16% | 10% |
| CCl4 | 19% | 12% |
| TCE | 15% | 10% |
| 112TCA | 18% | 12% |
| TOL | 15% | 10% |
| OCT | 20% | 13% |
| PCE | 16% | 11% |
| CIBENZ | 18% | 12% |
| 1112TetCA | 19% | 13% |
| EtBENZ | 18% | 12% |
| mpXYL | 18% | 12% |
| oXYL | 18% | 12% |
| 1122TetCA | 23% | 15% |
| 135TMB | 21% | 14% |
| 124TMB | 20% | 14% |
| 13DCB | 19% | 13% |
| 14DCB | 19% | 13% |
| 12DCB | 20% | 14% |
| NAPH | 21% | 14% |
| 2MeNAPH | 25% | 17% |

TABLE D
4 CONSTANT WATER CONCENTRATION UNCERTAINTY
ERROR IN CONCENTRATION REPORTING (1)

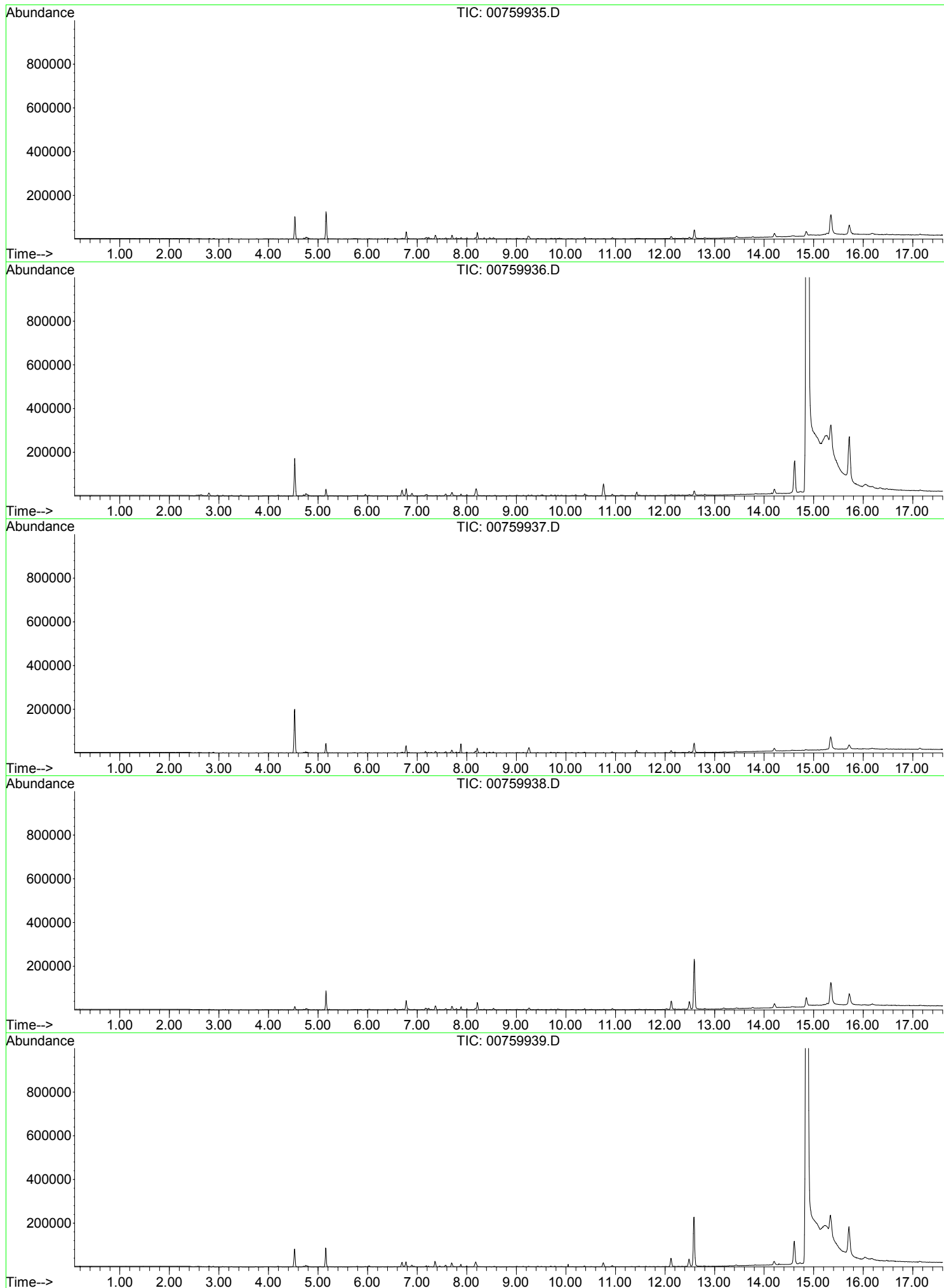
| | Average Error | Minimum Error | Maximum Error |
|-----------|--------------------------|--------------------------|--------------------------|
| MTBE | 6% | -12% | 12% |
| t12DCE | 11% | -26% | 21% |
| 11DCA | 8% | -19% | 13% |
| c12DCE | 9% | -19% | 15% |
| CHCl3 | 9% | -20% | 14% |
| 111TCA | 9% | -19% | 23% |
| 12DCA | 10% | -19% | 17% |
| BENZ | 8% | -18% | 13% |
| CCl4 | 10% | -23% | 22% |
| TCE | 10% | -21% | 14% |
| 112TCA | 11% | -21% | 21% |
| TOL | 7% | -17% | 14% |
| OCT | 20% | -41% | 42% |
| PCE | 10% | -24% | 15% |
| CIBENZ | 7% | -16% | 14% |
| 1112TetCA | 8% | -17% | 18% |
| EtBENZ | 6% | -19% | 14% |
| mpXYL | 7% | -22% | 13% |
| oXYL | 7% | -19% | 13% |
| 1122TetCA | 8% | -16% | 17% |
| 135TMB | 9% | -23% | 17% |
| 124TMB | 10% | -28% | 19% |
| 13DCB | 10% | -22% | 17% |
| 14DCB | 10% | -22% | 17% |
| 12DCB | 9% | -23% | 17% |
| NAPH | 10% | -24% | 21% |
| 2MeNAPH | 13% | -32% | 30% |

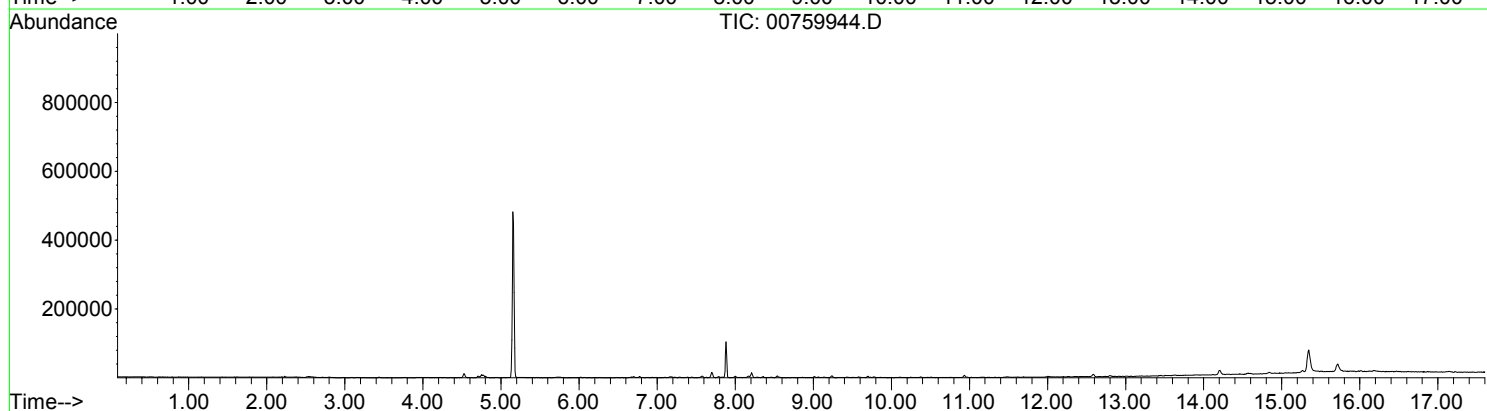
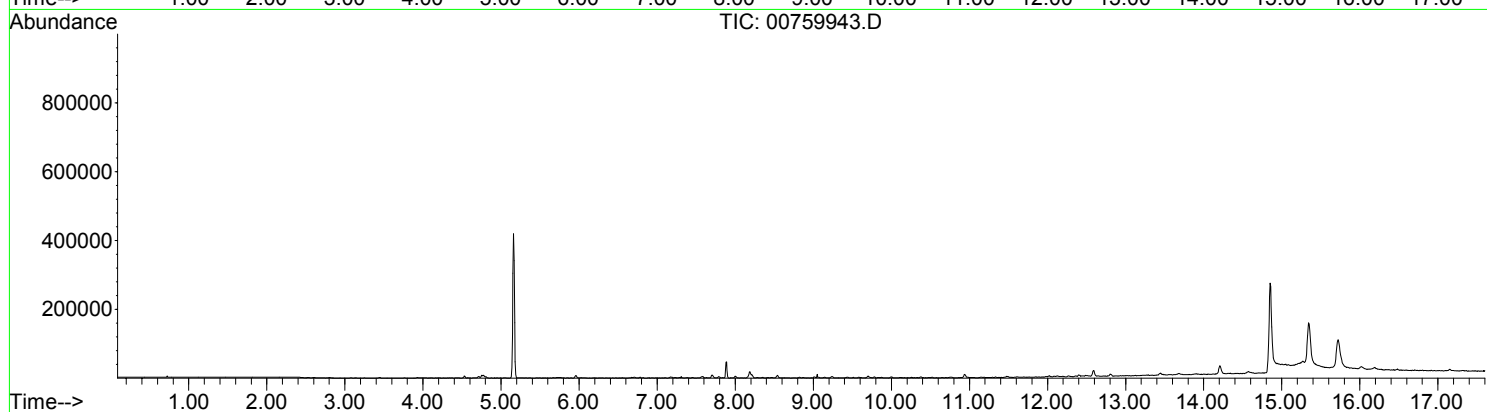
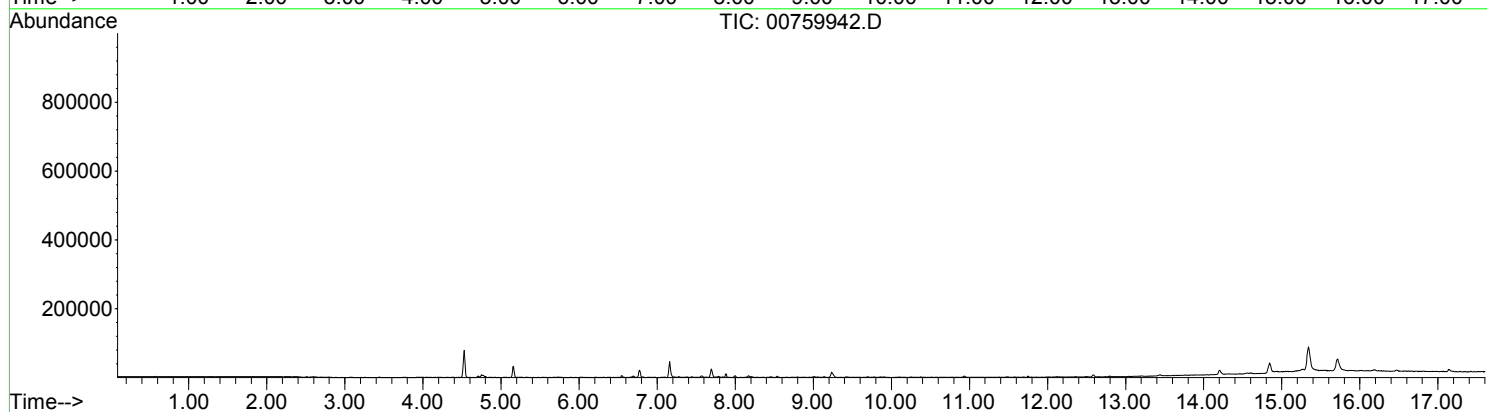
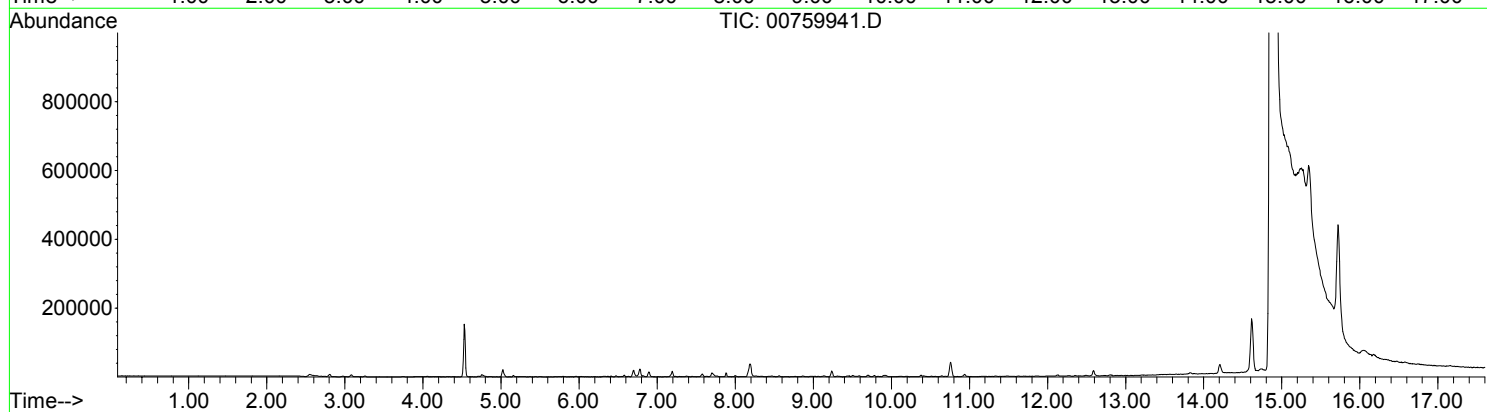
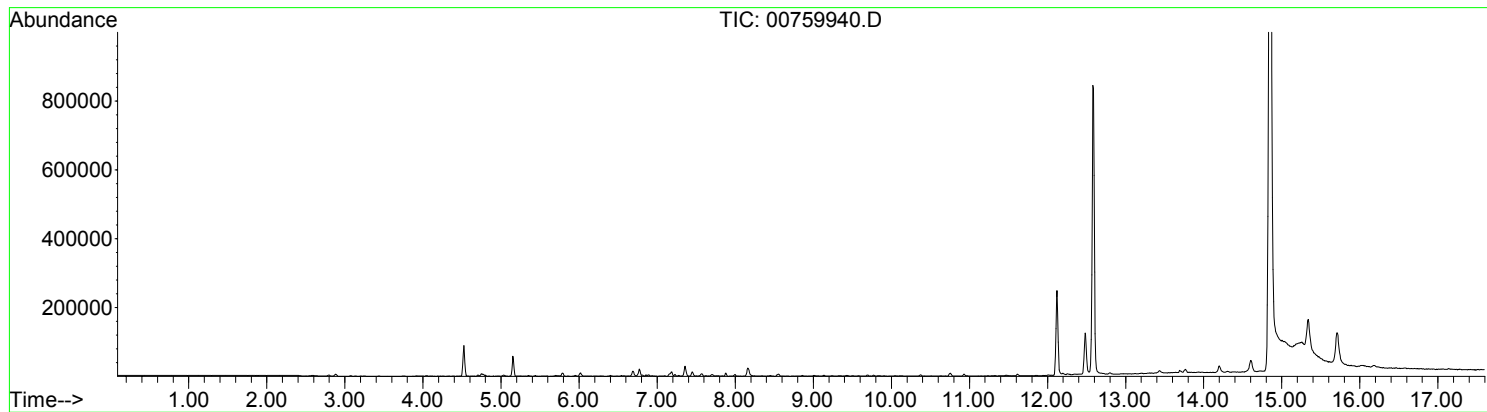
(1) For 1 hour exposure, includes error related to mass value from AGI analytical method 8260C

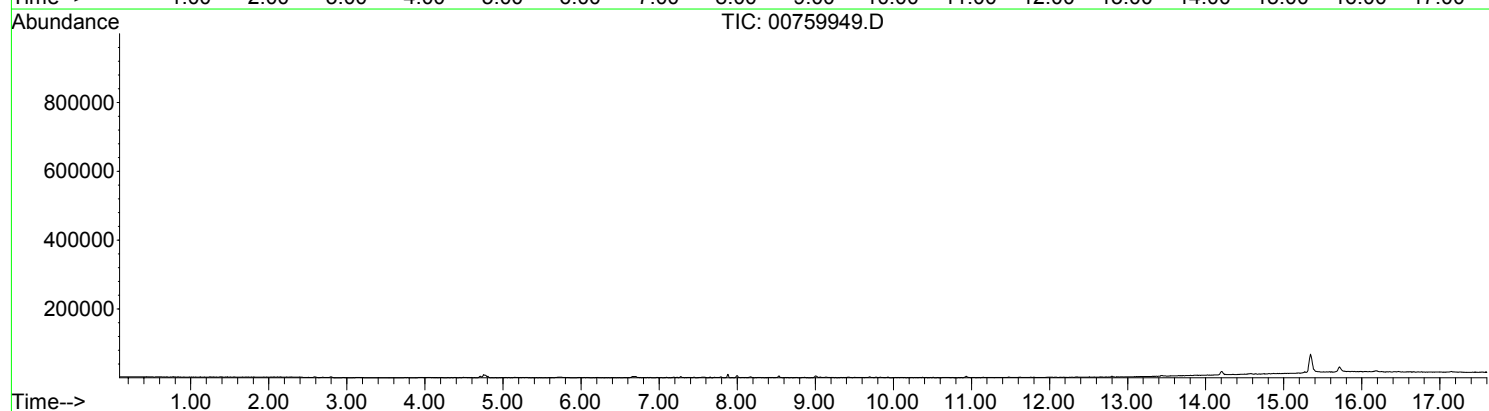
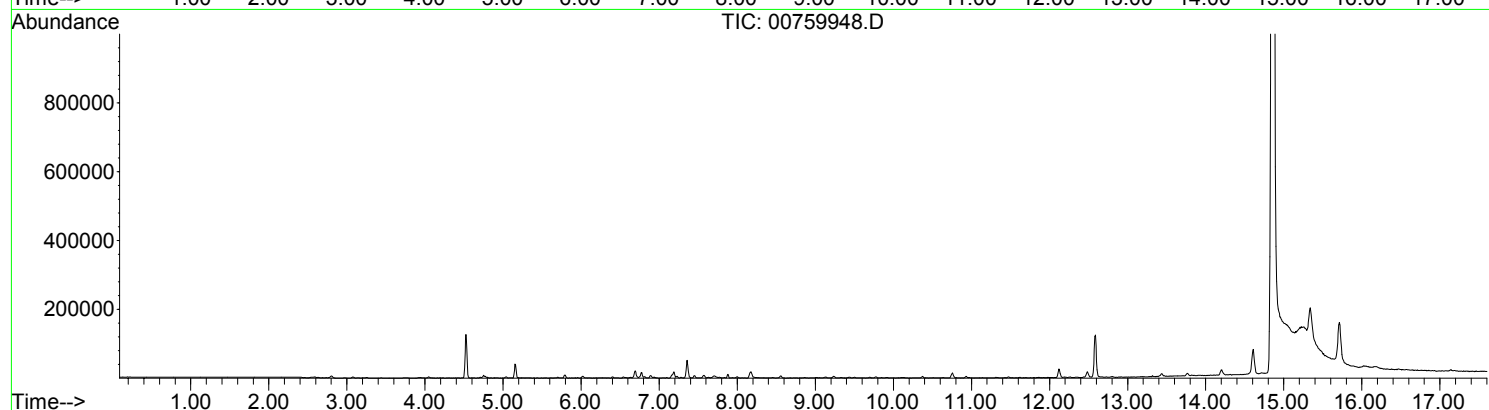
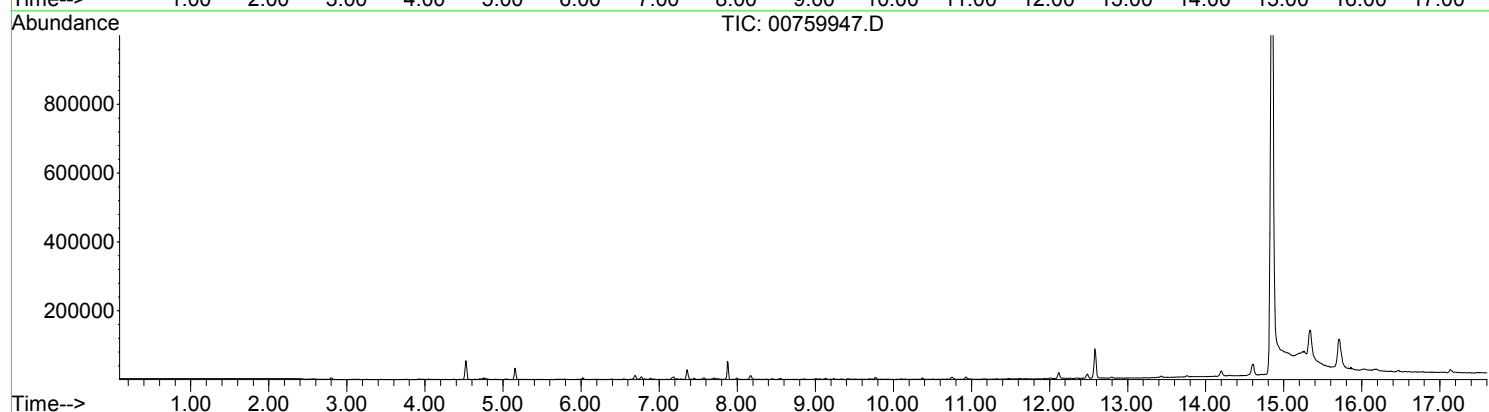
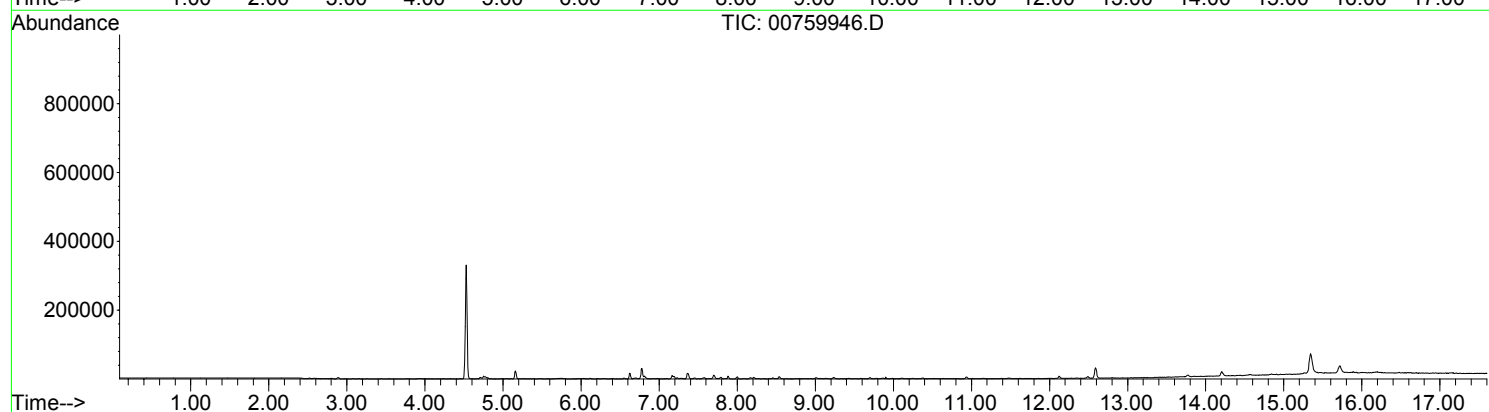
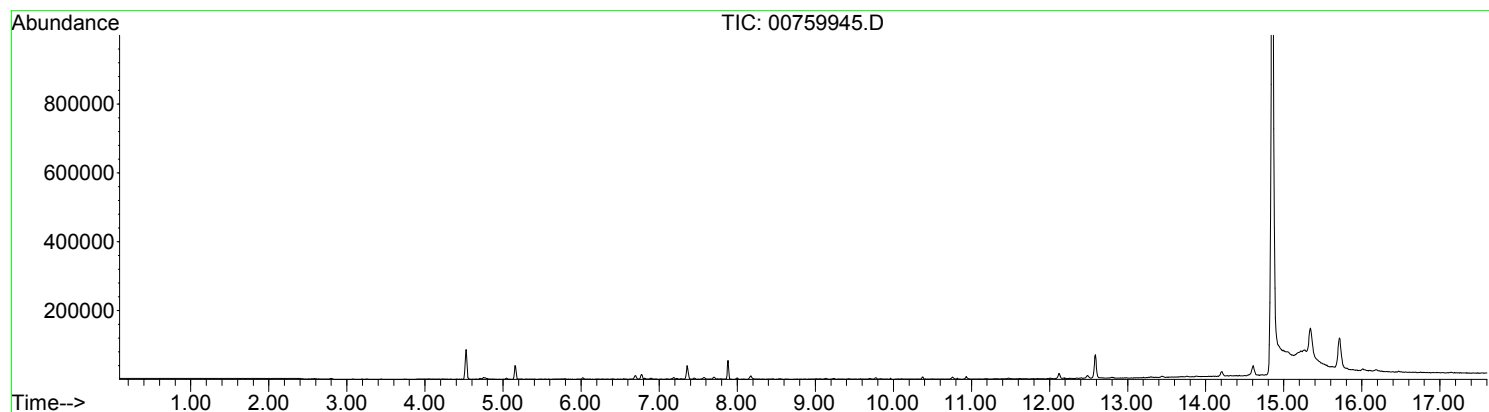
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AMPLIFIED
GEOCHEMICAL
IMAGING, LLC

Laboratory Report

Site: Comal & San Marcos Rivers
August Sampling

Prepared for:

SWCA Environmental Consultants
10245 Little York Road
Suite 600
San Antonio, TX
UNITED STATES

Prepared on:

September 10, 2015
(Revised September 15, 2015)

Project Summary and Objective

Amplified Geochemical Imaging, LLC. (AGI) provided the AGI Environmental Survey used at:

Comal & San Marcos Rivers August Sampling

The service provided by AGI included delivery of the required quantity of AGI Universal Samplers, analysis by the method described below for the requested organic compounds, reporting of the data, and contour mapping (as needed).

This report includes results for only the samples noted under the Laboratory Sample Report section. If contour maps are part of the project deliverable, the maps will be prepared and issued under a separate report cover, upon receipt of a usable sitemap (electronic) and compound choices for contouring.

Written/submitted by:

Jim E Whetzel

Project Manager

Reviewed/approved by:

Don D'Apolito

Project Manager

Analytical data approved by:

Ian McMullen

Chemist

Quality Assurance Statement

The AGI Laboratory, at Amplified Geochemical Imaging's facility in Newark, DE USA, operates under the guidelines of its ISO Standard 17025 DoD ELAP accreditation, and its Quality Assurance Manual, Operating Procedures, and Methods (SPG-SOP-0462).

For this project, the analytical method, results, and observations reported do [] do not [✓] fall within the scope of AGI's ISO 17025 accreditation.

Screening/Concentration Method

The AGI Universal Samplers are analyzed at AGI's fixed laboratory using thermal desorption-gas chromatography/mass spectrometry (TD-GC/MS) instrumentation following modified U.S. EPA Method 8260 (SPG-WI-0292) which includes the following:

- **BFB Tuning Frequency:** A BFB tune is analyzed at the start of each analytical run and after every 30 samples.
- **Initial Calibration:** A minimum of a five point calibration curve is analyzed prior to the analysis of samples.
- **Initial Calibration Verification (ICV):** Following the calibration a second-source reference standard is analyzed to verify the accuracy of the calibration. Acceptance criteria for the ICV is +/- 30%.
- **Linearity of Target Compounds:** If the RSD of any target analyte is less than or equal to 25% then average response factor can be used for quantitation. If the RSD exceeds 25% for a target compound a regression equation can be used for quantitation.
- **Continuing Calibration Verification:** After every 10 samples, and at the end of each analytical batch, a mid-level second-source Reference Standard is analyzed. The acceptance criteria for all target analytes in the reference standards are +/- 50% of the true value.
- **Method Blank:** Analyzed prior to the analysis of field samples and every 30 samples.

Note: Analyte levels reported for the field-deployed AGI Universal Samplers that exceed trip and method blank levels, and/or the reporting limit, are more likely to have originated from on-site sources.

| | |
|----------------------------|-------------------|
| Media Sampled: | WATER |
| Chemist - sample analysis: | Kelly J Stringham |
| Chemist - data processor: | Kelly J Stringham |
| Chemist - data review: | Ian McMullen |

Method Deviations: None

Please note that data file names ending with R are rerun samples using the second pair of sorbers, in which the original results were not reported. Data file names ending in D are duplicate analysis results for the second set of sorbers from the same sampler, and are reported.

Additional Report Information

- Comments
- Laboratory Sample Report
- Chain of Custody
- Installation and Retrieval Log
- Analytical Results and Key
- Concentration Calculation Method Summary
- Total Ion Chromatograms

Project Specific Comments

Due to an instrument malfunction, samples were analyzed on two different instruments. The second instrument had an elevated TPH background, therefore the TPH values for samples analyzed on the second instrument were found to have artificially higher TPH values reported. Samples analyzed on the second instrument included: 767203, -204, -205, -210, and -221. Relative impact of the elevated TPH background can be seen in sample 767221, which was an unused sampler analyzed on the second instrument and was found to have 0.65ug of TPH.

This report was revised at the request of SWCA to report only the trip blank designated on the chain of custody, 00767214.

Survey period ¹

Samplers were installed on August 13, 2015 and retrieved on August 27, 2015 for an exposure time of 14 days.

Tamper seal intact:

Yes

Date received:

8/28/15 11:00 am

By: Scott Kirlin

COC returned:

Yes

Comments:

None.

¹ - Installation start to end of retrieval, as reported. See installation and retrieval log for individual deployment and retrieval dates and times (i.e., sampler exposure time).

General Comments

Analytical QA/QC

Laboratory instrumentation consists of gas chromatographs equipped with mass selective detectors, coupled with automated thermal desorption units. Sample preparation involves cutting the tip off the bottom of the AGI Universal Sampler, and transferring one or more "sorbers" to a thermal desorption tube for analysis. The insertion/retrieval cord prevents soil, water and other interferences from coming in contact with the adsorbent. No further sample preparation is required. Any replicate sorbers not consumed in the initial analysis will be discarded fifteen (15) days from the date of the laboratory report.

Data are archived and stored in a secure manner as per AGI's Quality Assurance program (SPG-SOP-0462).

Total petroleum hydrocarbons (TPH), gasoline-range petroleum hydrocarbons (GRPH), and/or diesel range petroleum hydrocarbons (DRPH), when reported, are calculated using the area under the peaks observed in m/z 55 and 57 selected ion chromatograms. Quantitation of the mass values was performed using the response factor for a specific alkane (present in the calibration standards). TPH values include the entire chromatogram and provide estimates for aliphatic hydrocarbon ranges of C4 to C20. GRPH and DRPH include only the relevant regions of the chromatograms and provide estimates for C4 to C10 and C10 to C20 aliphatic hydrocarbons, respectively.

Trip blanks were provided to document potential exposures that were not part of the signal of interest (e.g., impact during sampler shipment, installation and/or retrieval, and storage). The trip blanks are identically manufactured and packaged AGI Universal Samplers to those samplers deployed in the field. The trip blanks remain unopened during all phases of the project. Levels reported on the trip blanks may indicate potential impact to the samplers other than the contaminant source of interest.

Unresolved peak envelopes (UPEs) are represented as a series of compound peaks clustered together around a central gas chromatograph elution time in the total ion chromatogram. UPEs may be indicative of complex fluid mixtures. UPEs observed early in the chromatograms are considered to indicate presence of more volatile fluids, while UPEs observed later in the chromatogram may indicate the presence of less volatile fluids. Multiple UPEs may indicate the presence of multiple complex fluids.

Total ion chromatograms (TICs) are included in the Attachments. The eight-digit serial number of each sampler is incorporated in the TIC identification (e.g., 12345678.D represents AGI Universal Sampler 12345678).

General Comments

Soil Gas Sampling

For soil gas sampling, the AGI Environmental Survey reports mass levels migrating through the open pore spaces of the soil and diffusing through the sampler membrane for sorption by the engineered, hydrophobic adsorbents, housed within the membrane tube. During the migration of the soil gas away from the source to the AGI Universal Sampler, the vapors are subject to a variety of attenuation factors. The soil gas masses reported on the samplers compare favorably with the concentrations reported in the soil or groundwater (e.g., where soil gas levels are reported at greater levels to other sampled locations on the site, the matrix data should reveal the same pattern, and vice versa). However, due to a variety of factors, a perfect comparison between matrix data and soil gas levels can rarely be achieved.

Soil gas concentrations ($\mu\text{g}/\text{m}^3$) are calculated following the method described in the Additional Report Information section.

Soil gas signals reported by this method cannot be correlated specifically to soil adsorbed, groundwater, and /or free-phase contamination. The soil gas signal reported from each AGI Universal Sampler can evolve from all of these sources. Differentiation between soil and groundwater contamination can only be achieved with prior knowledge of the site history (i.e., the site is known to have groundwater contamination only).

Air Sampling

For indoor, outdoor, and crawlspace air sampling, the AGI Environmental Survey reports mass levels present in the air and diffusing through the sampler membrane for sorption by the engineered adsorbents housed within the membrane tube.

Air concentrations ($\mu\text{g}/\text{m}^3$) are calculated following the method described in the Additional Report Information section.

Groundwater and Sediment Porewater Sampling

For groundwater and sediment porewater sampling, the AGI Environmental Survey reports the mass levels of compounds present in the water which, when coming in contact with the sampler membrane, partitions out of solution, and diffuses through the sampler membrane for sorption by the engineered adsorbents.

Water concentrations ($\mu\text{g}/\text{L}$) are calculated using the quantified mass, exposure period and the compound specific uptake rate. The rates were measured under controlled experimental conditions. The uptake rates are corrected for water pressure (depth of the AGI Universal Sampler below the water table), water temperature and the aquifer flow rate. For sediment porewater, the uptake rate is corrected for the reduced volume of water in the sediment, by multiplying the uptake rate by the pore water fraction.

LABORATORY SAMPLE REPORT

Project: ENV 01464

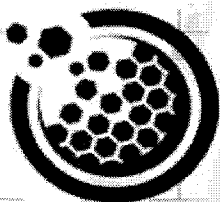
Site Name: Comal & San Marcos Rivers Aug.

Module Type: SPG0008

| Module ID | Sample Type | Field ID | |
|----------------------------|--------------------------|---------------------|-------------------|
| 00767200 | FIELD_SAMPLE | HCS430 | |
| 00767201 | FIELD_SAMPLE | HCS410 | |
| 00767202 | FIELD_SAMPLE | HCS420 | |
| 00767203 | FIELD_SAMPLE | HCS440 | |
| 00767204 | FIELD_SAMPLE | FDHCS440 | |
| 00767205 | FIELD_SAMPLE | HCS460 | |
| 00767206 | FIELD_SAMPLE | HSM410 | |
| 00767207 | FIELD_SAMPLE | HSM420 | |
| 00767208 | FIELD_SAMPLE | HSM430 | |
| 00767209 | FIELD_SAMPLE | FDHSM430 | |
| 00767210 | FIELD_SAMPLE | HSM440 | |
| 00767212 | FIELD_SAMPLE | HSM460 | |
| 00767213 | FIELD_SAMPLE | HSM470 | |
| 00767214 | TRIP_BLANK | TB10 | |
| 00767221 | UNUSED | NOT USED | |
| Total # "FIELD SAMPLES" | Total # "TRIP BLANKS" | Total # "UNUSED" | Total # "LOST" |
| 13 | 1 | 1 | 0 |

Duplicate samples: 0

Note: e o re r ed o e er e re e o W re ere o re or ed



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210 Executive Drive, Suite 1
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AGI Universal Passive Sampler Chain of Groundwater Sampling

Production Order #: **01464**

Customer Name: SWCA Environmental Consultants
Address: 6200 UTSA Boulevard
Suite 102
San Antonio, TX 78249
USA

Site Name: Comal & San Marcos Rivers Aug.
Site Address:
Project Manager:

| | | | | |
|------------------------------|--------------------------------|-------|------------------|---|
| Serial # of Samplers Shipped | # of Samplers for Installation | 14.00 | # of Trip Blanks | 2 |
| 00767200 - 00767214 | Total Samplers Shipped | 16.00 | Pieces | |
| 00767221 - 00767221 | Total Samplers Received | 16 | Pieces | |
| | Total Samplers Installed | 14 | Pieces | |

Serial # of Trip Blanks (Client Decides)

00767214

Please hold analysis of
00767211 until status
verified w/ our client

| | |
|--|---|
| Prepared By: <u>[Signature]</u> | Is Concurrent water sampling planned this monitoring period? YES <input type="radio"/> NO <input checked="" type="radio"/> |
| Verified By: <u>[Signature]</u> 7/22/15 | Scheduled Sampling Date: _____ |
| Installation Performed By: Name: <u>Jennifer Moreland</u> ^{gm} Guy <u>Rabin</u> Company: <u>SWCA</u> | Retrieval Performed By: Name: <u>Jennifer Moreland</u> Guy <u>Brittany</u> <u>Rabin</u> Company: <u>SWCA</u> |
| Installation Start Date / Time: <u>8/13/15 10:40</u> | Retrieval Start Date / Time: <u>8/27/15 948</u> |
| Installation Complete Date / Time: <u>8/13/15 14:38</u> | Retrieval Complete Date / Time: <u>8/27/15 1354</u> |
| Total Samplers Retrieved: <u>14</u> | |
| Total Samplers Lost In Field: <u>0</u> | |
| Total Unused Samplers Returned: <u>1</u> | |
| Relinquished By: <u>[Signature]</u> Date/Time: <u>7-22-15</u> Company: <u>AGI</u> 2:45PM | Received By: <u>[Signature]</u> Date/Time: <u>7/23/15</u> Company: <u>SWCA</u> 930AM |
| Relinquished By: <u>[Signature]</u> Date/Time: <u>8/27/15</u> Company: <u>SWCA</u> 16:51 | Received By: <u>[Signature]</u> Date/Time: <u>8-28-17</u> Company: <u>AGI</u> 11:00AM |



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IMAGING, LLC**

210 Executive Drive, Suite 1
Newark, DE USA 19702-3335
ph: 302-266-2428

AGI Project No.

ENV 01464

Site Name:

Comal & San Marcos Rivers Aug.

Site Location:

New Braunfels and San Marcos, TX

AGI Soil Gas Sampling

Installation & Retrieval Log

Company Name:

SWCA Environmental Consultants

Location:

San Antonio, TX

Samples collected by:

Jennifer Moreland, Guy Rubio, Brittany Rios

* Optional or as needed

| SAMPLER SERIAL NO. | WELL ID | SAMPLE TYPE (Field Sample, Trip Blank, Field Blank, etc.) | INSTALLATION DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 13:00 | RETRIEVAL DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 15:00 | WATER QUALITY MONITORING | | | |
|--------------------|----------|---|--|---|--|--|---|---|
| | | | | | INSTALLATION DEPTH FROM TOP OF CASING (feet) | DEPTH TO WATER FROM TOP OF CASING (feet) | SCREENED INTERVAL FROM TOP OF CASING (feet) | WATER TEMPERATURE AT MODULE DEPTH (deg C) |
| 00767200 | HCS430 | FIELD_SAMPLE | 8/13/15 10:40 | 8/27/15 12:55 | 3.25 | | | |
| 00767201 | HCS410 | FIELD_SAMPLE | 8/13/15 11:01 | 8/27/15 13:16 | 1.1 | | | |
| 00767202 | HCS420 | FIELD_SAMPLE | 8/13/15 11:08 | 8/27/15 13:25 | 2.1 | | | |
| 00767203 | HCS440 | FIELD_SAMPLE | 8/13/15 11:28 | 8/27/15 13:45 | 1.9 | | | |
| 00767204 | FDHCS440 | FIELD_SAMPLE | 8/13/15 11:28 | 8/27/15 13:45 | 1.9 | | | |
| 00767205 | HCS460 | FIELD_SAMPLE | 8/13/15 11:36 | 8/27/15 13:54 | 2.65 | | | |
| 00767206 | HSM410 | FIELD_SAMPLE | 8/13/15 13:08 | 8/27/15 9:48 | 1.5 | | | |
| 00767207 | HSM420 | FIELD_SAMPLE | 8/13/15 13:21 | 8/27/15 10:00 | 2.9 | | | |
| 00767208 | HSM430 | FIELD_SAMPLE | 8/13/15 13:27 | 8/27/15 10:08 | 0.55 | | | |
| 00767209 | FDHSM430 | FIELD_SAMPLE | 8/13/15 13:27 | 8/27/15 10:08 | 0.55 | | | |
| 00767210 | HSM440 | FIELD_SAMPLE | 8/13/15 13:39 | 8/27/15 10:19 | 5 | | | |
| 00767211 | HSM450 | FIELD_SAMPLE | 8/13/15 13:56 | 8/27/15 10:51 | 2.1 | | | |
| 00767212 | HSM460 | FIELD_SAMPLE | 8/13/15 14:25 | 8/27/15 11:06 | 0.8 | | | |
| 00767213 | HSM470 | FIELD_SAMPLE | 8/13/15 14:38 | 8/27/15 11:20 | 1.8 | | | |
| 00767214 | TB10 | TRIP_BLANK | 8/13/15 10:40 | 8/27/15 13:54 | NA | | | |
| 00767221 | NOT USED | TRIP_BLANK | NA | NA | NA | | | |



**AGI Soil Gas Sampling
Installation & Retrieval L**

* Optional or as needed

| SAMPLER SERIAL NO. | FLOW RATE THROUGH SCREEN: HIGH OR LOW (HIGH = flow > than 10 m/day; LOW = flow < 10 m/day) | OBSERVATIONS/COMMENTS* (e.g., observations, location description, missing, pulled from well, etc. - as needed) | YES / NO | |
|--------------------|--|--|--|--------|
| | | | EVIDENCE OF LIQUID PETROLEUM HYDROCARBONS? | ODOR ? |
| 00767200 | | | | |
| 00767201 | | | | |
| 00767202 | | | | |
| 00767203 | | | | |
| 00767204 | | | | |
| 00767205 | | | | |
| 00767206 | | | | |
| 00767207 | | | | |
| 00767208 | | | | |
| 00767209 | | | | |
| 00767210 | | | | |
| 00767211 | | PLEASE DO NOT ANALYZE THIS SAMPLER. IT WAS TAMPERED WITH IN THE FIELD AND REMOVED FROM THE WATER. | | |
| 00767212 | | | | |
| 00767213 | | | | |
| 00767214 | | | | |
| 00767221 | | | | |



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ph: +1-302-266-2428
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PROJECT NUMBER: ENV 01464
SITE NAME: Comal & San Marcos Rivers Aug.
SITE ADDRESS:

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00767200 FIELD_SAMPLE

Dilution Factor: 1 Field ID: HCS430

Installation Date: 8/13/2015 10:40:00AM

Retrieval Date: 8/27/2015 12:55:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Reviewer: Ian McMullen

Matrix: WATER

Product: SPG0008

Date Analyzed: 9/1/2015 2:19:00PM

Batch: ENV-150831-3

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.36 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767200 **FIELD_SAMPLE**

Dilution Factor: 1

Field ID: HCS430

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 10:40:00AM

Retrieval Date: 8/27/2015 12:55:00PM

Date Analyzed: 9/1/2015 2:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767201 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HCS410

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 11:01:00AM

Retrieval Date: 8/27/2015 1:16:00PM

Date Analyzed: 9/1/2015 3:50:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.09 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767201 **FIELD_SAMPLE**

Dilution Factor: 1

Field ID: HCS410

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 11:01:00AM

Retrieval Date: 8/27/2015 1:16:00PM

Date Analyzed: 9/1/2015 3:50:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767202 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS420

Installation Date: 8/13/2015 11:08:00AM

Retrieval Date: 8/27/2015 1:25:00PM

Date Analyzed: 9/1/2015 2:50:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.13 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767202 **FIELD_SAMPLE**

Dilution Factor: 1

Field ID: HCS420

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 11:08:00AM

Retrieval Date: 8/27/2015 1:25:00PM

Date Analyzed: 9/1/2015 2:50:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767203R FIELD_SAMPLE

Dilution Factor: 1

Field ID: HCS440

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 11:28:00AM

Retrieval Date: 8/27/2015 1:45:00PM

Date Analyzed: 9/3/2015 2:35:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150902-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.39 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767203R FIELD_SAMPLE

Dilution Factor: 1

Field ID: HCS440

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 11:28:00AM

Retrieval Date: 8/27/2015 1:45:00PM

Date Analyzed: 9/3/2015 2:35:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150902-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 1.23 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767204R FIELD_SAMPLE

Dilution Factor: 1

Field ID: FDHCS440

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 11:28:00AM

Retrieval Date: 8/27/2015 1:45:00PM

Date Analyzed: 9/3/2015 1:35:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150902-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.44 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767204R FIELD_SAMPLE

Dilution Factor: 1

Field ID: FDHCS440

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 11:28:00AM

Retrieval Date: 8/27/2015 1:45:00PM

Date Analyzed: 9/3/2015 1:35:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150902-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 1.44 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767205 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HCS460

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 11:36:00AM

Retrieval Date: 8/27/2015 1:54:00PM

Date Analyzed: 9/3/2015 11:04:00AM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150902-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|--------------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.31 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | 0.08 | 0.02 |
| o-Xylene | 95-47-6 | 0.04 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767205 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HCS460

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 11:36:00AM

Retrieval Date: 8/27/2015 1:54:00PM

Date Analyzed: 9/3/2015 11:04:00AM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150902-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 2.59 | 0.50 |
| BTEX | | 0.11 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767206 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM410

Installation Date: 8/13/2015 1:08:00PM

Retrieval Date: 8/27/2015 9:48:00AM

Date Analyzed: 9/1/2015 11:19:00AM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767206 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM410

Installation Date: 8/13/2015 1:08:00PM

Retrieval Date: 8/27/2015 9:48:00AM

Date Analyzed: 9/1/2015 11:19:00AM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767207 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM420

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 1:21:00PM

Retrieval Date: 8/27/2015 10:00:00AM

Date Analyzed: 9/1/2015 1:49:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.06 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767207 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM420

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 1:21:00PM

Retrieval Date: 8/27/2015 10:00:00AM

Date Analyzed: 9/1/2015 1:49:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 0.67 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767208 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM430

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 1:27:00PM

Retrieval Date: 8/27/2015 10:08:00AM

Date Analyzed: 9/1/2015 3:20:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.68 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767208 **FIELD_SAMPLE**

Dilution Factor: 1

Field ID: HSM430

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 1:27:00PM

Retrieval Date: 8/27/2015 10:08:00AM

Date Analyzed: 9/1/2015 3:20:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767209 FIELD_SAMPLE

Dilution Factor: 1

Field ID: FDHSM430

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 1:27:00PM

Retrieval Date: 8/27/2015 10:08:00AM

Date Analyzed: 9/1/2015 12:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.89 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767209 **FIELD_SAMPLE**

Dilution Factor: 1

Field ID: FDHSM430

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 1:27:00PM

Retrieval Date: 8/27/2015 10:08:00AM

Date Analyzed: 9/1/2015 12:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767210 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM440

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 1:39:00PM

Retrieval Date: 8/27/2015 10:19:00AM

Date Analyzed: 9/3/2015 12:34:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150902-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.11 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



AMPLIFIED GEOCHEMICAL IMAGING, LLC

210 Executive Drive, Suite 1
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www.agisurveys.net

PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767210 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM440

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 1:39:00PM

Retrieval Date: 8/27/2015 10:19:00AM

Date Analyzed: 9/3/2015 12:34:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150902-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | 1.18 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767212 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM460

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 2:25:00PM

Retrieval Date: 8/27/2015 11:06:00AM

Date Analyzed: 9/1/2015 11:49:00AM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.05 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767212 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM460

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 2:25:00PM

Retrieval Date: 8/27/2015 11:06:00AM

Date Analyzed: 9/1/2015 11:49:00AM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767213 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM470

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 2:38:00PM

Retrieval Date: 8/27/2015 11:20:00AM

Date Analyzed: 9/1/2015 1:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.05 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767213 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM470

Installation Date: 8/13/2015 2:38:00PM

Retrieval Date: 8/27/2015 11:20:00AM

Date Analyzed: 9/1/2015 1:19:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767214 TRIP_BLANK

Dilution Factor: 1

Field ID: TB10

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 10:40:00AM

Retrieval Date: 8/27/2015 1:54:00PM

Date Analyzed: 9/1/2015 12:49:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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www.agisurveys.net

PROJECT NUMBER: ENV 01464

SITE NAME: Comal & San Marcos Rivers Aug.

SITE ADDRESS:

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00767214 TRIP_BLANK

Dilution Factor: 1

Field ID: TB10

Matrix: WATER

Product: SPG0008

Installation Date: 8/13/2015 10:40:00AM

Retrieval Date: 8/27/2015 1:54:00PM

Date Analyzed: 9/1/2015 12:49:00PM

Analyst: Kelly J Stringham

Method: SPG-WI-0292

Batch: ENV-150831-3

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |

AMPLIFIED GEOCHEMICAL IMAGING, LLC
210 EXECUTIVE DRIVE, SUITE 1, NEWARK, DE 19702
SWCA ENVIRONMENTAL CONSULTANTS, SAN ANTONIO, TX
STANDARD TARGET VOCs/SVOCs
ESTIMATED WATER CONCENTRATIONS
COMAL & SAN MARCOS RIVERS
ORDER # 01464

| | | | | | | | | | estimated | | |
|----------------|--------------|------------|------------|------------|----|------------|----|----|-----------|------------|--------------|
| DATAFILE | FIELD | DATE/ TIME | DATE/ TIME | DATE/ TIME | | DATE/ TIME | | | | | |
| NAME | ID | INSTALLED | RETRIEVED | RECEIVED | | ANALYZED | | DF | TPH, ug/L | MTBE, ug/L | t12DCE, ug/L |
| Average RL = | | | | | | | | | 0.05 | 0.01 | 0.01 |
| 00767200 | HCS430 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/1/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |
| 00767201 | HCS410 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/1/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |
| 00767202 | HCS420 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/1/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |
| 00767203R | HCS440 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/3/2015 | ET | 1 | 0.08 | <0.01 | <0.01 |
| 00767204R | FDHCS440 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/3/2015 | ET | 1 | 0.08 | <0.01 | <0.01 |
| 00767205 | HCS460 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/3/2015 | ET | 1 | 0.11 | <0.01 | <0.01 |
| 00767206 | HSM410 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/1/2015 | ET | 1 | <0.06 | <0.01 | <0.01 |
| 00767207 | HSM420 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/1/2015 | ET | 1 | 0.06 | <0.01 | <0.01 |
| 00767208 | HSM430 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/1/2015 | ET | 1 | <0.06 | <0.01 | <0.01 |
| 00767209 | FDHSM430 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/1/2015 | ET | 1 | <0.06 | <0.01 | <0.01 |
| 00767210 | HSM440 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/3/2015 | ET | 1 | 0.08 | <0.01 | <0.01 |
| 00767212 | HSM460 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/1/2015 | ET | 1 | <0.06 | <0.01 | <0.01 |
| 00767213 | HSM470 | 8/13/2015 | 8/27/2015 | 8/28/2015 | ET | 9/1/2015 | ET | 1 | <0.06 | <0.01 | <0.01 |
| | | | | | | | | | | | |
| 00767214 | TB10 | | | 8/28/2015 | ET | 9/1/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |
| | | | | | | | | | | | |
| BLK_ENV-268217 | Method Blank | | | | | 9/1/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |
| BLK_ENV-268571 | Method Blank | | | | | 9/3/2015 | ET | 1 | <0.05 | <0.01 | <0.01 |

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ESTIMATED WATER CONCENTRATIONS
COMAL & SAN MARCOS RIVERS
ORDER # 01464

| DATAFILE | FIELD | | | | | | | | |
|----------------|--------------|-------------|--------------|-------------|--------------|-------------|------------|------------|-----------|
| NAME | ID | 11DCA, ug/L | c12DCE, ug/L | CHCl3, ug/L | 111TCA, ug/L | 12DCA, ug/L | BENZ, ug/L | CCl4, ug/L | TCE, ug/L |
| Average RL = | | 0.01 | 0.01 | 0.01 | 0.00 | 0.01 | 0.01 | 0.00 | 0.01 |
| 00767200 | HCS430 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 |
| 00767201 | HCS410 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 |
| 00767202 | HCS420 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 |
| 00767203R | HCS440 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 |
| 00767204R | FDHCS440 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 |
| 00767205 | HCS460 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 |
| 00767206 | HSM410 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 |
| 00767207 | HSM420 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 |
| 00767208 | HSM430 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 |
| 00767209 | FDHSM430 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 |
| 00767210 | HSM440 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 |
| 00767212 | HSM460 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 |
| 00767213 | HSM470 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 |
| | | | | | | | | | |
| 00767214 | TB10 | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 |
| | | | | | | | | | |
| BLK_ENV-268217 | Method Blank | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 |
| BLK_ENV-268571 | Method Blank | <0.01 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.00 | <0.01 |

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| DATAFILE | FIELD | | | | | | | | |
|----------------|--------------|--------------|-----------|-----------|-----------|--------------|-----------------|--------------|-------------|
| NAME | ID | 112TCA, ug/L | TOL, ug/L | OCT, ug/L | PCE, ug/L | CIBENZ, ug/L | 1112TetCA, ug/L | ETBENZ, ug/L | mpXYL, ug/L |
| Average RL = | | 0.01 | 0.01 | 0.00 | 0.01 | 0.01 | 0.01 | 0.01 | 0.01 |
| 00767200 | HCS430 | <0.01 | <0.01 | <0.00 | 0.07 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767201 | HCS410 | <0.01 | <0.01 | <0.00 | 0.02 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767202 | HCS420 | <0.01 | <0.01 | <0.00 | 0.03 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767203R | HCS440 | <0.01 | <0.01 | <0.00 | 0.07 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767204R | FDHCS440 | <0.01 | <0.01 | <0.00 | 0.08 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767205 | HCS460 | <0.01 | <0.01 | <0.00 | 0.06 | <0.01 | <0.01 | <0.01 | 0.02 |
| 00767206 | HSM410 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767207 | HSM420 | <0.01 | <0.01 | <0.00 | 0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767208 | HSM430 | <0.01 | <0.01 | <0.00 | 0.12 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767209 | FDHSM430 | <0.01 | <0.01 | <0.00 | 0.15 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767210 | HSM440 | <0.01 | <0.01 | <0.00 | 0.03 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767212 | HSM460 | <0.01 | <0.01 | <0.00 | 0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767213 | HSM470 | <0.01 | <0.01 | <0.00 | 0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| | | | | | | | | | |
| 00767214 | TB10 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| | | | | | | | | | |
| BLK_ENV-268217 | Method Blank | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| BLK_ENV-268571 | Method Blank | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |

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| DATAFILE | FIELD | | | | | | | |
|----------------|--------------|------------|-----------------|--------------|--------------|-------------|-------------|-------------|
| NAME | ID | oXYL, ug/L | 1122TetCA, ug/L | 135TMB, ug/L | 124TMB, ug/L | 13DCB, ug/L | 14DCB, ug/L | 12DCB, ug/L |
| Average RL = | | 0.01 | 0.01 | 0.00 | 0.01 | 0.01 | 0.01 | 0.01 |
| 00767200 | HCS430 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767201 | HCS410 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767202 | HCS420 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767203R | HCS440 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767204R | FDHCS440 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767205 | HCS460 | 0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767206 | HSM410 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767207 | HSM420 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767208 | HSM430 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767209 | FDHSM430 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767210 | HSM440 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767212 | HSM460 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| 00767213 | HSM470 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| | | | | | | | | |
| 00767214 | TB10 | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| | | | | | | | | |
| BLK_ENV-268217 | Method Blank | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |
| BLK_ENV-268571 | Method Blank | <0.01 | <0.01 | <0.00 | <0.01 | <0.01 | <0.01 | <0.01 |

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ORDER # 01464

| | | | | estimated | | estimated | estimated |
|----------------|--------------|-------------|------------|--------------|---------------|----------------------|----------------|
| DATAFILE | FIELD | | | | | | |
| NAME | ID | UNDEC, ug/L | NAPH, ug/L | TRIDEC, ug/L | 2MeNAPH, ug/L | Acenaphthylene, ug/L | PENTADEC, ug/L |
| Average RL = | | 0.02 | 0.02 | 0.02 | 0.01 | 0.01 | 0.02 |
| 00767200 | HCS430 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| 00767201 | HCS410 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| 00767202 | HCS420 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| 00767203R | HCS440 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| 00767204R | FDHCS440 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| 00767205 | HCS460 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| 00767206 | HSM410 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| 00767207 | HSM420 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| 00767208 | HSM430 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| 00767209 | FDHSM430 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| 00767210 | HSM440 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| 00767212 | HSM460 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| 00767213 | HSM470 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| | | | | | | | |
| 00767214 | TB10 | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| | | | | | | | |
| BLK_ENV-268217 | Method Blank | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |
| BLK_ENV-268571 | Method Blank | <0.02 | <0.02 | <0.02 | <0.01 | <0.01 | <0.02 |

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| | | estimated | estimated |
|----------------|--------------|--------------------|----------------|
| DATAFILE | FIELD | | |
| NAME | ID | Acenaphthene, ug/L | Fluorene, ug/L |
| Average RL = | | 0.01 | 0.01 |
| 00767200 | HCS430 | <0.01 | <0.01 |
| 00767201 | HCS410 | <0.01 | <0.01 |
| 00767202 | HCS420 | <0.01 | <0.01 |
| 00767203R | HCS440 | <0.01 | <0.01 |
| 00767204R | FDHCS440 | <0.01 | <0.01 |
| 00767205 | HCS460 | <0.01 | <0.01 |
| 00767206 | HSM410 | <0.01 | <0.01 |
| 00767207 | HSM420 | <0.01 | <0.01 |
| 00767208 | HSM430 | <0.01 | <0.01 |
| 00767209 | FDHSM430 | <0.01 | <0.01 |
| 00767210 | HSM440 | <0.01 | <0.01 |
| 00767212 | HSM460 | <0.01 | <0.01 |
| 00767213 | HSM470 | <0.01 | <0.01 |
| | | | |
| 00767214 | TB10 | <0.01 | <0.01 |
| | | | |
| BLK_ENV-268217 | Method Blank | <0.01 | <0.01 |
| BLK_ENV-268571 | Method Blank | <0.01 | <0.01 |

KEY TO DATA TABLE

UNITS

| | |
|-------------------|--|
| µg | micrograms, relative mass value |
| µg/m ³ | micrograms per cubic meter; estimated soil gas concentration |
| µg/L | micrograms per Liter; calculated water concentration |

DATA QUALIFIERS

| | |
|---|--|
| > | greater than; value exceeds calibration range, estimated value |
| < | less than; compound value is below the LOD and RL |
| J | mass value below LOQ or RL, but above LOD, estimated mass value |
| E | mass value exceeds upper calibration level, estimated mass value |
| Q | one or more quality control parameters failed for the compound |

ABBREVIATIONS

| | |
|--------|--|
| AVG RL | average reporting limit; calculated based on individual field sample RLs |
| LOD | limit of detection |
| LOQ | limit of quantification |
| MDL | method detection limit |
| RL | reporting limit |

| | | | |
|-------------|---|----------|--|
| 1112TetCA | 1,1,1,2-tetrachloroethane | CIBENZ | chlorobenzene |
| 111TCA | 1,1,1-trichloroethane | ct12DCE | cis- & trans-1,2-dichloroethene |
| 1122TetCA | 1,1,2,2-tetrachloroethane | EtBENZ | ethylbenzene |
| 112TCA | 1,1,2-trichloroethane | mpXYL | m-, p-xylene |
| 11DCA | 1,1-dichloroethane | MTBE | methyl t-butyl ether |
| 11DCE | 1,1-dichloroethene | NAPH | naphthalene |
| 124TMB | 1,2,4-trimethylbenzene | OCT | octane |
| 12DCA | 1,2-dichloroethane | oXYL | o-xylene |
| 12DCB | 1,2-dichlorobenzene | PCE | tetrachloroethene |
| 135TMB | 1,3,5-trimethylbenzene | PENTADEC | pentadecane |
| 13DCB | 1,3-dichlorobenzene | PHEN | phenanthrene |
| 14DCB | 1,4-dichlorobenzene | t12DCE | trans-1,2-dichloroethene |
| 2MeNAPH | 2-methyl naphthalene | TCE | trichloroethene |
| BENZ | benzene | TMBs | combined masses of 1,3,5-trimethylbenzene and 1,2,4-trimethylbenzene |
| BTEX | combined masses of benzene, toluene, ethylbenzene, and total xylenes (Gasoline Range Aromatics) | TOL | toluene |
| C11,C13&C15 | combined masses of undecane, tridecane, and pentadecane (C11+C13+C15) (Diesel Range Alkanes) | TPH | total petroleum hydrocarbons |
| c12DCE | cis-1,2-dichloroethene | TRIDEC | tridecane |
| CCl4 | carbon tetrachloride | UNDEC | undecane |
| CHC13 | chloroform | VC | vinyl chloride |

SUMMARY OF SAMPLING RATE CALIBRATION FOR AGI UNIVERSAL SAMPLER IN AQUEOUS PHASES

INTRODUCTION:

The Amplified Geochemical Imaging, LLC (AGI) passive vapor sampler is designed to be used for soil gas, water, sediment pore water, and air sampling. This document describes the process used to calibrate the sampler's compound specific sampling or uptake rates in aqueous phases.

Sampling rates are measured following AGI's "Standard Practice for Determining the Sampling Rate of Passive Diffusion Samplers in Various Environmental Media": SPG-SOP-0493. Rates are used to calculate dissolved phase concentrations of volatile and semi-volatile contaminants in water. The calibration process is summarized in three parts: Part 1: shallow water, Part 2: deep water, and Part 3: sediment.

PURPOSE:

The purpose of this document is to:

1. Summarize the test protocol,
2. Summarize the methodology for analysis of data,
3. Present general results for generating concentration calibration of the AGI Universal Sampler

Principle of Operation of the AGI Sampler

The AGI Universal Sampler is designed with solid adsorbents enclosed inside a tubular microporous PTFE membrane. When placed in water, the pores and hydrophobic nature of the PTFE keep liquid water from entering the membrane until a water head of about 34 feet is reached. The membrane will not keep water vapor from entering but the adsorbents are very hydrophobic and through testing validated to be unaffected by this moisture vapor. In shallow water, <34', volatile and semi-volatile compounds will partition from the dissolved water into the air phase in the PTFE membrane according to Henry's Law. This partitioning is instantaneous and within seconds-minutes, the compound is adsorbed by the adsorbent inside the sealed tube. Because the diffusivity in air is about 10,000 times higher than the diffusivity in water, the sampling rate is controlled by the water contact area of the membrane that allows the Henry's Law effect to occur. This contact area is set by the membrane diameter and length of the sealed tube, which is fixed in AGI's manufacturing process.

Henry's law as well as diffusivity, which are fundamentally incorporated into the sampling rate, are affected by temperature, T , and follow an Arrhenius equation $H_T = H_r \times \exp\left(\frac{-E_a/R}{1/T_r - 1/T}\right)$. Because a 5°C temperature change can make a 15% change in sampling rate, the temperature of the sampled water should be known to get the most precise concentration.

The membrane pore size is also small enough that colloidal particles and microbes cannot pass through the membrane. This keeps the adsorbent from getting contaminated and eliminates any need to add preservative or chilling during storage or transportation.

When the water pressure exceeds the water entry pressure of the membrane, about 34 feet of water, the water becomes in contact with the solid adsorbent. Under this condition the compounds in the water will partition from the water into the solid. The partitioning coefficient, K_{AW} , can be approximated by the octanol-water coefficient, K_{OW} , but has been measured more precisely in the lab for AGI's specific solid adsorbent. The sampling rate is the product of the sampling rate at <34' of water and the K_{AW} .

In sediment, the sampler measures pore-water concentration, which is generally agreed to be the preferred measurement as it is more indicative of bioavailability. In sediment the volumetric

availability of water to the sampler is reduced by the volume fraction solids in the sediment, which typically varies from zero to 35%, but can be as high as 73% in well packed and broad particle size distribution sediments. As a result, sampling rates in sediment are multiplied by the fraction pore water in the sediment to determine concentration.

PART 1: Calibration in shallow water

Part 1 summarizes the work in shallow water generating calibration data, evaluating the physical and chemical factors affecting the sampling rate, and measurement of the actual sampling rates or regression calibration equations needed to determine concentrations.

Sample Generation in water

In this calibration work, solutions of analytes at known concentrations were formulated in clean 4 liter smoked glass jugs by injecting microliter measured amounts of environmental standards using a calibrated syringe into pure or deionized water and stirring for a minimum of 2 hours but generally overnight. Headspace in the jugs was minimized and generally less than 1% by volume during the tests. Jugs were temperature controlled by placing them in a water filled cooler, chilled via a copper tubing loop in the cooler. Temperature was measured with a certified digital temperature gauge and an average value used for each temperature experiment.

AGI samplers were weighted so they won't float and placed in the jugs at time zero. They were removed at various intervals to generate samples along with duplicates that showed mass increasing with exposure time. The sampler exposure time was selected to span minutes to hours and was generally reduced for high concentration tests to maintain uptake with time in roughly the linear dynamic range. Samplers were removed and dried with a paper towel and returned to their original container for analysis. They were analyzed by AGI's 8260C (SPG-WI-318 or SPG-WI-10028) method in duplicate, which is based on EPA SW846 Method 8260C.

Water samples were also taken and measured at an outside accredited lab using EPA SW846 Method 8260B. The concentrations agreed well with the calculated concentrations based on the standard certification, jug volume, and syringe injection. The variability of the outside lab 8260B values were found to be high, so for the sampling rate calculations we used the concentrations based on syringe dosing.

Calibrations were run at five concentrations, nominally at 6, 24, 118, 590, 1420 ug/L and five temperatures nominally at 5, 10, 15, 20, and 25 degrees centigrade. Samples were taken at 4 different exposure times. Samples were run in duplicate. A total of 176 data points were generated using 28 compounds from AGI's standard compounds list. Tridecane and pentadecane were not evaluated due to their very low solubility in water. In addition, another 23 compounds were tested using an 8260 liquid standard at nominal concentrations of 0.5, 1.0, 5.0, 15, 95, and 470 ug/L at a temperature typical of groundwater, 15°C. This is a living calibration and as additional data are generated, they may be qualified and added to this data set to improve the precision of the sampling rate calibration and broaden the compound list.

Key Variable Effects

As expected from theory, at short to moderate exposure times, mass will increase roughly linearly proportional to exposure time, as well as proportional to concentration, and exponentially with temperature following Arrhenius law. Temperature affects the Henry's law as well as diffusivity in water. Sampling rate is generally independent of concentration and time at mass values significantly below saturation. In the following sections we have characterized the sampling rate for each compound as affected by temperature and also developed calibrations using regression which account for the minor impact of time, and mass.

Concentration using Simple Sampling Rate Determination

A simple way to determine concentration is to measure mass on the AGI sampler, divide by exposure time, and divide by sampling rate, SR.

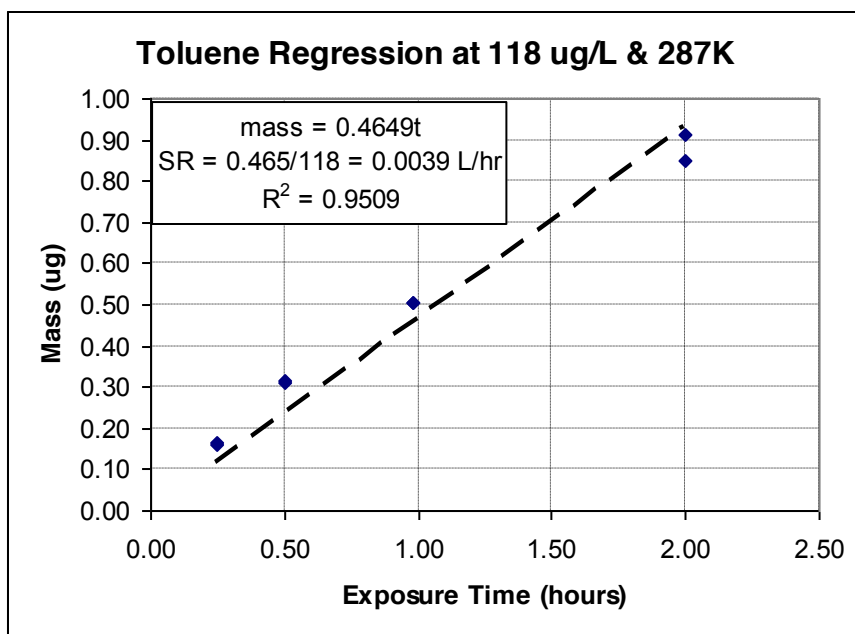
$$\text{Conc [ug/L]} = \text{mass/time/SR} \quad (1)$$

The sampling rate can be determined via measurements of mass versus time at a known concentration and temperature according to the following modification of equation (1).

$$\text{SR} = \text{mass/time/concentration} \quad (2)$$

Sampling rates in L/hr were determined by measuring the trend or regression mass uptake versus time and dividing by the concentration. A measurement like this will use 8 data points (4 times x 2 samples). Such a sampling rate can be measured at any concentration and temperature.

The chart to the right shows a plot of mass versus time for water at 118 ug/L and 287K (actual data from a single run). Slope of 0.465 ug/hr divided by the concentration of 118 ug/L yields a sampling rate, SR, of 0.0039 L/hr.



SR's typically range from about 0.004 to 0.007 L/hr at 15°C. Table A shows SR's measured for our standard compound list at 5 temperatures.

Rigorous Concentration using Regression

A preferred method for determining concentration that will yield improved accuracy over a wide range of concentrations, exposure times, and temperatures is to use all data in a regression analysis, which allows adjustments for the minor non-linear influences of mass and time as well as the effects of temperature. This step is done by regressing equation (1) or a universal version of equation (1):

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRo} * \exp(-\text{Ea}/\text{R}/\text{T})] \quad (3)$$

The subtle non-linear effects of mass and time will be evident in the deviation of coefficients b and d from 1.0. This regression generates four constants b, d, SRo, and $-\text{Ea}/\text{R}$ by regressing $\ln(\text{conc})$ versus $\ln(\text{mass})$, $\ln(\text{time})$, $1/\text{temp}$. These four constants can be used to determine concentration via the equation:

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRo} * \exp(-\text{Ea}/\text{R}(1/\text{T}))] \quad (4)$$

Where conc is in ug/L, mass is in ug, time in hours, T in degrees Kelvin.

Equation (4) can be also expressed at a reference temperature, Tr , such as 15°C by

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRr} * \exp(-\text{Ea}/\text{R}(1/\text{Tr}-1/\text{T}))] \quad (5)$$

This step allows sampling rates, SRr, at any reference temperature, Tr , and for any analyte to easily be compared. The values of SRr at 293.14K can be found in Table A.

When sampling times are between 0 and 4 hours, using the 4 constant equation (5) is preferred. For concentrations from about 5 to 1500 ug/L one hour exposure times generally give the lowest error, typically with average error of 6-20% and with total error range of 12%-32%. For low concentrations where sampling times are greater than 4 hours, it is preferred to use equation (1) to avoid unrealistic effects from the coefficient d or to set d to 1.0. In such a case SR in equation (1) can be substituted with $[\text{SRr} * \exp(-\text{Ea}/\text{R}(1/\text{Tr}-1/\text{T}))]$ to use an SR representative of the well temperature, T.

The chart to the right is a plot of the calculated concentration from the 4 constant regression compared to the dosed concentration. Agreement is excellent for the 176 data points.

However, there does appear to be a slight high bias of 8.6% over the full range of this data, although it is well within acceptable limits of variability.

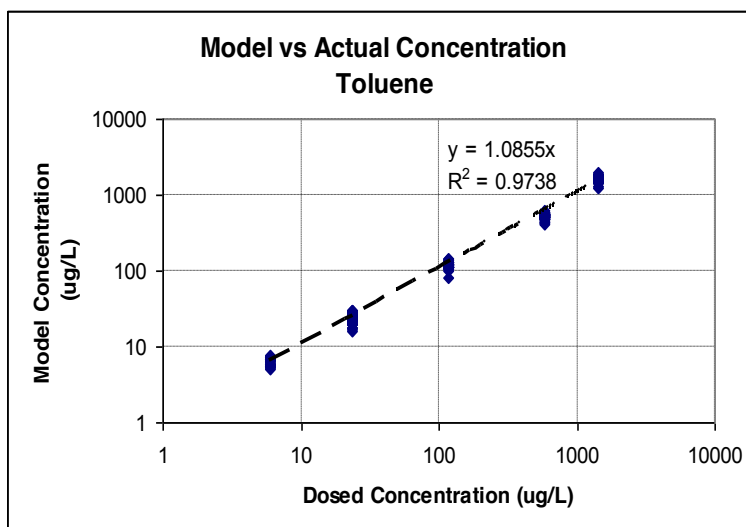


Table B shows the tabulated summary of the 4 constants regression with Rsq values and error estimates for the 4 constants for each analyte. Most regression Rsq values are 0.99 or greater for each analyte. In general, $-E_a/R$ is about 2400 \pm 400, b is about 0.9, d is about -0.75, and SR(15°C) ranges from .004 L/hr to 0.007 L/hr increasing with MW of the compound.

Error Estimates

The error in the water concentration values will depend on both the error in mass from the analytical method as well as the error in the concentration calibration. Table C shows the error in the mass values from the 8260C low sensitivity method.

The standard error of the regression and standard errors of the constants can be found in table B. For each compound we have measured the error between the derived concentration and the actual concentration. The error tends to be lowest at our recommended exposure time of one hour as shown by the example for Toluene to the right.

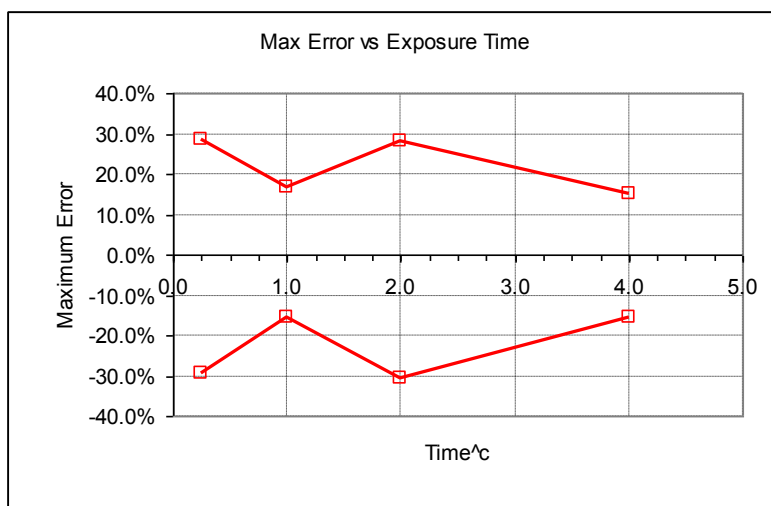
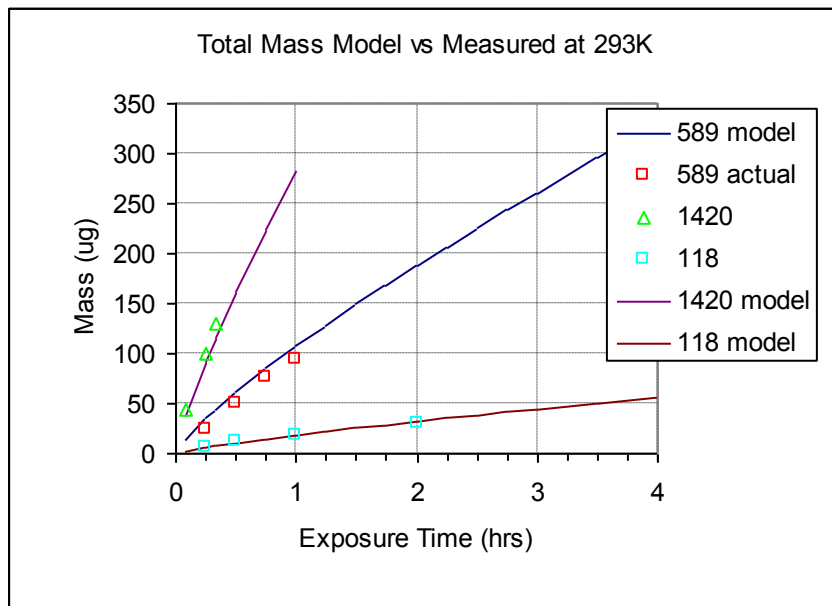


Table D shows the total average error in water concentration by compound as well as the low and high error. The average ranges from about 6% to 20%, which is similar to the analytical method errors. The low and high errors range from 12% to 32% and include contribution from measurement errors in both time and temperature.

Sorbent Saturation

As mass increases on a solid sorbent and approaches saturation, reverse diffusion can occur causing the sampling rate to drop. Eventually the mass level will reach a maximum steady state value at any concentration. A rate of mass uptake with time that deviates significantly from linear, indicates that sorbent saturation could be an issue. When using equation (1), staying in the linear range to avoid the effects of adsorbent saturation is important. We recommend keeping the total mass on the sampler below 50 ug or flagging when this is exceeded.

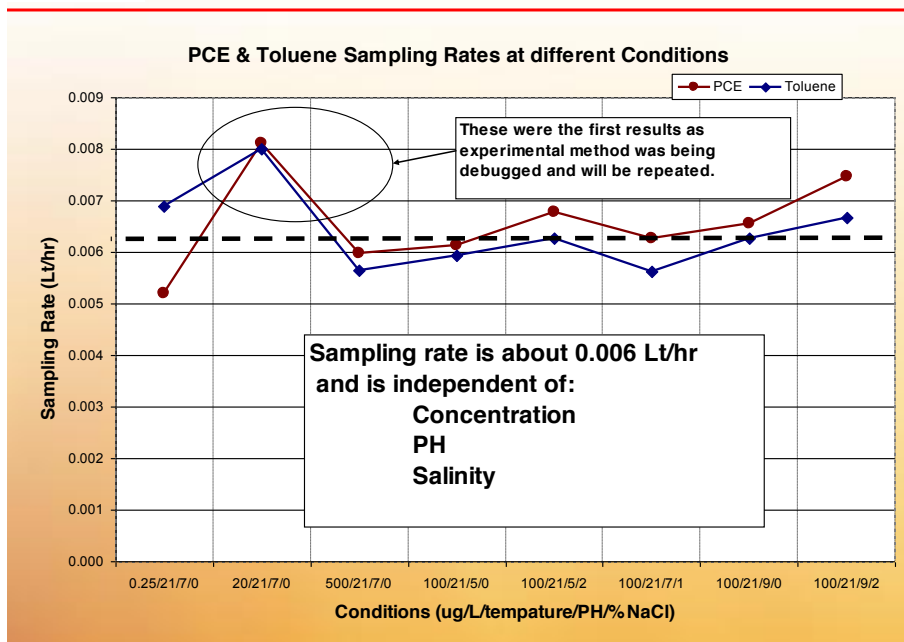
The 4 constant regression accounts for some of the non linearity allowing good accuracy at higher mass levels. From the experimental data we have found this safe range can be extended to 100 ug or higher as shown in the chart below. This chart compares total mass of all compounds (excluding heavy alkanes, which have solubility issues) versus time in comparison to that predicted from the 4-constant concentration equation.



Effect of PH and Salinity

Because neither PH nor salinity is known to have a significant impact on Henry's law or diffusivity in water, we did not expect them to have a significant impact on sampling rate. To confirm this, experiments were run varying PH from 5 to 9 and NaCl content from 0 to 2%. The chart below shows no significant impact for combinations of PH and NaCl content over this range on the sampling rate of toluene in water at 21°C.

Checked for Effects of PH & Salinity

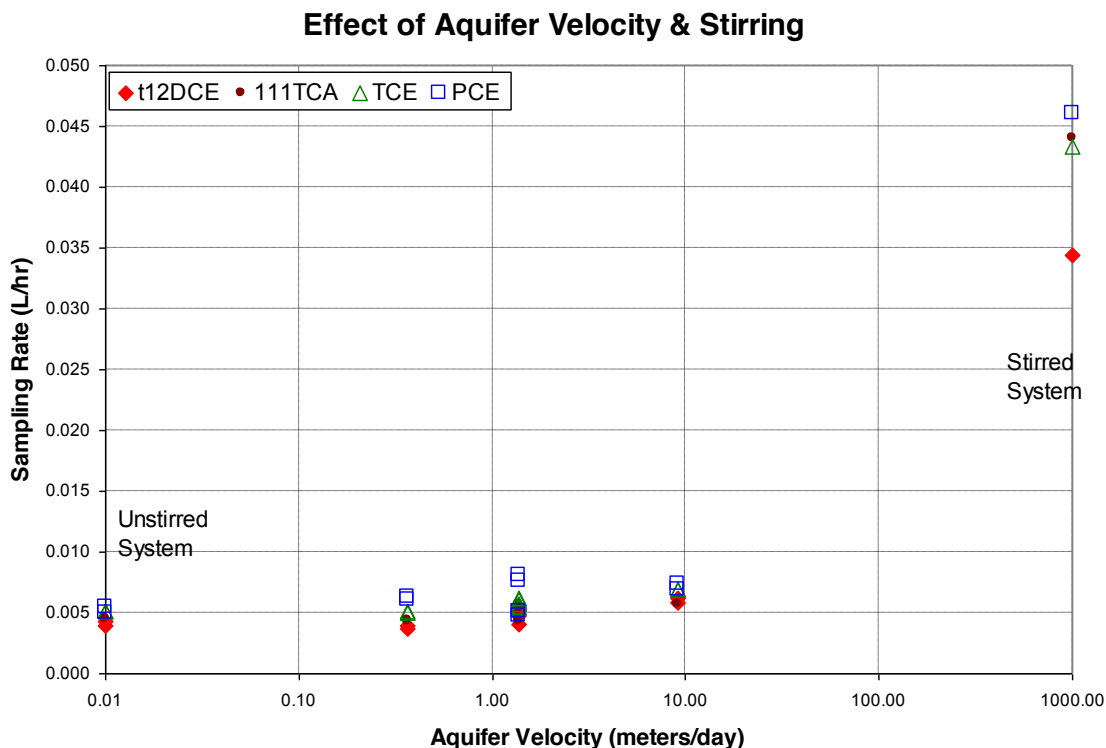


Impact of Aquifer Velocity

The velocity in most aquifers is quite slow, typically a meter/day or less. Occasionally water flow could be much higher such as encountered in karst aquifers, streams or rivers. Mass transfer coefficients are higher in high flow conditions, which will lead to higher sampling rates. We validated that a highly stirred system had sampling rates about 10 times higher than those that were non-stirred. We decided to evaluate the effect of aquifer velocity.

A test apparatus was built comprising a 3" PVC pipe tee filled with clean sand in each of the horizontal straight legs and screened to leave the center open. A test solution was run through this system using a variable flow pump and AGI samplers were placed into the simulated well through the vertical leg of the tee. Tests were run to examine the effect of velocity by varying the pumping rate and hence water velocity.

The chart below shows no significant effect of aquifer velocity up to a speed of about 10 meters/day. At velocities significantly above this, similar to a stirred system, sampling rates are about 10 times higher.

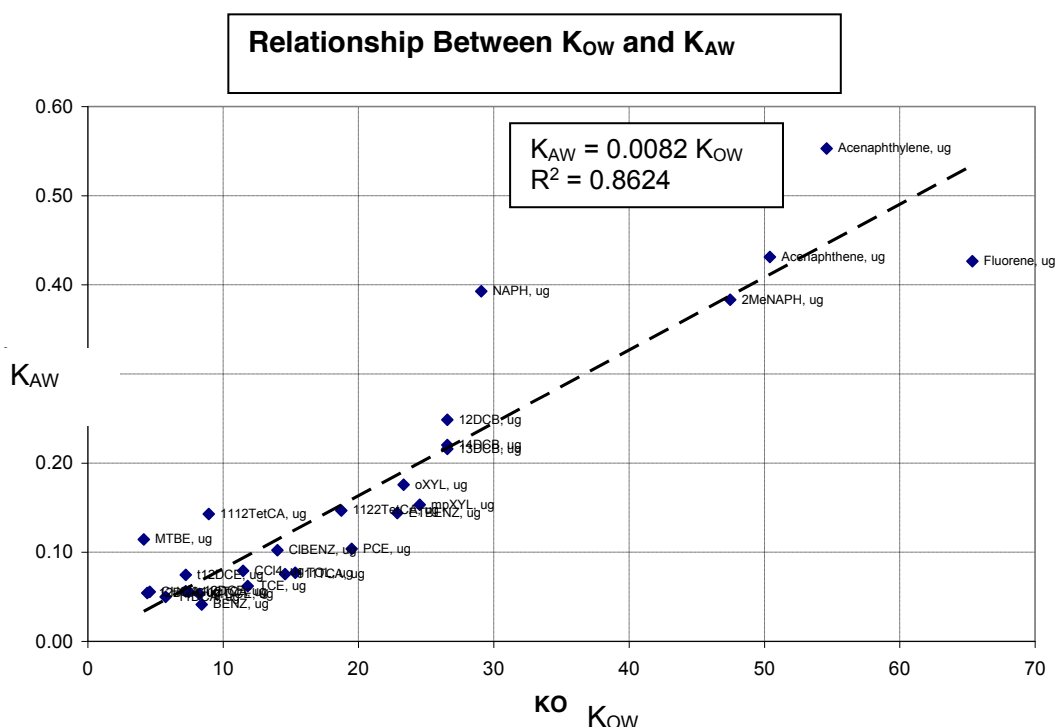


Part 2: Calibration in Deep (>34') water

Part 2 describes the effect of deep water on the AGI sampler and summarizes the effects on sampling rate and concentration measurement.

When the water pressure exceeds the water entry pressure of the membrane, about 34 feet of water, the water becomes in direct contact with the solid adsorbent. Under this condition the compounds in the water will partition from the water into the solid. The partitioning coefficient is closely related to the octanol-water coefficient, K_{OW} , but has been measured more precisely in the lab for AGI's specific solid adsorbent, K_{AW} . The sampling rate for deep water is the product of the sampling rate at <34' of water and the K_{AW} .

Measurement of the K_{AW} was done in a one liter stainless steel vessel pressurized with nitrogen to simulate water heads above 34' of water. Pressures of up to 465 psig or 200' of water head were used. The sampling rate change was the same at all pressures above 34' of water. The K_{AW} was determined as the ratio between the mass or sampling rate above 34' of head to the rate at <34' of head and is shown in the chart below.



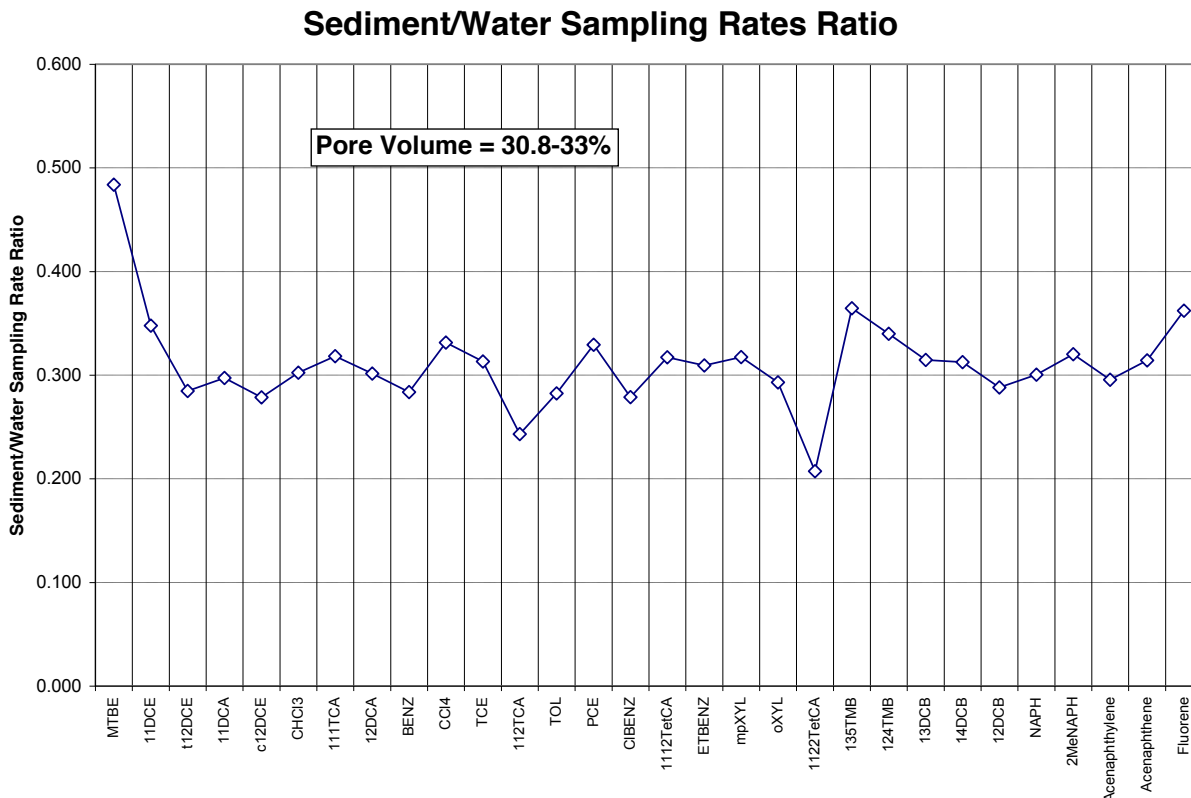
Part 3: Calibration in Sediment

Part 3 describes the effect of sediment solids or sediment pore volume on the sampling rate and concentration measurement.

In sediment, the sampler measures pore water concentration, which is generally agreed to be the preferred measurement as it is more indicative of bioavailability. In sediment the volumetric availability of water to the sampler is reduced by the volume fraction solids in the sediment. As a result sampling rates in sediment are multiplied by the fraction pore water to determine

concentration. Pore water fraction can range from 1.0 for water without sediment to as low as 0.25. Typically most sediments have pore fractions of 0.9 to 0.65.

A sampling rate study was done with water and with water added into a well-packed sorted sand. Pore water fraction in this test was measured between 30.8% and 33% by volume. Below is a plot of the ratio of sampling rates measured in the sediment to open water. The average ratio is equal to the pore water fraction confirming that sampling rate in sediment is on average equal to the product of pore water fraction times the sampling rate in water.



Summary

The AGI Sampler can be used to determine the concentration of volatile and semi-volatile compounds in a water phase. This requires knowing the exposure time and water temperature. It also requires knowing if the sample is above or below 34' of water head and if the water has a velocity above 10 meters/day. Regressions of large amounts of data were used to generate a four constant equation to generate concentration values in water. Potential error in the concentration values is excellent typically less than 25%.

TABLE A
WATER SAMPLING RATES STANDARD LIST

| | SRr 293.14 | SR @ 277.54 | SR @ 282.44 | SR @ 287.84 | SR @ 293.24 | SR @ 298.94 |
|-------------------|----------------------|-----------------------|-----------------------|-----------------------|-----------------------|-----------------------|
| MTBE | 0.0025 | 0.0014 | 0.0016 | 0.0018 | 0.0022 | 0.0029 |
| t12DCE | 0.0043 | 0.0028 | 0.0028 | 0.0027 | 0.0037 | 0.0048 |
| 11DCA | 0.0047 | 0.0031 | 0.0033 | 0.0033 | 0.0039 | 0.0052 |
| c12DCE | 0.0046 | 0.0030 | 0.0031 | 0.0031 | 0.0038 | 0.0051 |
| CHCl3 | 0.0046 | 0.0030 | 0.0031 | 0.0031 | 0.0038 | 0.0051 |
| 111TCA | 0.0066 | 0.0043 | 0.0047 | 0.0047 | 0.0056 | 0.0076 |
| 12DCA | 0.0045 | 0.0029 | 0.0029 | 0.0030 | 0.0036 | 0.0050 |
| BENZ | 0.0050 | 0.0031 | 0.0034 | 0.0035 | 0.0042 | 0.0056 |
| CCl4 | 0.0068 | 0.0044 | 0.0048 | 0.0047 | 0.0058 | 0.0080 |
| TCE | 0.0052 | 0.0030 | 0.0034 | 0.0034 | 0.0043 | 0.0058 |
| 112TCA | 0.0043 | 0.0027 | 0.0027 | 0.0028 | 0.0034 | 0.0048 |
| TOL | 0.0056 | 0.0034 | 0.0039 | 0.0039 | 0.0047 | 0.0062 |
| OCT | 0.0064 | 0.0046 | 0.0050 | 0.0040 | 0.0058 | 0.0089 |
| PCE | 0.0061 | 0.0036 | 0.0043 | 0.0043 | 0.0051 | 0.0069 |
| CIBENZ | 0.0054 | 0.0033 | 0.0039 | 0.0040 | 0.0045 | 0.0059 |
| 1112TetCA | 0.0061 | 0.0037 | 0.0042 | 0.0044 | 0.0050 | 0.0065 |
| EtBENZ | 0.0060 | 0.0037 | 0.0045 | 0.0044 | 0.0052 | 0.0069 |
| mpXYL | 0.0064 | 0.0039 | 0.0048 | 0.0046 | 0.0055 | 0.0072 |
| oXYL | 0.0066 | 0.0041 | 0.0050 | 0.0048 | 0.0057 | 0.0074 |
| 1122TetCA | 0.0044 | 0.0027 | 0.0029 | 0.0031 | 0.0036 | 0.0046 |
| 135TMB | 0.0079 | 0.0046 | 0.0059 | 0.0056 | 0.0071 | 0.0093 |
| 124TMB | 0.0078 | 0.0046 | 0.0060 | 0.0055 | 0.0071 | 0.0092 |
| 13DCB | 0.0072 | 0.0041 | 0.0055 | 0.0053 | 0.0063 | 0.0080 |
| 14DCB | 0.0071 | 0.0040 | 0.0054 | 0.0052 | 0.0062 | 0.0079 |
| 12DCB | 0.0070 | 0.0040 | 0.0053 | 0.0051 | 0.0060 | 0.0076 |
| UNDEC | | 0.0026 | 0.0024 | 0.0020 | 0.0031 | 0.0029 |
| NAPH | | 0.0041 | 0.0056 | 0.0054 | 0.0064 | 0.0081 |
| TRIDEC | | | | | | |
| 2MeNAPH | | 0.0043 | 0.0066 | 0.0066 | 0.0080 | 0.0108 |
| PENTADEC | | | | | | |
| Total mass | 0.1177 | 0.0822 | 0.1339 | 0.1334 | 0.1773 | 0.1981 |

Notes:

Values in L/hr

Total mass does not include UNDEC, TRIDEC, PENTADEC (28 compounds)

TABLE B
4 CONSTANT REGRESSION OUTPUT

| | Adjusted Rsq | Standard Error | ln(SR0) | b | -Ea/R | d | Std Error ln(SR0) | Std Error b | Std Error - Ea/R | Std Error d |
|------------|-------------------------|---------------------------|----------------|----------|--------------|----------|----------------------------------|----------------------------|-------------------------------------|----------------------------|
| MTBE | 0.997 | 0.0960 | -3.217 | 0.981 | 2704 | -0.709 | 0.2881 | 0.0062 | 83 | 0.0082 |
| t12DCE | 0.992 | 0.1659 | -1.877 | 0.905 | 2147 | -0.760 | 0.4971 | 0.0100 | 144 | 0.0138 |
| 11DCA | 0.995 | 0.1272 | -1.346 | 0.916 | 1965 | -0.746 | 0.3809 | 0.0077 | 110 | 0.0106 |
| c12DCE | 0.995 | 0.1299 | -1.905 | 0.911 | 2137 | -0.751 | 0.3892 | 0.0078 | 112 | 0.0109 |
| CHCl3 | 0.996 | 0.1260 | -1.841 | 0.912 | 2118 | -0.748 | 0.3776 | 0.0076 | 109 | 0.0105 |
| 111TCA | 0.995 | 0.1279 | -2.684 | 0.902 | 2259 | -0.761 | 0.3836 | 0.0076 | 111 | 0.0106 |
| 12DCA | 0.995 | 0.1263 | -2.161 | 0.908 | 2218 | -0.746 | 0.3786 | 0.0076 | 109 | 0.0106 |
| BENZ | 0.995 | 0.1323 | -2.207 | 0.920 | 2198 | -0.754 | 0.3965 | 0.0080 | 114 | 0.0110 |
| CCl4 | 0.994 | 0.1405 | -3.121 | 0.889 | 2379 | -0.776 | 0.4220 | 0.0083 | 122 | 0.0116 |
| TCE | 0.992 | 0.1655 | -3.338 | 0.900 | 2522 | -0.772 | 0.4969 | 0.0099 | 144 | 0.0137 |
| 112TCA | 0.995 | 0.1264 | -2.412 | 0.896 | 2302 | -0.724 | 0.3790 | 0.0075 | 109 | 0.0107 |
| TOL | 0.994 | 0.1426 | -2.873 | 0.916 | 2364 | -0.756 | 0.4281 | 0.0087 | 124 | 0.0119 |
| OCT | 0.938 | 0.4698 | -5.984 | 0.822 | 3235 | -0.827 | 1.4231 | 0.0277 | 412 | 0.0388 |
| PCE | 0.991 | 0.1773 | -3.780 | 0.877 | 2601 | -0.775 | 0.5329 | 0.0103 | 154 | 0.0147 |
| CIBENZ | 0.994 | 0.1457 | -2.601 | 0.911 | 2292 | -0.747 | 0.4370 | 0.0088 | 126 | 0.0122 |
| 1112TetCA | 0.996 | 0.1235 | -2.676 | 0.898 | 2281 | -0.725 | 0.3705 | 0.0073 | 107 | 0.0104 |
| EtBENZ | 0.993 | 0.1597 | -2.930 | 0.918 | 2357 | -0.752 | 0.4794 | 0.0097 | 138 | 0.0134 |
| mpXYL | 0.992 | 0.1678 | -3.036 | 0.909 | 2372 | -0.749 | 0.5037 | 0.0101 | 145 | 0.0140 |
| oXYL | 0.993 | 0.1555 | -2.862 | 0.911 | 2312 | -0.740 | 0.4667 | 0.0094 | 135 | 0.0131 |
| 1122TetCA | 0.996 | 0.1118 | -1.971 | 0.913 | 2167 | -0.691 | 0.3351 | 0.0067 | 97 | 0.0096 |
| 135TMB | 0.988 | 0.2024 | -4.435 | 0.897 | 2720 | -0.738 | 0.6093 | 0.0121 | 176 | 0.0170 |
| 124TMB | 0.989 | 0.1997 | -4.126 | 0.890 | 2631 | -0.731 | 0.6009 | 0.0118 | 173 | 0.0169 |
| 13DCB | 0.991 | 0.1832 | -3.422 | 0.888 | 2449 | -0.730 | 0.5503 | 0.0108 | 159 | 0.0155 |
| 14DCB | 0.991 | 0.1802 | -3.263 | 0.892 | 2408 | -0.724 | 0.5413 | 0.0107 | 156 | 0.0153 |
| 12DCB | 0.992 | 0.1697 | -2.970 | 0.894 | 2327 | -0.716 | 0.5092 | 0.0101 | 147 | 0.0144 |
| UNDEC | 0.694 | 0.374 | -1.406 | 0.426 | 1708 | -0.806 | 1.792 | 0.028 | 517 | 0.053 |
| NAPH | 0.992 | 0.166 | -3.374 | 0.915 | 2430 | -0.671 | 0.497 | 0.010 | 144 | 0.014 |
| TRIDEC | | | | | | | | | | |
| 2MeNAPH | 0.984 | 0.238 | -5.498 | 0.869 | 2990 | -0.689 | 0.72 | 0.014 | 208 | 0.021 |
| PENTADEC | | | | | | | | | | |
| Total mass | 0.993 | 0.1543 | -6.111 | 0.907 | 2419 | -0.732 | 0.4666 | 0.0093 | 134 | 0.0130 |

TABLE C
8260C MASS UNCERTAINTY

**AGI 8260C Method for Mass using SPG-0008
Samplers**

| | 99% Uncertainty Range +/- | 95% Uncertainty Range +/- |
|-----------|---------------------------------|---------------------------------|
| MTBE | 20% | 14% |
| t12DCE | 22% | 15% |
| 11DCA | 18% | 12% |
| c12DCE | 18% | 12% |
| CHCl3 | 16% | 11% |
| 111TCA | 18% | 12% |
| 12DCA | 20% | 13% |
| BENZ | 16% | 10% |
| CCl4 | 19% | 12% |
| TCE | 15% | 10% |
| 112TCA | 18% | 12% |
| TOL | 15% | 10% |
| OCT | 20% | 13% |
| PCE | 16% | 11% |
| CIBENZ | 18% | 12% |
| 1112TetCA | 19% | 13% |
| EtBENZ | 18% | 12% |
| mpXYL | 18% | 12% |
| oXYL | 18% | 12% |
| 1122TetCA | 23% | 15% |
| 135TMB | 21% | 14% |
| 124TMB | 20% | 14% |
| 13DCB | 19% | 13% |
| 14DCB | 19% | 13% |
| 12DCB | 20% | 14% |
| NAPH | 21% | 14% |
| 2MeNAPH | 25% | 17% |

TABLE D
4 CONSTANT WATER CONCENTRATION UNCERTAINTY
ERROR IN CONCENTRATION REPORTING (1)

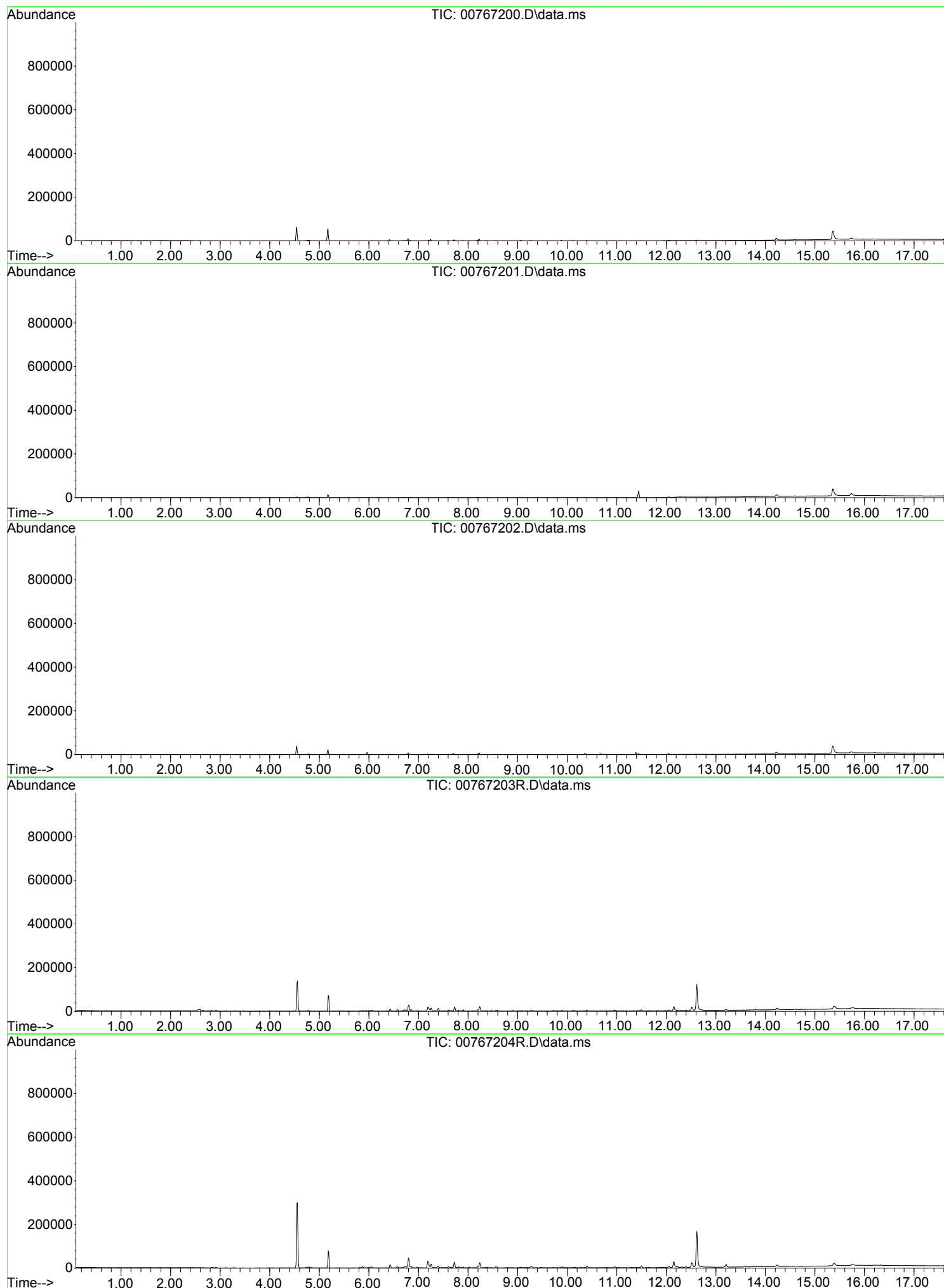
| | Average Error | Minimum Error | Maximum Error |
|-----------|--------------------------|--------------------------|--------------------------|
| MTBE | 6% | -12% | 12% |
| t12DCE | 11% | -26% | 21% |
| 11DCA | 8% | -19% | 13% |
| c12DCE | 9% | -19% | 15% |
| CHCl3 | 9% | -20% | 14% |
| 111TCA | 9% | -19% | 23% |
| 12DCA | 10% | -19% | 17% |
| BENZ | 8% | -18% | 13% |
| CCl4 | 10% | -23% | 22% |
| TCE | 10% | -21% | 14% |
| 112TCA | 11% | -21% | 21% |
| TOL | 7% | -17% | 14% |
| OCT | 20% | -41% | 42% |
| PCE | 10% | -24% | 15% |
| CIBENZ | 7% | -16% | 14% |
| 1112TetCA | 8% | -17% | 18% |
| EtBENZ | 6% | -19% | 14% |
| mpXYL | 7% | -22% | 13% |
| oXYL | 7% | -19% | 13% |
| 1122TetCA | 8% | -16% | 17% |
| 135TMB | 9% | -23% | 17% |
| 124TMB | 10% | -28% | 19% |
| 13DCB | 10% | -22% | 17% |
| 14DCB | 10% | -22% | 17% |
| 12DCB | 9% | -23% | 17% |
| NAPH | 10% | -24% | 21% |
| 2MeNAPH | 13% | -32% | 30% |

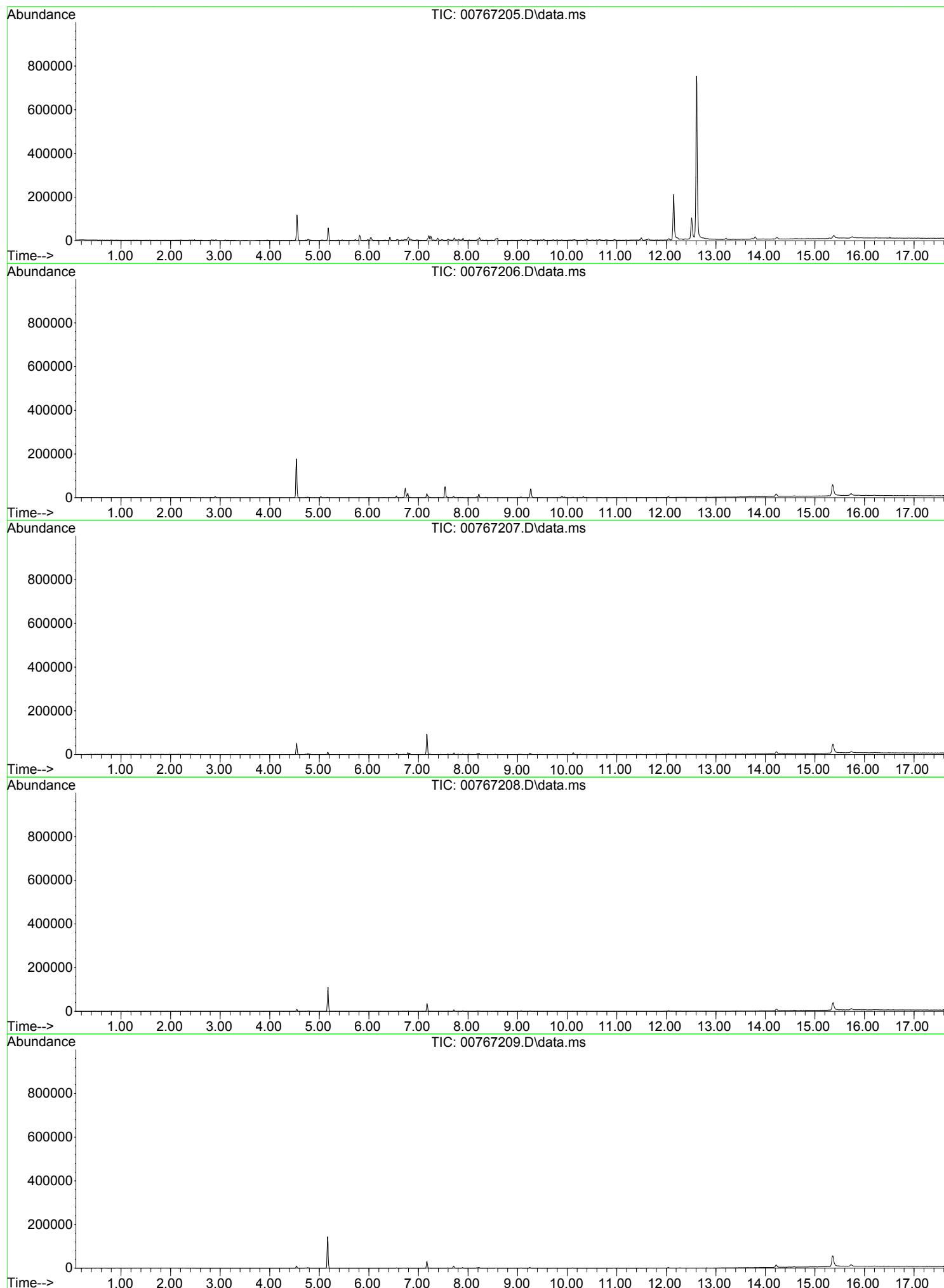
(1) For 1 hour exposure, includes error related to mass value from AGI analytical method 8260C

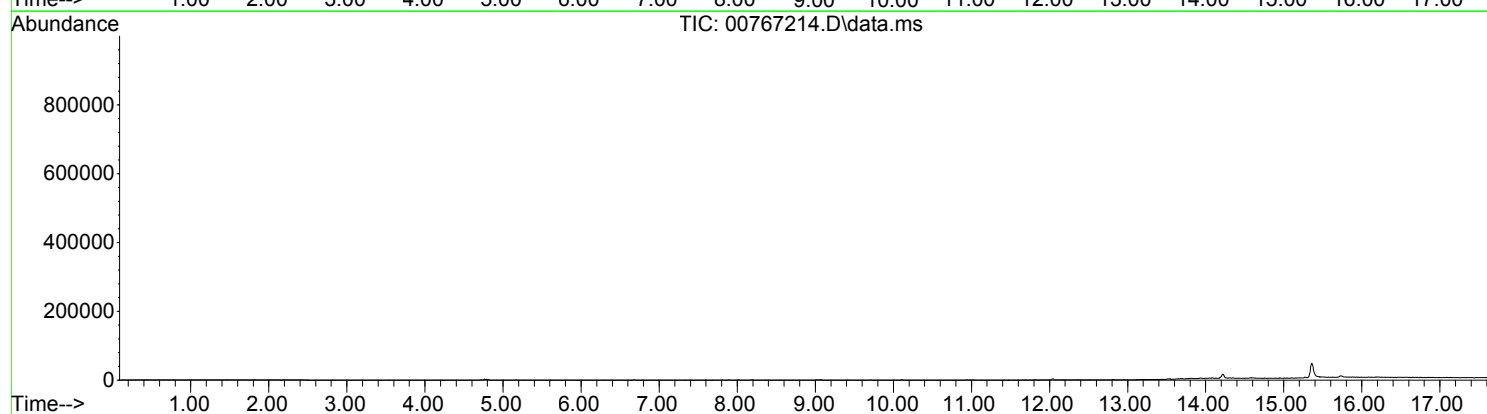
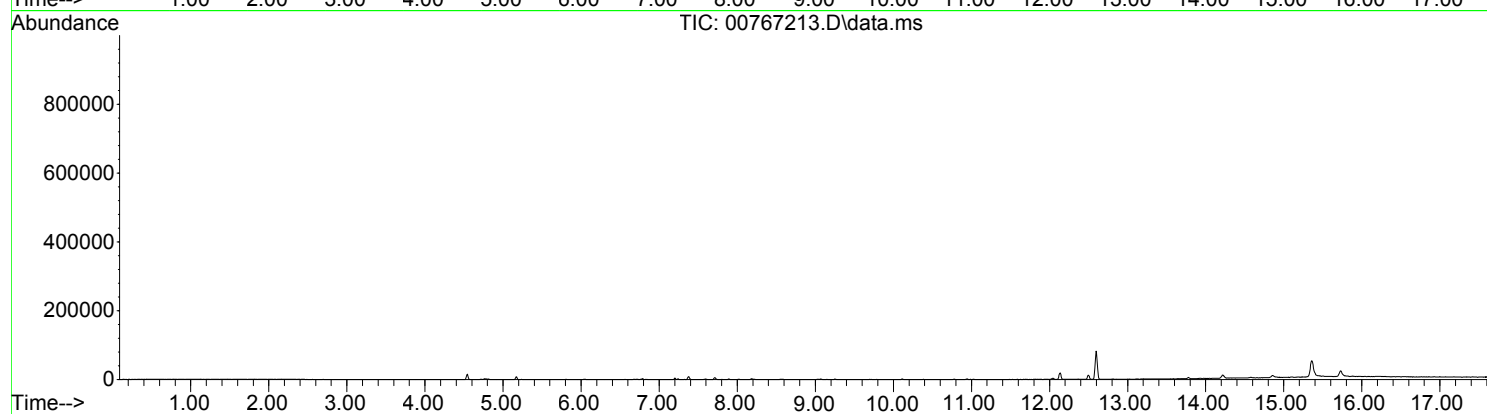
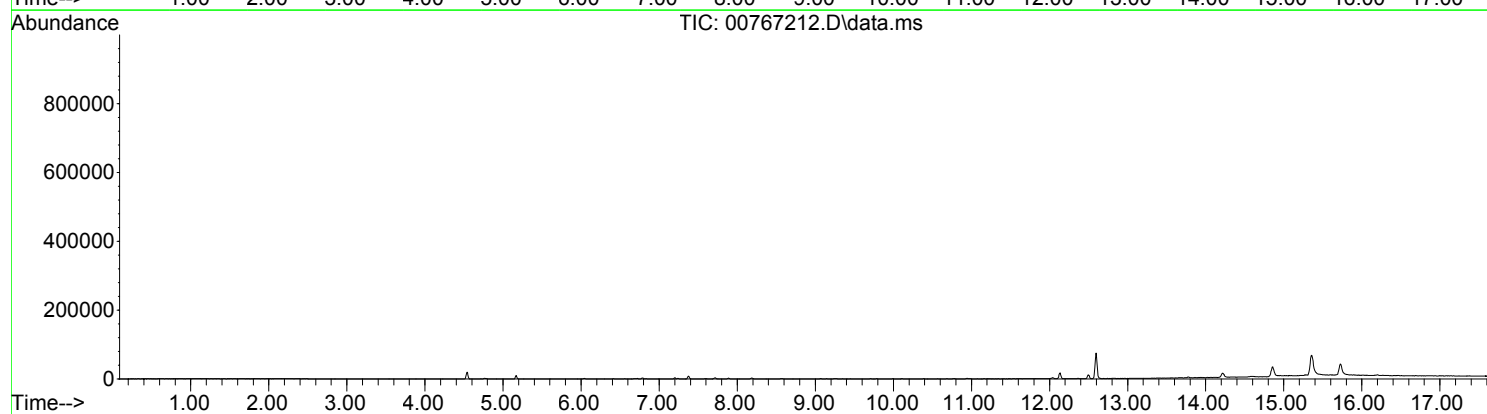
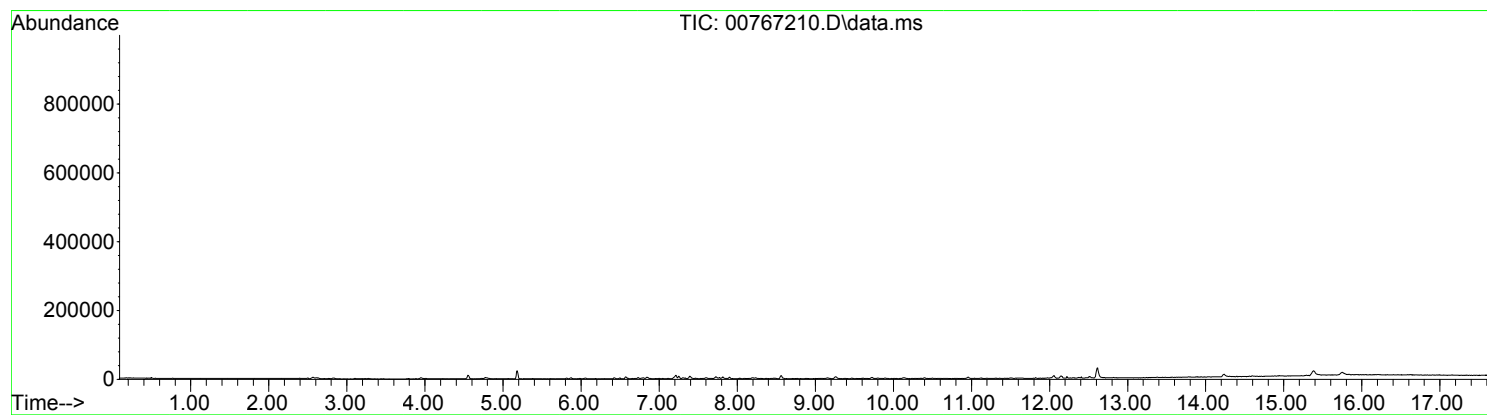
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www.agisurveys.net



AMPLIFIED
GEOCHEMICAL
IMAGING, LLC

Laboratory Report

Site: Comal & San Marcos Rivers October Sampling

Prepared for:

SWCA Environmental Consultants
10245 Little York Road
Suite 600
San Antonio, TX
UNITED STATES

Prepared on:
October 30, 2015

Project Summary and Objective

Amplified Geochemical Imaging, LLC. (AGI) provided the AGI Environmental Survey used at:

Comal & San Marcos Rivers

October Sampling

The service provided by AGI included delivery of the required quantity of AGI Universal Samplers, analysis by the method described below for the requested organic compounds, reporting of the data, and contour mapping (as needed).

This report includes results for only the samples noted under the Laboratory Sample Report section. If contour maps are part of the project deliverable, the maps will be prepared and issued under a separate report cover, upon receipt of a usable sitemap (electronic) and compound choices for contouring.

Written/submitted by:

Jim E Whetzel

Project Manager

Reviewed/approved by:

Don D'Apolito

Project Manager

Analytical data approved by:

Ian McMullen

Chemist

Quality Assurance Statement

The AGI Laboratory, at Amplified Geochemical Imaging's facility in Newark, DE USA, operates under the guidelines of its ISO Standard 17025 DoD ELAP accreditation, and its Quality Assurance Manual, Operating Procedures, and Methods (SPG-SOP-0462).

For this project, the analytical method, results, and observations reported do [] do not [✓] fall within the scope of AGI's ISO 17025 accreditation.

Screening/Concentration Method

The AGI Universal Samplers are analyzed at AGI's fixed laboratory using thermal desorption-gas chromatography/mass spectrometry (TD-GC/MS) instrumentation following modified U.S. EPA Method 8260 (SPG-WI-0292) which includes the following:

- **BFB Tuning Frequency:** A BFB tune is analyzed at the start of each analytical run and after every 30 samples.
- **Initial Calibration:** A minimum of a five point calibration curve is analyzed prior to the analysis of samples.
- **Initial Calibration Verification (ICV):** Following the calibration a second-source reference standard is analyzed to verify the accuracy of the calibration. Acceptance criteria for the ICV is +/- 30%.
- **Linearity of Target Compounds:** If the RSD of any target analyte is less than or equal to 25% then average response factor can be used for quantitation. If the RSD exceeds 25% for a target compound a regression equation can be used for quantitation.
- **Continuing Calibration Verification:** After every 10 samples, and at the end of each analytical batch, a mid-level second-source Reference Standard is analyzed. The acceptance criteria for all target analytes in the reference standards are +/- 50% of the true value.
- **Method Blank:** Analyzed prior to the analysis of field samples and every 30 samples.

Note: Analyte levels reported for the field-deployed AGI Universal Samplers that exceed trip and method blank levels, and/or the reporting limit, are more likely to have originated from on-site sources.

| | |
|----------------------------|---------------|
| Media Sampled: | WATER |
| Chemist - sample analysis: | Jasmine Smith |
| Chemist - data processor: | Jasmine Smith |
| Chemist - data review: | Ian McMullen |

Methodology: See

Please note that data file names ending with R are rerun samples using the second pair of sorbers, in which the original results were not reported. Data file names ending in D are duplicate analysis results for the second set of sorbers from the same sampler, and are reported.

Additional Report Information

- Comments
- Laboratory Sample Report
- Chain of Custody
- Installation and Retrieval Log
- e and Key
- Concentration Calculation Method Summary
- Total Ion Chromatograms

Project Specific Comments

o e

Survey period ¹

er ere ed o O o er d re re ed o O o er
or e o re er od o d

Tamper seal intact:

Yes

Date received:

10/21/15 1 : 5 am

By: Darlene Yellowdy

COC returned:

Yes

Comments:

o e

1 - Installation start to end of retrieval, as reported. See installation and retrieval log for individual deployment and retrieval dates and times (i.e., sampler exposure time).

General Comments

Analytical QA/QC

Laboratory instrumentation consists of gas chromatographs equipped with mass selective detectors, coupled with automated thermal desorption units. Sample preparation involves cutting the tip off the bottom of the AGI Universal Sampler, and transferring one or more "sorbents" to a thermal desorption tube for analysis. The insertion/retrieval cord prevents soil, water and other interferences from coming in contact with the adsorbent. No further sample preparation is required. Any replicate sorbents not consumed in the initial analysis will be discarded fifteen (15) days from the date of the laboratory report.

Data are archived and stored in a secure manner as per AGI's Quality Assurance program (SPG-SOP-0462).

Total petroleum hydrocarbons (TPH), gasoline-range petroleum hydrocarbons (GRPH), and/or diesel range petroleum hydrocarbons (DRPH), when reported, are calculated using the area under the peaks observed in m/z 55 and 57 selected ion chromatograms. Quantitation of the mass values was performed using the response factor for a specific alkane (present in the calibration standards). TPH values include the entire chromatogram and provide estimates for aliphatic hydrocarbon ranges of C4 to C20. GRPH and DRPH include only the relevant regions of the chromatograms and provide estimates for C4 to C10 and C10 to C20 aliphatic hydrocarbons, respectively.

Trip blanks were provided to document potential exposures that were not part of the signal of interest (e.g., impact during sampler shipment, installation and/or retrieval, and storage). The trip blanks are identically manufactured and packaged AGI Universal Samplers to those samplers deployed in the field. The trip blanks remain unopened during all phases of the project. Levels reported on the trip blanks may indicate potential impact to the samplers other than the contaminant source of interest.

Unresolved peak envelopes (UPEs) are represented as a series of compound peaks clustered together around a central gas chromatograph elution time in the total ion chromatogram. UPEs may be indicative of complex fluid mixtures. UPEs observed early in the chromatograms are considered to indicate presence of more volatile fluids, while UPEs observed later in the chromatogram may indicate the presence of less volatile fluids. Multiple UPEs may indicate the presence of multiple complex fluids.

Total ion chromatograms (TICs) are included in the Attachments. The eight-digit serial number of each sampler is incorporated in the TIC identification (e.g., 12345678.D represents AGI Universal Sampler 12345678).

General Comments

Soil Gas Sampling

For soil gas sampling, the AGI Environmental Survey reports mass levels migrating through the open pore spaces of the soil and diffusing through the sampler membrane for sorption by the engineered, hydrophobic adsorbents, housed within the membrane tube. During the migration of the soil gas away from the source to the AGI Universal Sampler, the vapors are subject to a variety of attenuation factors. The soil gas masses reported on the samplers compare favorably with the concentrations reported in the soil or groundwater (e.g., where soil gas levels are reported at greater levels to other sampled locations on the site, the matrix data should reveal the same pattern, and vice versa). However, due to a variety of factors, a perfect comparison between matrix data and soil gas levels can rarely be achieved.

Soil gas concentrations ($\mu\text{g}/\text{m}^3$) are calculated following the method described in the Additional Report Information section.

Soil gas signals reported by this method cannot be correlated specifically to soil adsorbed, groundwater, and/or free-phase contamination. The soil gas signal reported from each AGI Universal Sampler can evolve from all of these sources. Differentiation between soil and groundwater contamination can only be achieved with prior knowledge of the site history (i.e., the site is known to have groundwater contamination only).

Air Sampling

For indoor, outdoor, and crawlspace air sampling, the AGI Environmental Survey reports mass levels present in the air and diffusing through the sampler membrane for sorption by the engineered adsorbents housed within the membrane tube.

Air concentrations ($\mu\text{g}/\text{m}^3$) are calculated following the method described in the Additional Report Information section.

Groundwater and Sediment Porewater Sampling

For groundwater and sediment porewater sampling, the AGI Environmental Survey reports the mass levels of compounds present in the water which, when coming in contact with the sampler membrane, partitions out of solution, and diffuses through the sampler membrane for sorption by the engineered adsorbents.

Water concentrations ($\mu\text{g}/\text{L}$) are calculated using the quantified mass, exposure period and the compound specific uptake rate. The rates were measured under controlled experimental conditions. The uptake rates are corrected for water pressure (depth of the AGI Universal Sampler below the water table), water temperature and the aquifer flow rate. For sediment porewater, the uptake rate is corrected for the reduced volume of water in the sediment, by multiplying the uptake rate by the pore water fraction.

LABORATORY SAMPLE REPORT

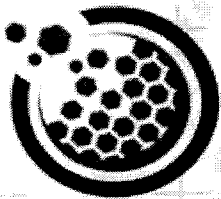
Project: ENV 01509

Site Name: Comal & San Marcos Rivers

Module Type: SPG0008

| Module ID | Sample Type | Field ID | |
|----------------------------|--------------------------|---------------------|-------------------|
| 00770685 | FIELD_SAMPLE | HCS410 | |
| 00770686 | FIELD_SAMPLE | HCS420 | |
| 00770687 | FIELD_SAMPLE | HCS430 | |
| 00770688 | FIELD_SAMPLE | HCS440 | |
| 00770689 | FIELD_SAMPLE | FDHCS440 | |
| 00770690 | FIELD_SAMPLE | HCS460 | |
| 00770691 | FIELD_SAMPLE | HSM410 | |
| 00770692 | FIELD_SAMPLE | HSM420 | |
| 00770693 | FIELD_SAMPLE | HSM430 | |
| 00770694 | FIELD_SAMPLE | FDHSM430 | |
| 00770695 | FIELD_SAMPLE | HSM440 | |
| 00770696 | FIELD_SAMPLE | HSM450 | |
| 00770697 | FIELD_SAMPLE | HSM460 | |
| 00770698 | FIELD_SAMPLE | HSM470 | |
| 00770699 | TRIP_BLANK | TB13 | |
| Total # "FIELD SAMPLES" | Total # "TRIP BLANKS" | Total # "UNUSED" | Total # "LOST" |
| 14 | 1 | 0 | 0 |

Duplicate samples: 0



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Newark, DE 19702-3335 USA
ph: +1-302-266-2428
www.agisurveys.net

AGI Universal Passive Sampler Chain of Groundwater Sampling

Production Order #: **01509**

Customer Name: SWCA Environmental Consultants
Address: 6200 UTSA Boulevard
Suite 102

San Antonio, TX 78249
USA

Site Name: Comal & San Marcos Rivers oct
Site Address:

Project Manager:

Serial # of Samplers Shipped
00770685 - 00770699

| | | | |
|--------------------------------|-------|------------------|---|
| # of Samplers for Installation | 14.00 | # of Trip Blanks | 1 |
| Total Samplers Shipped | 15.00 | Pieces | |
| Total Samplers Received | 15 | Pieces | |
| Total Samplers Installed | 15 | Pieces | |

Serial # of Trip Blanks (Client Decides)

| | | |
|----------|--|--|
| 00770699 | | |
|----------|--|--|

| | |
|--|--|
| Prepared By: <u>[Signature]</u> | Is Concurrent water sampling planned this monitoring period? YES <input type="radio"/> NO <input checked="" type="radio"/> |
| Verified By: <u>Darkene Galloway</u> | Scheduled Sampling Date: _____ |
| Installation Performed By: Name: <u>Jennifer Moreland Brittany Rios</u> Company: <u>SWCA</u> | Retrieval Performed By: Name: <u>Jennifer Moreland Guy Rubio</u> Company: <u>SWCA</u> |
| Installation Start Date / Time: <u>10/6/15 1156</u> | Retrieval Start Date / Time: <u>10/20/15 1303</u> |
| Installation Complete Date / Time: <u>10/6/15 1512</u> | Retrieval Complete Date / Time: <u>10/20/15 1303</u> |
| Total Samplers Retrieved: <u>14</u> | |
| Total Samplers Lost In Field: <u>0</u> | |
| Total Unused Samplers Returned: <u>0</u> | |
| Relinquished By: <u>[Signature]</u> Date/Time: <u>9-29-15</u> Company: <u>AGI</u> 1:30PM | Received By: <u>[Signature]</u> Date/Time: <u>10/1/15</u> Company: <u>SWCA</u> 15:08 |
| Relinquished By: <u>[Signature]</u> Date/Time: <u>10/20/15</u> Company: <u>SWCA</u> 17:00 | Received By: <u>Darkene Galloway</u> Date/Time: <u>10/21/15</u> Company: <u>AGI</u> 10:45 AM |



210 Executive Drive, Suite 1
Newark, DE USA 19702-3335
ph: 302-266-2428

AGI Project No.

ENV 01509

Site Name:

Comal & San Marcos Rivers oct

Site Location:

**AGI Soil Gas Sampling
Installation & Retrieval Log**

Company Name:

SWCA Environmental Consultants

Location:

Samples collected by:

Jennifer Moreland, Brittany Rios, Guy Rubio

* Optional or as needed

| SAMPLER SERIAL NO. | WELL ID | SAMPLE TYPE (Field Sample, Trip Blank, Field Blank, etc.) | INSTALLATION DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 13:00 | RETRIEVAL DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 15:00 | WATER QUALITY M | | |
|--------------------|----------|---|--|---|--|--|---|
| | | | | | INSTALLATION DEPTH FROM TOP OF CASING (feet) | DEPTH TO WATER FROM TOP OF CASING (feet) | SCREENED INTERVAL FROM TOP OF CASING (feet) |
| 00770685 | HCS410 | FIELD_SAMPLE | 10/6/15 12:19 | 10/20/15 10:32 | 3.25 | | |
| 00770686 | HCS420 | FIELD_SAMPLE | 10/6/15 12:37 | 10/20/15 10:42 | 1.1 | | |
| 00770687 | HCS430 | FIELD_SAMPLE | 10/6/15 11:56 | 10/20/15 10:16 | 2.1 | | |
| 00770688 | HCS440 | FIELD_SAMPLE | 10/6/15 12:47 | 10/20/15 10:54 | 1.9 | | |
| 00770689 | FDHCS440 | FIELD_SAMPLE | 10/6/15 12:47 | 10/20/15 10:54 | 1.9 | | |
| 00770690 | HCS460 | FIELD_SAMPLE | 10/6/15 13:03 | 10/20/15 11:03 | 2.65 | | |
| 00770691 | HSM410 | FIELD_SAMPLE | 10/6/15 13:45 | 10/20/15 11:54 | 1.5 | | |
| 00770692 | HSM420 | FIELD_SAMPLE | 10/6/15 13:59 | 10/20/15 12:04 | 2.9 | | |
| 00770693 | HSM430 | FIELD_SAMPLE | 10/6/15 14:13 | 10/20/15 12:10 | 0.55 | | |
| 00770694 | FDHSM430 | FIELD_SAMPLE | 10/6/15 14:13 | 10/20/15 12:10 | 0.55 | | |
| 00770695 | HSM440 | FIELD_SAMPLE | 10/6/15 14:28 | 10/20/15 12:22 | 5 | | |
| 00770696 | HSM450 | FIELD_SAMPLE | 10/6/15 14:44 | 10/20/15 12:38 | 2.1 | | |
| 00770697 | HSM460 | FIELD_SAMPLE | 10/6/15 14:55 | 10/20/15 12:53 | 0.8 | | |
| 00770698 | HSM470 | FIELD_SAMPLE | 10/6/15 15:12 | 10/20/15 13:03 | 1.8 | | |
| 00770699 | TB13 | TRIP_BLANK | 10/6/15 12:19 | 10/20/15 13:03 | NA | | |



**AGI Soil Gas Sampling
Installation & Retrieval L**

* Optional or as needed

| MONITORING | | | OBSERVATIONS/COMMENTS* (e.g., observations, location description, missing, pulled from well, etc. - as needed) | YES / NO | |
|--------------------|--|--|--|--|--------|
| SAMPLER SERIAL NO. | WATER TEMPERATURE AT MODULE DEPTH (deg C) | FLOW RATE THROUGH SCREEN: HIGH OR LOW (HIGH = flow > than 10 m/day; LOW = flow < 10 m/day) | | EVIDENCE OF LIQUID PETROLEUM HYDROCARBONS? | ODOR ? |
| 00770685 | 23.5 | | | | |
| 00770686 | 23.5 | | | | |
| 00770687 | 23.5 | | | | |
| 00770688 | 23.5 | | | | |
| 00770689 | 23.5 | | | | |
| 00770690 | 23.5 | | | | |
| 00770691 | 22 | | | | |
| 00770692 | 22 | | | | |
| 00770693 | 22 | | | | |
| 00770694 | 22 | | | | |
| 00770695 | 22 | | | | |
| 00770696 | 22 | | | | |
| 00770697 | 22 | | | | |
| 00770698 | 22 | | | | |
| 00770699 | | | | | |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770685 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 **Field ID:** HCS410

Installation Date: 10/6/2015 12:19:00PM

Retrieval Date: 10/20/2015 10:32:00AM

Date Analyzed: 10/27/2015 6:55:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | 0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.07 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770685 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS410

Installation Date: 10/6/2015 12:19:00PM

Retrieval Date: 10/20/2015 10:32:00AM

Date Analyzed: 10/27/2015 6:55:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770686 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS420

Installation Date: 10/6/2015 12:37:00PM

Retrieval Date: 10/20/2015 10:42:00AM

Date Analyzed: 10/27/2015 5:25:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.21 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770686 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS420

Installation Date: 10/6/2015 12:37:00PM

Retrieval Date: 10/20/2015 10:42:00AM

Date Analyzed: 10/27/2015 5:25:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770687 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS430

Installation Date: 10/6/2015 11:56:00AM

Retrieval Date: 10/20/2015 10:16:00AM

Date Analyzed: 10/27/2015 8:56:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.55 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770687 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS430

Installation Date: 10/6/2015 11:56:00AM

Retrieval Date: 10/20/2015 10:16:00AM

Date Analyzed: 10/27/2015 8:56:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770688 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS440

Installation Date: 10/6/2015 12:47:00PM

Retrieval Date: 10/20/2015 10:54:00AM

Date Analyzed: 10/27/2015 2:54:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.47 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770688 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS440

Installation Date: 10/6/2015 12:47:00PM

Retrieval Date: 10/20/2015 10:54:00AM

Date Analyzed: 10/27/2015 2:54:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770689 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: FDHCS440

Installation Date: 10/6/2015 12:47:00PM

Retrieval Date: 10/20/2015 10:54:00AM

Date Analyzed: 10/27/2015 3:24:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.46 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770689 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: FDHCS440

Installation Date: 10/6/2015 12:47:00PM

Retrieval Date: 10/20/2015 10:54:00AM

Date Analyzed: 10/27/2015 3:24:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770690 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS460

Installation Date: 10/6/2015 1:03:00PM

Retrieval Date: 10/20/2015 11:03:00AM

Date Analyzed: 10/27/2015 6:25:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.30 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770690 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS460

Installation Date: 10/6/2015 1:03:00PM

Retrieval Date: 10/20/2015 11:03:00AM

Date Analyzed: 10/27/2015 6:25:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770691 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM410

Installation Date: 10/6/2015 1:45:00PM

Retrieval Date: 10/20/2015 11:54:00AM

Date Analyzed: 10/27/2015 4:55:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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www.agisurveys.net

PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770691 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 **Field ID:** HSM410

Installation Date: 10/6/2015 1:45:00PM

Retrieval Date: 10/20/2015 11:54:00AM

Date Analyzed: 10/27/2015 4:55:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770692 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM420

Installation Date: 10/6/2015 1:59:00PM

Retrieval Date: 10/20/2015 12:04:00PM

Date Analyzed: 10/27/2015 8:26:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.08 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770692 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM420

Installation Date: 10/6/2015 1:59:00PM

Retrieval Date: 10/20/2015 12:04:00PM

Date Analyzed: 10/27/2015 8:26:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770693 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM430

Installation Date: 10/6/2015 2:13:00PM

Retrieval Date: 10/20/2015 12:10:00PM

Date Analyzed: 10/27/2015 2:23:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.94 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770693 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM430

Installation Date: 10/6/2015 2:13:00PM

Retrieval Date: 10/20/2015 12:10:00PM

Date Analyzed: 10/27/2015 2:23:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770694 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: FDHSM430

Installation Date: 10/6/2015 2:13:00PM

Retrieval Date: 10/20/2015 12:10:00PM

Date Analyzed: 10/27/2015 9:26:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 1.01 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770694 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: FDHSM430

Installation Date: 10/6/2015 2:13:00PM

Retrieval Date: 10/20/2015 12:10:00PM

Date Analyzed: 10/27/2015 9:26:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00770695 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM440

Installation Date: 10/6/2015 2:28:00PM

Retrieval Date: 10/20/2015 12:22:00PM

Date Analyzed: 10/27/2015 5:55:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.11 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770695 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM440

Installation Date: 10/6/2015 2:28:00PM

Retrieval Date: 10/20/2015 12:22:00PM

Date Analyzed: 10/27/2015 5:55:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

**FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA**

SAMPLER ID: 00770696 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM450

Installation Date: 10/6/2015 2:44:00PM

Retrieval Date: 10/20/2015 12:38:00PM

Date Analyzed: 10/27/2015 3:54:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.04 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



AMPLIFIED GEOCHEMICAL IMAGING, LLC

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Newark, DE 19702-3335 USA
ph: +1-302-266-2428
www.agisurveys.net

PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770696 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM450

Installation Date: 10/6/2015 2:44:00PM

Retrieval Date: 10/20/2015 12:38:00PM

Date Analyzed: 10/27/2015 3:54:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770697 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM460

Installation Date: 10/6/2015 2:55:00PM

Retrieval Date: 10/20/2015 12:53:00PM

Date Analyzed: 10/27/2015 10:26:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.08 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770697 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM460

Installation Date: 10/6/2015 2:55:00PM

Retrieval Date: 10/20/2015 12:53:00PM

Date Analyzed: 10/27/2015 10:26:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770698 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM470

Installation Date: 10/6/2015 3:12:00PM

Retrieval Date: 10/20/2015 1:03:00PM

Date Analyzed: 10/27/2015 9:56:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.09 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | 0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770698 **FIELD_SAMPLE**

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM470

Installation Date: 10/6/2015 3:12:00PM

Retrieval Date: 10/20/2015 1:03:00PM

Date Analyzed: 10/27/2015 9:56:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770699 TRIP_BLANK

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: TB13

Installation Date: 10/6/2015 12:19:00PM

Retrieval Date: 10/20/2015 1:03:00PM

Date Analyzed: 10/27/2015 4:25:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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www.agisurveys.net

PROJECT NUMBER: ENV 01509
SITE NAME: Comal & San Marcos Rivers
SITE ADDRESS: October Sampling

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770699 TRIP_BLANK

Matrix: WATER

Product: SPG0008

Dilution Factor: 1 Field ID: TB13

Installation Date: 10/6/2015 12:19:00PM

Retrieval Date: 10/20/2015 1:03:00PM

Date Analyzed: 10/27/2015 4:25:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151026-1

Reviewer: Ian McMullen

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |

AMPLIFIED GEOCHEMICAL IMAGING ANALYTICAL RESULTS
 210 EXECUTIVE DRIVE, SUITE 1, NEWARK, DE
 SWCA, SAN ANTONIO, TX
 AGI STANDARD TARGET COMPOUNDS
 ESTIMATED WATER CONCENTRATIONS
 COMAL AND SAN MARCOS RIVERS OCT 2015
 ORDER #01509

| DATAFILE | FIELD | DATE/ TIME | | DATE/ TIME | | DATE/ TIME | | DATE/ TIME | | | estimated | |
|----------------|----------|------------|-----|------------|-----|------------|-----|------------|-----|----|-----------|------------|
| NAME | ID | INSTALLED | | RETRIEVED | | RECEIVED | | ANALYZED | | DF | TPH, ug/L | MTBE, ug/L |
| Average RL= | | | | | | | | | | | 0.054 | 0.013 |
| 00770685 | HCS410 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.053 | <0.013 |
| 00770686 | HCS420 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.053 | <0.013 |
| 00770687 | HCS430 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.053 | <0.013 |
| 00770688 | HCS440 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.053 | <0.013 |
| 00770689 | FDHCS440 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.053 | <0.013 |
| 00770690 | HCS460 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.053 | <0.013 |
| 00770691 | HSM410 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.055 | <0.013 |
| 00770692 | HSM420 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.055 | <0.013 |
| 00770693 | HSM430 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.055 | <0.013 |
| 00770694 | FDHSM430 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.055 | <0.013 |
| 00770695 | HSM440 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.055 | <0.013 |
| 00770696 | HSM450 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.055 | <0.013 |
| 00770697 | HSM460 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.055 | <0.013 |
| 00770698 | HSM470 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.055 | <0.013 |
| 00770699 | TB13 | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.054 | <0.013 |
| BLK_ENV-273757 | | 10/6/2015 | CDT | 10/20/2015 | CDT | 10/21/2015 | EST | 10/27/2015 | EST | 1 | <0.054 | <0.013 |

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| DATAFILE | | | | | | | | | |
|----------------|--------------|-------------|--------------|-------------|--------------|-------------|------------|------------|-----------|
| NAME | t12DCE, ug/L | 11DCA, ug/L | c12DCE, ug/L | CHCl3, ug/L | 111TCA, ug/L | 12DCA, ug/L | BENZ, ug/L | CCl4, ug/L | TCE, ug/L |
| Average RL= | 0.008 | 0.007 | 0.007 | 0.007 | 0.005 | 0.008 | 0.006 | 0.005 | 0.006 |
| 00770685 | <0.007 | <0.007 | <0.007 | 0.008 | <0.005 | <0.008 | <0.006 | <0.004 | <0.006 |
| 00770686 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.004 | <0.006 |
| 00770687 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.004 | <0.006 |
| 00770688 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.004 | <0.006 |
| 00770689 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.004 | <0.006 |
| 00770690 | <0.007 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.004 | <0.006 |
| 00770691 | <0.008 | <0.007 | <0.007 | <0.008 | <0.005 | <0.008 | <0.006 | <0.005 | <0.006 |
| 00770692 | <0.008 | <0.007 | <0.007 | <0.008 | <0.005 | <0.008 | <0.006 | <0.005 | <0.006 |
| 00770693 | <0.008 | <0.007 | <0.007 | <0.008 | <0.005 | <0.008 | <0.006 | <0.005 | <0.006 |
| 00770694 | <0.008 | <0.007 | <0.007 | <0.008 | <0.005 | <0.008 | <0.006 | <0.005 | <0.006 |
| 00770695 | <0.008 | <0.007 | <0.007 | <0.008 | <0.005 | <0.008 | <0.006 | <0.005 | <0.006 |
| 00770696 | <0.008 | <0.007 | <0.007 | <0.008 | <0.005 | <0.008 | <0.006 | <0.005 | <0.006 |
| 00770697 | <0.008 | <0.007 | <0.007 | <0.008 | <0.005 | <0.008 | <0.006 | <0.005 | <0.006 |
| 00770698 | <0.008 | <0.007 | <0.007 | <0.008 | <0.005 | <0.008 | <0.006 | <0.005 | <0.006 |
| 00770699 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | <0.006 |
| BLK_ENV-273757 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | <0.005 | <0.006 |

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| DATAFILE | | | | | | | | | |
|----------------|--------------|-----------|-----------|-----------|--------------|-----------------|--------------|-------------|------------|
| NAME | 112TCA, ug/L | TOL, ug/L | OCT, ug/L | PCE, ug/L | CIBENZ, ug/L | 1112TetCA, ug/L | ETBENZ, ug/L | mpXYL, ug/L | oXYL, ug/L |
| Average RL= | 0.010 | 0.006 | 0.005 | 0.005 | 0.006 | 0.007 | 0.005 | 0.005 | 0.005 |
| 00770685 | <0.009 | <0.006 | <0.004 | 0.015 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 |
| 00770686 | <0.009 | <0.006 | <0.004 | 0.040 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 |
| 00770687 | <0.009 | <0.006 | <0.004 | 0.095 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 |
| 00770688 | <0.009 | <0.006 | <0.004 | 0.083 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 |
| 00770689 | <0.009 | <0.006 | <0.004 | 0.081 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 |
| 00770690 | <0.009 | <0.006 | <0.004 | 0.055 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 |
| 00770691 | <0.010 | <0.006 | <0.005 | <0.005 | <0.006 | <0.007 | <0.005 | <0.005 | <0.006 |
| 00770692 | <0.010 | <0.006 | <0.005 | 0.019 | <0.006 | <0.007 | <0.005 | <0.005 | <0.006 |
| 00770693 | <0.010 | <0.006 | <0.005 | 0.160 | <0.006 | <0.007 | <0.005 | <0.005 | <0.006 |
| 00770694 | <0.010 | <0.006 | <0.005 | 0.171 | <0.006 | <0.007 | <0.005 | <0.005 | <0.006 |
| 00770695 | <0.010 | <0.006 | <0.005 | 0.025 | <0.006 | <0.007 | <0.005 | <0.005 | <0.006 |
| 00770696 | <0.010 | <0.006 | <0.005 | 0.009 | <0.006 | <0.007 | <0.005 | <0.005 | <0.006 |
| 00770697 | <0.010 | <0.006 | <0.005 | 0.019 | <0.006 | <0.007 | <0.005 | <0.005 | <0.006 |
| 00770698 | <0.010 | <0.006 | <0.005 | 0.020 | <0.006 | <0.007 | <0.005 | <0.005 | <0.006 |
| 00770699 | <0.010 | <0.006 | <0.005 | <0.005 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 |
| BLK_ENV-273757 | <0.010 | <0.006 | <0.005 | <0.005 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 |

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| DATAFILE | | | | | | | | |
|----------------|-----------------|--------------|--------------|-------------|-------------|-------------|-------------|------------|
| NAME | 1122TetCA, ug/L | 135TMB, ug/L | 124TMB, ug/L | 13DCB, ug/L | 14DCB, ug/L | 12DCB, ug/L | UNDEC, ug/L | NAPH, ug/L |
| Average RL= | 0.011 | 0.005 | 0.005 | 0.006 | 0.006 | 0.006 | 0.020 | 0.016 |
| 00770685 | <0.010 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.020 | <0.016 |
| 00770686 | <0.010 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.020 | <0.016 |
| 00770687 | <0.010 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.020 | <0.016 |
| 00770688 | <0.010 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.020 | <0.016 |
| 00770689 | <0.010 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.020 | <0.016 |
| 00770690 | <0.010 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.020 | <0.016 |
| 00770691 | <0.011 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.021 | <0.017 |
| 00770692 | <0.011 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.021 | <0.017 |
| 00770693 | <0.011 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.021 | <0.017 |
| 00770694 | <0.011 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.021 | <0.017 |
| 00770695 | <0.011 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.021 | <0.017 |
| 00770696 | <0.011 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.021 | <0.017 |
| 00770697 | <0.011 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.021 | <0.017 |
| 00770698 | <0.011 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.021 | 0.017 |
| 00770699 | <0.011 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.020 | <0.016 |
| BLK_ENV-273757 | <0.011 | <0.005 | <0.005 | <0.006 | <0.006 | <0.006 | <0.020 | <0.016 |

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| | estimated | | estimated | estimated | estimated | estimated |
|----------------|--------------|---------------|----------------------|----------------|--------------------|----------------|
| DATAFILE | | | | | | |
| NAME | TRIDEC, ug/L | 2MeNAPH, ug/L | Acenaphthylene, ug/L | PENTADEC, ug/L | Acenaphthene, ug/L | Fluorene, ug/L |
| Average RL= | 0.020 | 0.014 | 0.014 | 0.020 | 0.014 | 0.014 |
| 00770685 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00770686 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00770687 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00770688 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00770689 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00770690 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00770691 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770692 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770693 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770694 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770695 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770696 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770697 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770698 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770699 | <0.020 | <0.014 | <0.014 | <0.020 | <0.014 | <0.014 |
| BLK_ENV-273757 | <0.020 | <0.014 | <0.014 | <0.020 | <0.014 | <0.014 |

KEY TO DATA TABLE

UNITS

| | |
|-------------------|--|
| µg | micrograms, relative mass value |
| µg/m ³ | micrograms per cubic meter; estimated soil gas concentration |
| µg/L | micrograms per Liter; calculated water concentration |

DATA QUALIFIERS

| | |
|---|--|
| > | greater than; value exceeds calibration range, estimated value |
| < | less than; compound value is below the LOD and RL |
| J | mass value below LOQ or RL, but above LOD, estimated mass value |
| E | mass value exceeds upper calibration level, estimated mass value |
| Q | one or more quality control parameters failed for the compound |

ABBREVIATIONS

| | |
|--------|--|
| AVG RL | average reporting limit; calculated based on individual field sample RLs |
| LOD | limit of detection |
| LOQ | limit of quantification |
| MDL | method detection limit |
| RL | reporting limit |

| | | | |
|-------------|---|----------|--|
| 1112TetCA | 1,1,1,2-tetrachloroethane | CIBENZ | chlorobenzene |
| 111TCA | 1,1,1-trichloroethane | ct12DCE | cis- & trans-1,2-dichloroethene |
| 1122TetCA | 1,1,2,2-tetrachloroethane | EtBENZ | ethylbenzene |
| 112TCA | 1,1,2-trichloroethane | mpXYL | m-, p-xylene |
| 11DCA | 1,1-dichloroethane | MTBE | methyl t-butyl ether |
| 11DCE | 1,1-dichloroethene | NAPH | naphthalene |
| 124TMB | 1,2,4-trimethylbenzene | OCT | octane |
| 12DCA | 1,2-dichloroethane | oXYL | o-xylene |
| 12DCB | 1,2-dichlorobenzene | PCE | tetrachloroethene |
| 135TMB | 1,3,5-trimethylbenzene | PENTADEC | pentadecane |
| 13DCB | 1,3-dichlorobenzene | PHEN | phenanthrene |
| 14DCB | 1,4-dichlorobenzene | t12DCE | trans-1,2-dichloroethene |
| 2MeNAPH | 2-methyl naphthalene | TCE | trichloroethene |
| BENZ | benzene | TMBs | combined masses of 1,3,5-trimethylbenzene and 1,2,4-trimethylbenzene |
| BTEX | combined masses of benzene, toluene, ethylbenzene, and total xylenes (Gasoline Range Aromatics) | TOL | toluene |
| C11,C13&C15 | combined masses of undecane, tridecane, and pentadecane (C11+C13+C15) (Diesel Range Alkanes) | TPH | total petroleum hydrocarbons |
| c12DCE | cis-1,2-dichloroethene | TRIDEC | tridecane |
| CCl4 | carbon tetrachloride | UNDEC | undecane |
| CHC13 | chloroform | VC | vinyl chloride |

SUMMARY OF SAMPLING RATE CALIBRATION FOR AGI UNIVERSAL SAMPLER IN AQUEOUS PHASES

INTRODUCTION:

The Amplified Geochemical Imaging, LLC (AGI) passive vapor sampler is designed to be used for soil gas, water, sediment pore water, and air sampling. This document describes the process used to calibrate the sampler's compound specific sampling or uptake rates in aqueous phases.

Sampling rates are measured following AGI's "Standard Practice for Determining the Sampling Rate of Passive Diffusion Samplers in Various Environmental Media": SPG-SOP-0493. Rates are used to calculate dissolved phase concentrations of volatile and semi-volatile contaminants in water. The calibration process is summarized in three parts: Part 1: shallow water, Part 2: deep water, and Part 3: sediment.

PURPOSE:

The purpose of this document is to:

1. Summarize the test protocol,
2. Summarize the methodology for analysis of data,
3. Present general results for generating concentration calibration of the AGI Universal Sampler

Principle of Operation of the AGI Sampler

The AGI Universal Sampler is designed with solid adsorbents enclosed inside a tubular microporous PTFE membrane. When placed in water, the pores and hydrophobic nature of the PTFE keep liquid water from entering the membrane until a water head of about 34 feet is reached. The membrane will not keep water vapor from entering but the adsorbents are very hydrophobic and through testing validated to be unaffected by this moisture vapor. In shallow water, <34', volatile and semi-volatile compounds will partition from the dissolved water into the air phase in the PTFE membrane according to Henry's Law. This partitioning is instantaneous and within seconds-minutes, the compound is adsorbed by the adsorbent inside the sealed tube. Because the diffusivity in air is about 10,000 times higher than the diffusivity in water, the sampling rate is controlled by the water contact area of the membrane that allows the Henry's Law effect to occur. This contact area is set by the membrane diameter and length of the sealed tube, which is fixed in AGI's manufacturing process.

Henry's law as well as diffusivity, which are fundamentally incorporated into the sampling rate, are affected by temperature, T , and follow an Arrhenius equation $H_T = H_r \times \exp\left(\frac{-E_a/R}{1/T_r - 1/T}\right)$. Because a 5°C temperature change can make a 15% change in sampling rate, the temperature of the sampled water should be known to get the most precise concentration.

The membrane pore size is also small enough that colloidal particles and microbes cannot pass through the membrane. This keeps the adsorbent from getting contaminated and eliminates any need to add preservative or chilling during storage or transportation.

When the water pressure exceeds the water entry pressure of the membrane, about 34 feet of water, the water becomes in contact with the solid adsorbent. Under this condition the compounds in the water will partition from the water into the solid. The partitioning coefficient, K_{AW} , can be approximated by the octanol-water coefficient, K_{OW} , but has been measured more precisely in the lab for AGI's specific solid adsorbent. The sampling rate is the product of the sampling rate at <34' of water and the K_{AW} .

In sediment, the sampler measures pore-water concentration, which is generally agreed to be the preferred measurement as it is more indicative of bioavailability. In sediment the volumetric

availability of water to the sampler is reduced by the volume fraction solids in the sediment, which typically varies from zero to 35%, but can be as high as 73% in well packed and broad particle size distribution sediments. As a result, sampling rates in sediment are multiplied by the fraction pore water in the sediment to determine concentration.

PART 1: Calibration in shallow water

Part 1 summarizes the work in shallow water generating calibration data, evaluating the physical and chemical factors affecting the sampling rate, and measurement of the actual sampling rates or regression calibration equations needed to determine concentrations.

Sample Generation in water

In this calibration work, solutions of analytes at known concentrations were formulated in clean 4 liter smoked glass jugs by injecting microliter measured amounts of environmental standards using a calibrated syringe into pure or deionized water and stirring for a minimum of 2 hours but generally overnight. Headspace in the jugs was minimized and generally less than 1% by volume during the tests. Jugs were temperature controlled by placing them in a water filled cooler, chilled via a copper tubing loop in the cooler. Temperature was measured with a certified digital temperature gauge and an average value used for each temperature experiment.

AGI samplers were weighted so they won't float and placed in the jugs at time zero. They were removed at various intervals to generate samples along with duplicates that showed mass increasing with exposure time. The sampler exposure time was selected to span minutes to hours and was generally reduced for high concentration tests to maintain uptake with time in roughly the linear dynamic range. Samplers were removed and dried with a paper towel and returned to their original container for analysis. They were analyzed by AGI's 8260C (SPG-WI-318 or SPG-WI-10028) method in duplicate, which is based on EPA SW846 Method 8260C.

Water samples were also taken and measured at an outside accredited lab using EPA SW846 Method 8260B. The concentrations agreed well with the calculated concentrations based on the standard certification, jug volume, and syringe injection. The variability of the outside lab 8260B values were found to be high, so for the sampling rate calculations we used the concentrations based on syringe dosing.

Calibrations were run at five concentrations, nominally at 6, 24, 118, 590, 1420 ug/L and five temperatures nominally at 5, 10, 15, 20, and 25 degrees centigrade. Samples were taken at 4 different exposure times. Samples were run in duplicate. A total of 176 data points were generated using 28 compounds from AGI's standard compounds list. Tridecane and pentadecane were not evaluated due to their very low solubility in water. In addition, another 23 compounds were tested using an 8260 liquid standard at nominal concentrations of 0.5, 1.0, 5.0, 15, 95, and 470 ug/L at a temperature typical of groundwater, 15°C. This is a living calibration and as additional data are generated, they may be qualified and added to this data set to improve the precision of the sampling rate calibration and broaden the compound list.

Key Variable Effects

As expected from theory, at short to moderate exposure times, mass will increase roughly linearly proportional to exposure time, as well as proportional to concentration, and exponentially with temperature following Arrhenius law. Temperature affects the Henry's law as well as diffusivity in water. Sampling rate is generally independent of concentration and time at mass values significantly below saturation. In the following sections we have characterized the sampling rate for each compound as affected by temperature and also developed calibrations using regression which account for the minor impact of time, and mass.

Concentration using Simple Sampling Rate Determination

A simple way to determine concentration is to measure mass on the AGI sampler, divide by exposure time, and divide by sampling rate, SR.

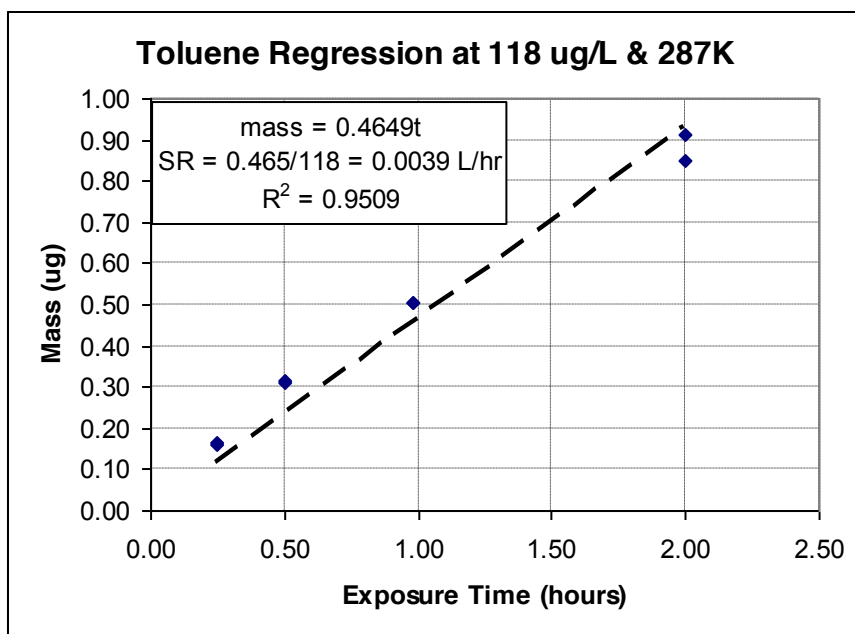
$$\text{Conc [ug/L]} = \text{mass/time/SR} \quad (1)$$

The sampling rate can be determined via measurements of mass versus time at a known concentration and temperature according to the following modification of equation (1).

$$\text{SR} = \text{mass/time/concentration} \quad (2)$$

Sampling rates in L/hr were determined by measuring the trend or regression mass uptake versus time and dividing by the concentration. A measurement like this will use 8 data points (4 times x 2 samples). Such a sampling rate can be measured at any concentration and temperature.

The chart to the right shows a plot of mass versus time for water at 118 ug/L and 287K (actual data from a single run). Slope of 0.465 ug/hr divided by the concentration of 118 ug/L yields a sampling rate, SR, of 0.0039 L/hr.



SR's typically range from about 0.004 to 0.007 L/hr at 15°C. Table A shows SR's measured for our standard compound list at 5 temperatures.

Rigorous Concentration using Regression

A preferred method for determining concentration that will yield improved accuracy over a wide range of concentrations, exposure times, and temperatures is to use all data in a regression analysis, which allows adjustments for the minor non-linear influences of mass and time as well as the effects of temperature. This step is done by regressing equation (1) or a universal version of equation (1):

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRo} * \exp(-\text{Ea}/\text{R}/\text{T})] \quad (3)$$

The subtle non-linear effects of mass and time will be evident in the deviation of coefficients b and d from 1.0. This regression generates four constants b, d, SRo, and $-\text{Ea}/\text{R}$ by regressing $\ln(\text{conc})$ versus $\ln(\text{mass})$, $\ln(\text{time})$, $1/\text{temp}$. These four constants can be used to determine concentration via the equation:

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRo} * \exp(-\text{Ea}/\text{R}(1/\text{T}))] \quad (4)$$

Where conc is in ug/L, mass is in ug, time in hours, T in degrees Kelvin.

Equation (4) can be also expressed at a reference temperature, Tr , such as 15°C by

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRr} * \exp(-\text{Ea}/\text{R}(1/\text{Tr} - 1/\text{T}))] \quad (5)$$

This step allows sampling rates, SRr, at any reference temperature, Tr , and for any analyte to easily be compared. The values of SRr at 293.14K can be found in Table A.

When sampling times are between 0 and 4 hours, using the 4 constant equation (5) is preferred. For concentrations from about 5 to 1500 ug/L one hour exposure times generally give the lowest error, typically with average error of 6-20% and with total error range of 12%-32%. For low concentrations where sampling times are greater than 4 hours, it is preferred to use equation (1) to avoid unrealistic effects from the coefficient d or to set d to 1.0. In such a case SR in equation (1) can be substituted with $[\text{SRr} * \exp(-\text{Ea}/\text{R}(1/\text{Tr} - 1/\text{T}))]$ to use an SR representative of the well temperature, T.

The chart to the right is a plot of the calculated concentration from the 4 constant regression compared to the dosed concentration. Agreement is excellent for the 176 data points.

However, there does appear to be a slight high bias of 8.6% over the full range of this data, although it is well within acceptable limits of variability.

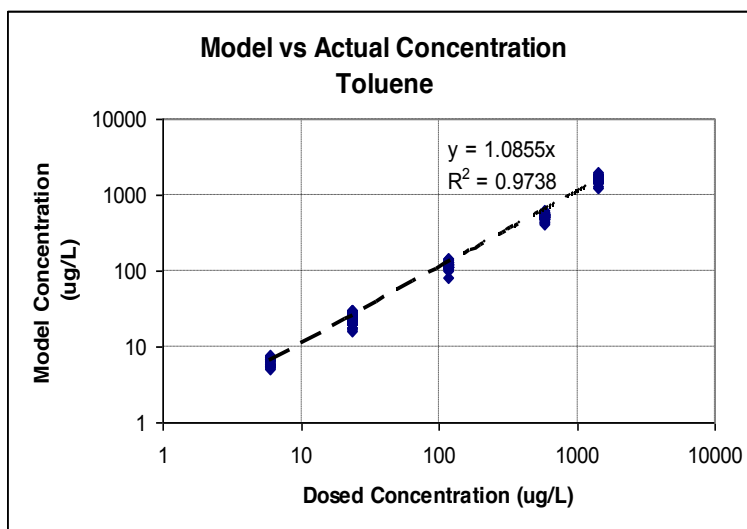


Table B shows the tabulated summary of the 4 constants regression with Rsq values and error estimates for the 4 constants for each analyte. Most regression Rsq values are 0.99 or greater for each analyte. In general, $-E_a/R$ is about 2400 \pm 400, b is about 0.9, d is about -0.75, and SR(15°C) ranges from .004 L/hr to 0.007 L/hr increasing with MW of the compound.

Error Estimates

The error in the water concentration values will depend on both the error in mass from the analytical method as well as the error in the concentration calibration. Table C shows the error in the mass values from the 8260C low sensitivity method.

The standard error of the regression and standard errors of the constants can be found in table B. For each compound we have measured the error between the derived concentration and the actual concentration. The error tends to be lowest at our recommended exposure time of one hour as shown by the example for Toluene to the right.

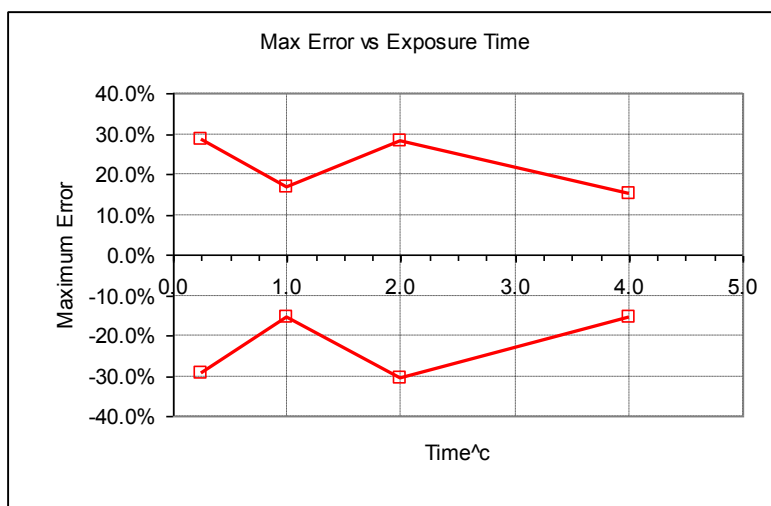
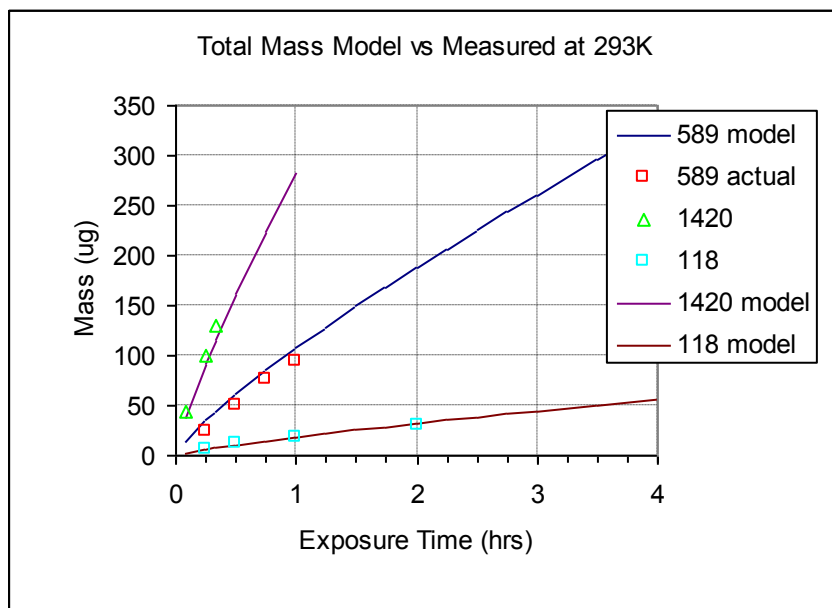


Table D shows the total average error in water concentration by compound as well as the low and high error. The average ranges from about 6% to 20%, which is similar to the analytical method errors. The low and high errors range from 12% to 32% and include contribution from measurement errors in both time and temperature.

Sorbent Saturation

As mass increases on a solid sorbent and approaches saturation, reverse diffusion can occur causing the sampling rate to drop. Eventually the mass level will reach a maximum steady state value at any concentration. A rate of mass uptake with time that deviates significantly from linear, indicates that sorbent saturation could be an issue. When using equation (1), staying in the linear range to avoid the effects of adsorbent saturation is important. We recommend keeping the total mass on the sampler below 50 ug or flagging when this is exceeded.

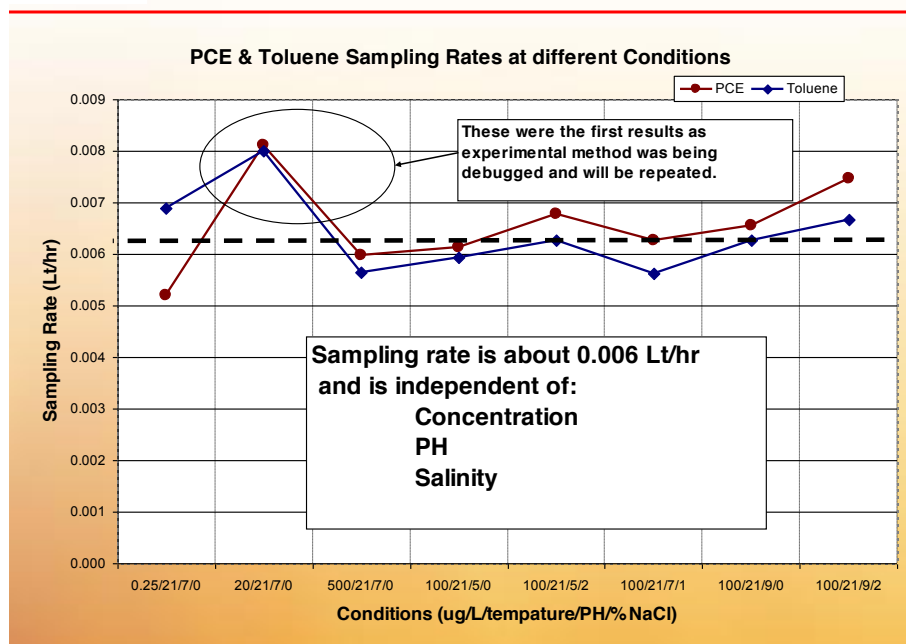
The 4 constant regression accounts for some of the non linearity allowing good accuracy at higher mass levels. From the experimental data we have found this safe range can be extended to 100 ug or higher as shown in the chart below. This chart compares total mass of all compounds (excluding heavy alkanes, which have solubility issues) versus time in comparison to that predicted from the 4-constant concentration equation.



Effect of PH and Salinity

Because neither PH nor salinity is known to have a significant impact on Henry's law or diffusivity in water, we did not expect them to have a significant impact on sampling rate. To confirm this, experiments were run varying PH from 5 to 9 and NaCl content from 0 to 2%. The chart below shows no significant impact for combinations of PH and NaCl content over this range on the sampling rate of toluene in water at 21°C.

Checked for Effects of PH & Salinity

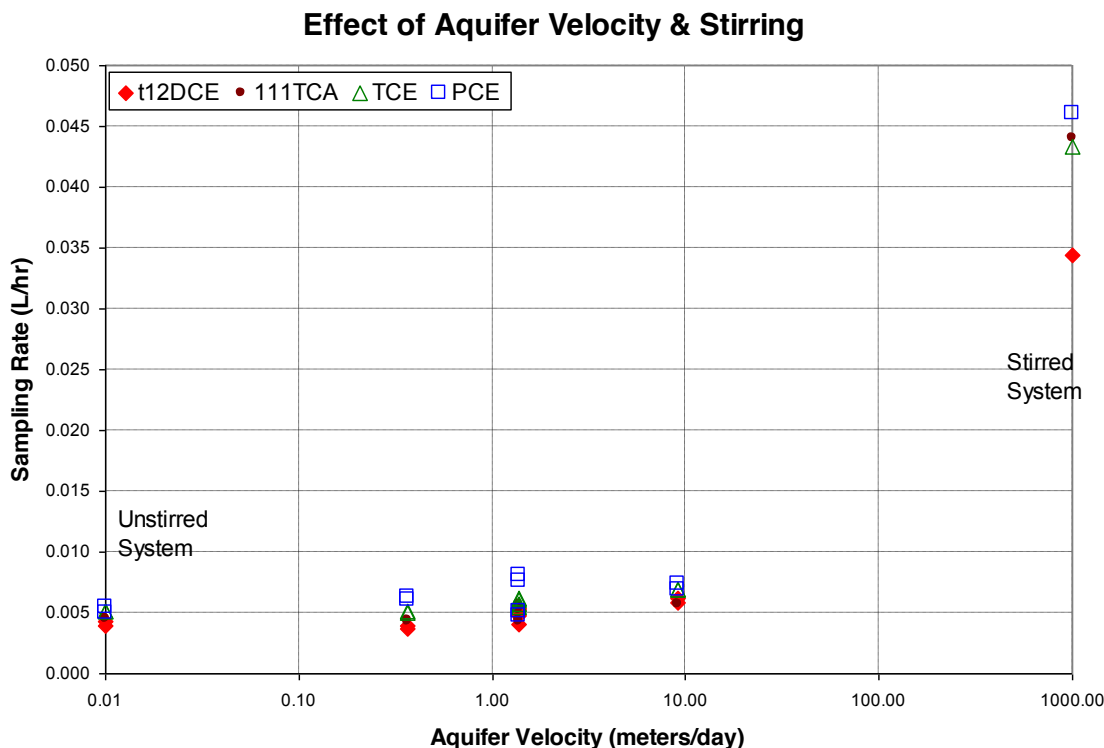


Impact of Aquifer Velocity

The velocity in most aquifers is quite slow, typically a meter/day or less. Occasionally water flow could be much higher such as encountered in karst aquifers, streams or rivers. Mass transfer coefficients are higher in high flow conditions, which will lead to higher sampling rates. We validated that a highly stirred system had sampling rates about 10 times higher than those that were non-stirred. We decided to evaluate the effect of aquifer velocity.

A test apparatus was built comprising a 3" PVC pipe tee filled with clean sand in each of the horizontal straight legs and screened to leave the center open. A test solution was run through this system using a variable flow pump and AGI samplers were placed into the simulated well through the vertical leg of the tee. Tests were run to examine the effect of velocity by varying the pumping rate and hence water velocity.

The chart below shows no significant effect of aquifer velocity up to a speed of about 10 meters/day. At velocities significantly above this, similar to a stirred system, sampling rates are about 10 times higher.

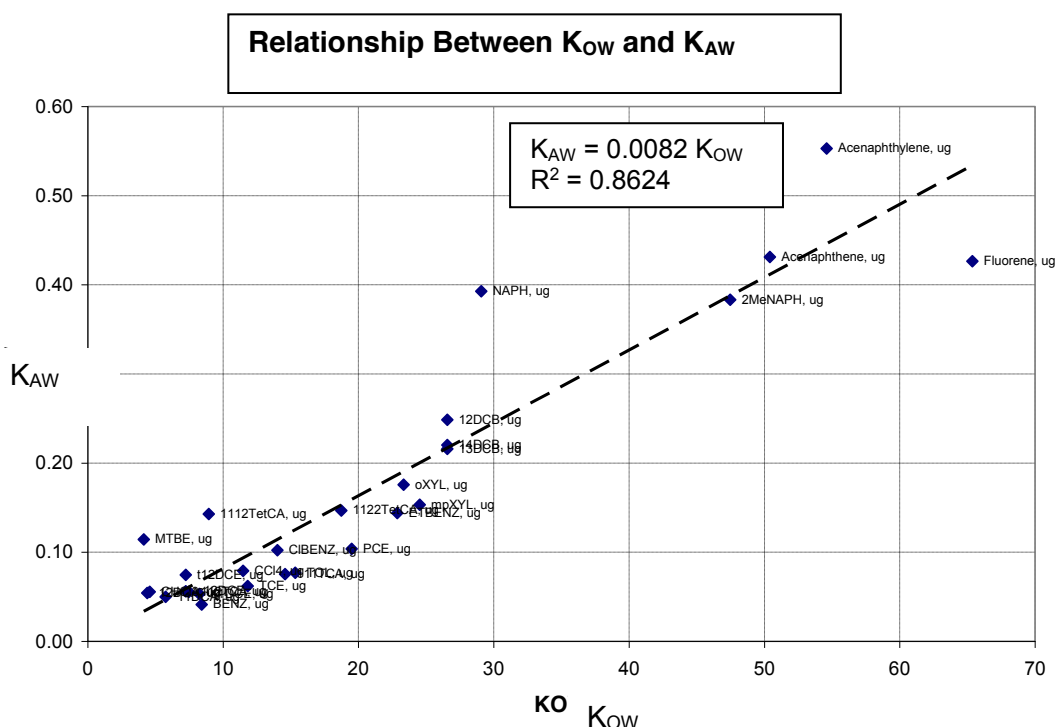


Part 2: Calibration in Deep (>34') water

Part 2 describes the effect of deep water on the AGI sampler and summarizes the effects on sampling rate and concentration measurement.

When the water pressure exceeds the water entry pressure of the membrane, about 34 feet of water, the water becomes in direct contact with the solid adsorbent. Under this condition the compounds in the water will partition from the water into the solid. The partitioning coefficient is closely related to the octanol-water coefficient, K_{OW} , but has been measured more precisely in the lab for AGI's specific solid adsorbent, K_{AW} . The sampling rate for deep water is the product of the sampling rate at <34' of water and the K_{AW} .

Measurement of the K_{AW} was done in a one liter stainless steel vessel pressurized with nitrogen to simulate water heads above 34' of water. Pressures of up to 465 psig or 200' of water head were used. The sampling rate change was the same at all pressures above 34' of water. The K_{AW} was determined as the ratio between the mass or sampling rate above 34' of head to the rate at <34' of head and is shown in the chart below.



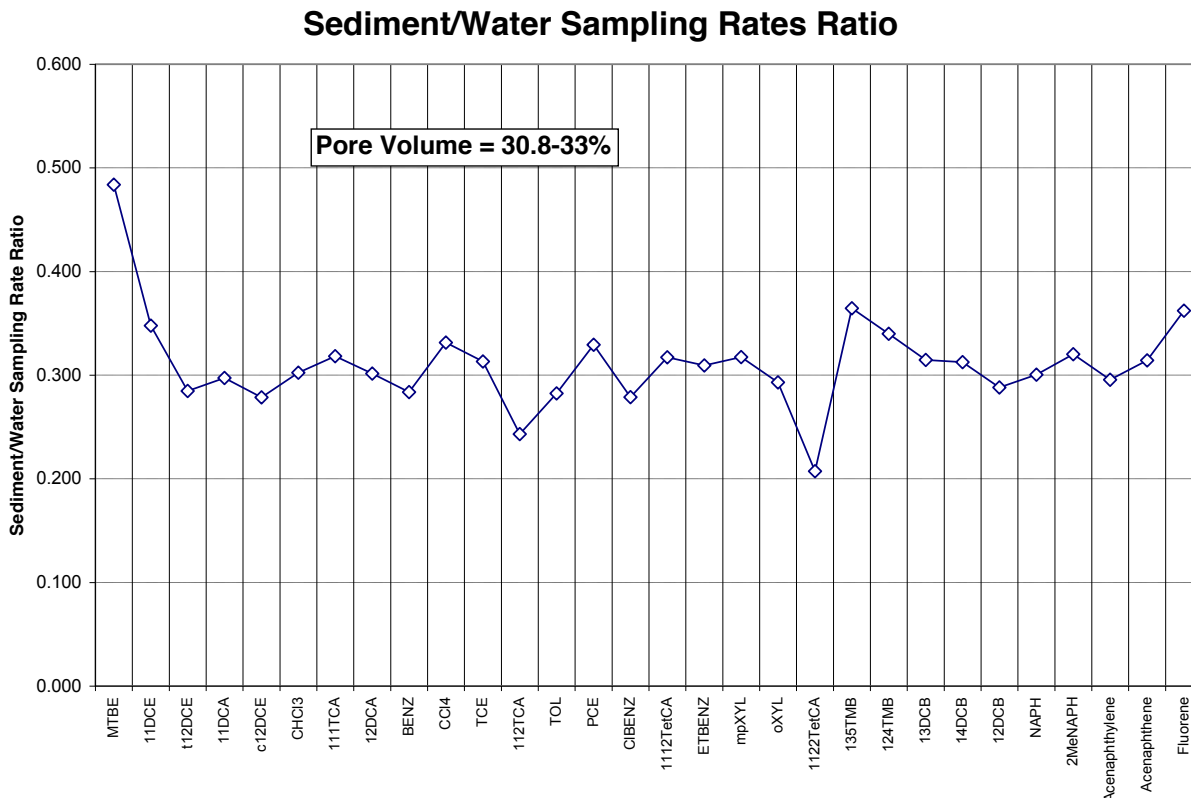
Part 3: Calibration in Sediment

Part 3 describes the effect of sediment solids or sediment pore volume on the sampling rate and concentration measurement.

In sediment, the sampler measures pore water concentration, which is generally agreed to be the preferred measurement as it is more indicative of bioavailability. In sediment the volumetric availability of water to the sampler is reduced by the volume fraction solids in the sediment. As a result sampling rates in sediment are multiplied by the fraction pore water to determine

concentration. Pore water fraction can range from 1.0 for water without sediment to as low as 0.25. Typically most sediments have pore fractions of 0.9 to 0.65.

A sampling rate study was done with water and with water added into a well-packed sorted sand. Pore water fraction in this test was measured between 30.8% and 33% by volume. Below is a plot of the ratio of sampling rates measured in the sediment to open water. The average ratio is equal to the pore water fraction confirming that sampling rate in sediment is on average equal to the product of pore water fraction times the sampling rate in water.



Summary

The AGI Sampler can be used to determine the concentration of volatile and semi-volatile compounds in a water phase. This requires knowing the exposure time and water temperature. It also requires knowing if the sample is above or below 34' of water head and if the water has a velocity above 10 meters/day. Regressions of large amounts of data were used to generate a four constant equation to generate concentration values in water. Potential error in the concentration values is excellent typically less than 25%.

TABLE A
WATER SAMPLING RATES STANDARD LIST

| | SRr 293.14 | SR @ 277.54 | SR @ 282.44 | SR @ 287.84 | SR @ 293.24 | SR @ 298.94 |
|-------------------|----------------------|-----------------------|-----------------------|-----------------------|-----------------------|-----------------------|
| MTBE | 0.0025 | 0.0014 | 0.0016 | 0.0018 | 0.0022 | 0.0029 |
| t12DCE | 0.0043 | 0.0028 | 0.0028 | 0.0027 | 0.0037 | 0.0048 |
| 11DCA | 0.0047 | 0.0031 | 0.0033 | 0.0033 | 0.0039 | 0.0052 |
| c12DCE | 0.0046 | 0.0030 | 0.0031 | 0.0031 | 0.0038 | 0.0051 |
| CHCl3 | 0.0046 | 0.0030 | 0.0031 | 0.0031 | 0.0038 | 0.0051 |
| 111TCA | 0.0066 | 0.0043 | 0.0047 | 0.0047 | 0.0056 | 0.0076 |
| 12DCA | 0.0045 | 0.0029 | 0.0029 | 0.0030 | 0.0036 | 0.0050 |
| BENZ | 0.0050 | 0.0031 | 0.0034 | 0.0035 | 0.0042 | 0.0056 |
| CCl4 | 0.0068 | 0.0044 | 0.0048 | 0.0047 | 0.0058 | 0.0080 |
| TCE | 0.0052 | 0.0030 | 0.0034 | 0.0034 | 0.0043 | 0.0058 |
| 112TCA | 0.0043 | 0.0027 | 0.0027 | 0.0028 | 0.0034 | 0.0048 |
| TOL | 0.0056 | 0.0034 | 0.0039 | 0.0039 | 0.0047 | 0.0062 |
| OCT | 0.0064 | 0.0046 | 0.0050 | 0.0040 | 0.0058 | 0.0089 |
| PCE | 0.0061 | 0.0036 | 0.0043 | 0.0043 | 0.0051 | 0.0069 |
| CIBENZ | 0.0054 | 0.0033 | 0.0039 | 0.0040 | 0.0045 | 0.0059 |
| 1112TetCA | 0.0061 | 0.0037 | 0.0042 | 0.0044 | 0.0050 | 0.0065 |
| EtBENZ | 0.0060 | 0.0037 | 0.0045 | 0.0044 | 0.0052 | 0.0069 |
| mpXYL | 0.0064 | 0.0039 | 0.0048 | 0.0046 | 0.0055 | 0.0072 |
| oXYL | 0.0066 | 0.0041 | 0.0050 | 0.0048 | 0.0057 | 0.0074 |
| 1122TetCA | 0.0044 | 0.0027 | 0.0029 | 0.0031 | 0.0036 | 0.0046 |
| 135TMB | 0.0079 | 0.0046 | 0.0059 | 0.0056 | 0.0071 | 0.0093 |
| 124TMB | 0.0078 | 0.0046 | 0.0060 | 0.0055 | 0.0071 | 0.0092 |
| 13DCB | 0.0072 | 0.0041 | 0.0055 | 0.0053 | 0.0063 | 0.0080 |
| 14DCB | 0.0071 | 0.0040 | 0.0054 | 0.0052 | 0.0062 | 0.0079 |
| 12DCB | 0.0070 | 0.0040 | 0.0053 | 0.0051 | 0.0060 | 0.0076 |
| UNDEC | | 0.0026 | 0.0024 | 0.0020 | 0.0031 | 0.0029 |
| NAPH | | 0.0041 | 0.0056 | 0.0054 | 0.0064 | 0.0081 |
| TRIDEC | | | | | | |
| 2MeNAPH | | 0.0043 | 0.0066 | 0.0066 | 0.0080 | 0.0108 |
| PENTADEC | | | | | | |
| Total mass | 0.1177 | 0.0822 | 0.1339 | 0.1334 | 0.1773 | 0.1981 |

Notes:

Values in L/hr

Total mass does not include UNDEC, TRIDEC, PENTADEC (28 compounds)

TABLE B
4 CONSTANT REGRESSION OUTPUT

| | Adjusted Rsq | Standard Error | ln(SR0) | b | -Ea/R | d | Std Error ln(SR0) | Std Error b | Std Error - Ea/R | Std Error d |
|-------------------|-------------------------|---------------------------|----------------|----------|--------------|----------|----------------------------------|----------------------------|-------------------------------------|----------------------------|
| MTBE | 0.997 | 0.0960 | -3.217 | 0.981 | 2704 | -0.709 | 0.2881 | 0.0062 | 83 | 0.0082 |
| t12DCE | 0.992 | 0.1659 | -1.877 | 0.905 | 2147 | -0.760 | 0.4971 | 0.0100 | 144 | 0.0138 |
| 11DCA | 0.995 | 0.1272 | -1.346 | 0.916 | 1965 | -0.746 | 0.3809 | 0.0077 | 110 | 0.0106 |
| c12DCE | 0.995 | 0.1299 | -1.905 | 0.911 | 2137 | -0.751 | 0.3892 | 0.0078 | 112 | 0.0109 |
| CHCl3 | 0.996 | 0.1260 | -1.841 | 0.912 | 2118 | -0.748 | 0.3776 | 0.0076 | 109 | 0.0105 |
| 111TCA | 0.995 | 0.1279 | -2.684 | 0.902 | 2259 | -0.761 | 0.3836 | 0.0076 | 111 | 0.0106 |
| 12DCA | 0.995 | 0.1263 | -2.161 | 0.908 | 2218 | -0.746 | 0.3786 | 0.0076 | 109 | 0.0106 |
| BENZ | 0.995 | 0.1323 | -2.207 | 0.920 | 2198 | -0.754 | 0.3965 | 0.0080 | 114 | 0.0110 |
| CCl4 | 0.994 | 0.1405 | -3.121 | 0.889 | 2379 | -0.776 | 0.4220 | 0.0083 | 122 | 0.0116 |
| TCE | 0.992 | 0.1655 | -3.338 | 0.900 | 2522 | -0.772 | 0.4969 | 0.0099 | 144 | 0.0137 |
| 112TCA | 0.995 | 0.1264 | -2.412 | 0.896 | 2302 | -0.724 | 0.3790 | 0.0075 | 109 | 0.0107 |
| TOL | 0.994 | 0.1426 | -2.873 | 0.916 | 2364 | -0.756 | 0.4281 | 0.0087 | 124 | 0.0119 |
| OCT | 0.938 | 0.4698 | -5.984 | 0.822 | 3235 | -0.827 | 1.4231 | 0.0277 | 412 | 0.0388 |
| PCE | 0.991 | 0.1773 | -3.780 | 0.877 | 2601 | -0.775 | 0.5329 | 0.0103 | 154 | 0.0147 |
| CIBENZ | 0.994 | 0.1457 | -2.601 | 0.911 | 2292 | -0.747 | 0.4370 | 0.0088 | 126 | 0.0122 |
| 1112TetCA | 0.996 | 0.1235 | -2.676 | 0.898 | 2281 | -0.725 | 0.3705 | 0.0073 | 107 | 0.0104 |
| EtBENZ | 0.993 | 0.1597 | -2.930 | 0.918 | 2357 | -0.752 | 0.4794 | 0.0097 | 138 | 0.0134 |
| mpXYL | 0.992 | 0.1678 | -3.036 | 0.909 | 2372 | -0.749 | 0.5037 | 0.0101 | 145 | 0.0140 |
| oXYL | 0.993 | 0.1555 | -2.862 | 0.911 | 2312 | -0.740 | 0.4667 | 0.0094 | 135 | 0.0131 |
| 1122TetCA | 0.996 | 0.1118 | -1.971 | 0.913 | 2167 | -0.691 | 0.3351 | 0.0067 | 97 | 0.0096 |
| 135TMB | 0.988 | 0.2024 | -4.435 | 0.897 | 2720 | -0.738 | 0.6093 | 0.0121 | 176 | 0.0170 |
| 124TMB | 0.989 | 0.1997 | -4.126 | 0.890 | 2631 | -0.731 | 0.6009 | 0.0118 | 173 | 0.0169 |
| 13DCB | 0.991 | 0.1832 | -3.422 | 0.888 | 2449 | -0.730 | 0.5503 | 0.0108 | 159 | 0.0155 |
| 14DCB | 0.991 | 0.1802 | -3.263 | 0.892 | 2408 | -0.724 | 0.5413 | 0.0107 | 156 | 0.0153 |
| 12DCB | 0.992 | 0.1697 | -2.970 | 0.894 | 2327 | -0.716 | 0.5092 | 0.0101 | 147 | 0.0144 |
| UNDEC | 0.694 | 0.374 | -1.406 | 0.426 | 1708 | -0.806 | 1.792 | 0.028 | 517 | 0.053 |
| NAPH | 0.992 | 0.166 | -3.374 | 0.915 | 2430 | -0.671 | 0.497 | 0.010 | 144 | 0.014 |
| TRIDEC | | | | | | | | | | |
| 2MeNAPH | 0.984 | 0.238 | -5.498 | 0.869 | 2990 | -0.689 | 0.72 | 0.014 | 208 | 0.021 |
| PENTADEC | | | | | | | | | | |
| Total mass | 0.993 | 0.1543 | -6.111 | 0.907 | 2419 | -0.732 | 0.4666 | 0.0093 | 134 | 0.0130 |

TABLE C
8260C MASS UNCERTAINTY

**AGI 8260C Method for Mass using SPG-0008
Samplers**

| | 99% Uncertainty Range +/- | 95% Uncertainty Range +/- |
|-----------|---------------------------------|---------------------------------|
| MTBE | 20% | 14% |
| t12DCE | 22% | 15% |
| 11DCA | 18% | 12% |
| c12DCE | 18% | 12% |
| CHCl3 | 16% | 11% |
| 111TCA | 18% | 12% |
| 12DCA | 20% | 13% |
| BENZ | 16% | 10% |
| CCl4 | 19% | 12% |
| TCE | 15% | 10% |
| 112TCA | 18% | 12% |
| TOL | 15% | 10% |
| OCT | 20% | 13% |
| PCE | 16% | 11% |
| CIBENZ | 18% | 12% |
| 1112TetCA | 19% | 13% |
| EtBENZ | 18% | 12% |
| mpXYL | 18% | 12% |
| oXYL | 18% | 12% |
| 1122TetCA | 23% | 15% |
| 135TMB | 21% | 14% |
| 124TMB | 20% | 14% |
| 13DCB | 19% | 13% |
| 14DCB | 19% | 13% |
| 12DCB | 20% | 14% |
| NAPH | 21% | 14% |
| 2MeNAPH | 25% | 17% |

TABLE D
4 CONSTANT WATER CONCENTRATION UNCERTAINTY
ERROR IN CONCENTRATION REPORTING (1)

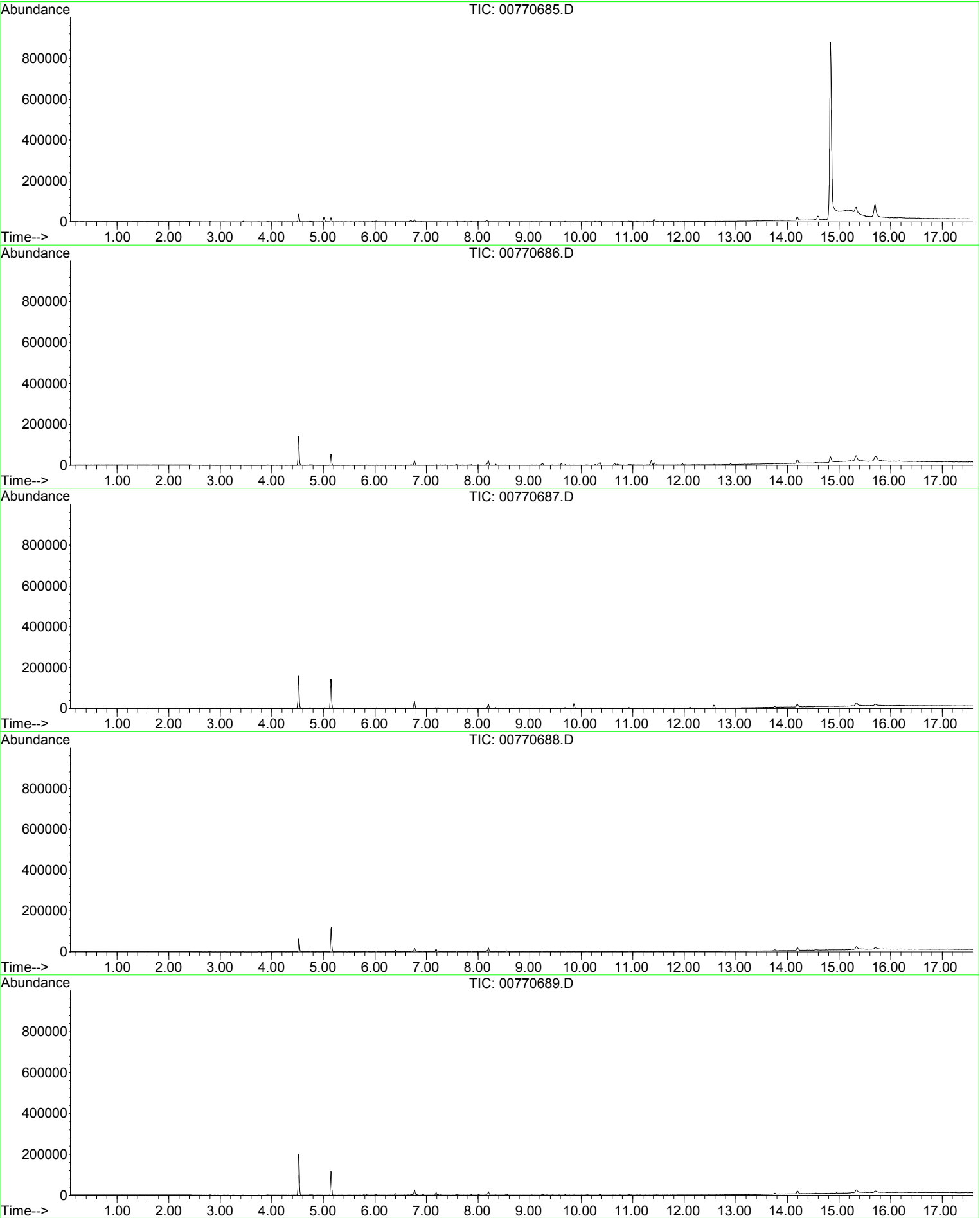
| | Average Error | Minimum Error | Maximum Error |
|-----------|--------------------------|--------------------------|--------------------------|
| MTBE | 6% | -12% | 12% |
| t12DCE | 11% | -26% | 21% |
| 11DCA | 8% | -19% | 13% |
| c12DCE | 9% | -19% | 15% |
| CHCl3 | 9% | -20% | 14% |
| 111TCA | 9% | -19% | 23% |
| 12DCA | 10% | -19% | 17% |
| BENZ | 8% | -18% | 13% |
| CCl4 | 10% | -23% | 22% |
| TCE | 10% | -21% | 14% |
| 112TCA | 11% | -21% | 21% |
| TOL | 7% | -17% | 14% |
| OCT | 20% | -41% | 42% |
| PCE | 10% | -24% | 15% |
| CIBENZ | 7% | -16% | 14% |
| 1112TetCA | 8% | -17% | 18% |
| EtBENZ | 6% | -19% | 14% |
| mpXYL | 7% | -22% | 13% |
| oXYL | 7% | -19% | 13% |
| 1122TetCA | 8% | -16% | 17% |
| 135TMB | 9% | -23% | 17% |
| 124TMB | 10% | -28% | 19% |
| 13DCB | 10% | -22% | 17% |
| 14DCB | 10% | -22% | 17% |
| 12DCB | 9% | -23% | 17% |
| NAPH | 10% | -24% | 21% |
| 2MeNAPH | 13% | -32% | 30% |

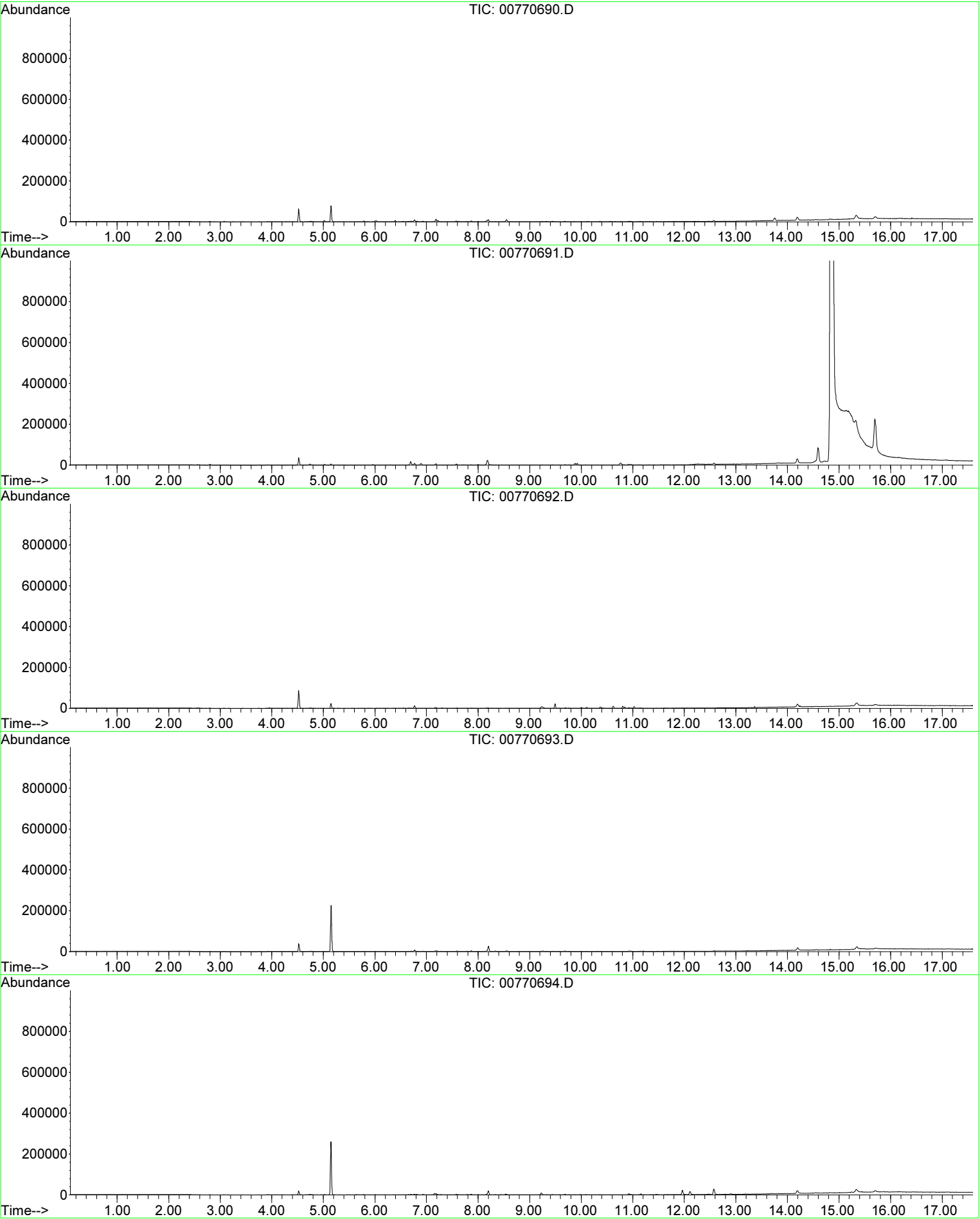
(1) For 1 hour exposure, includes error related to mass value from AGI analytical method 8260C

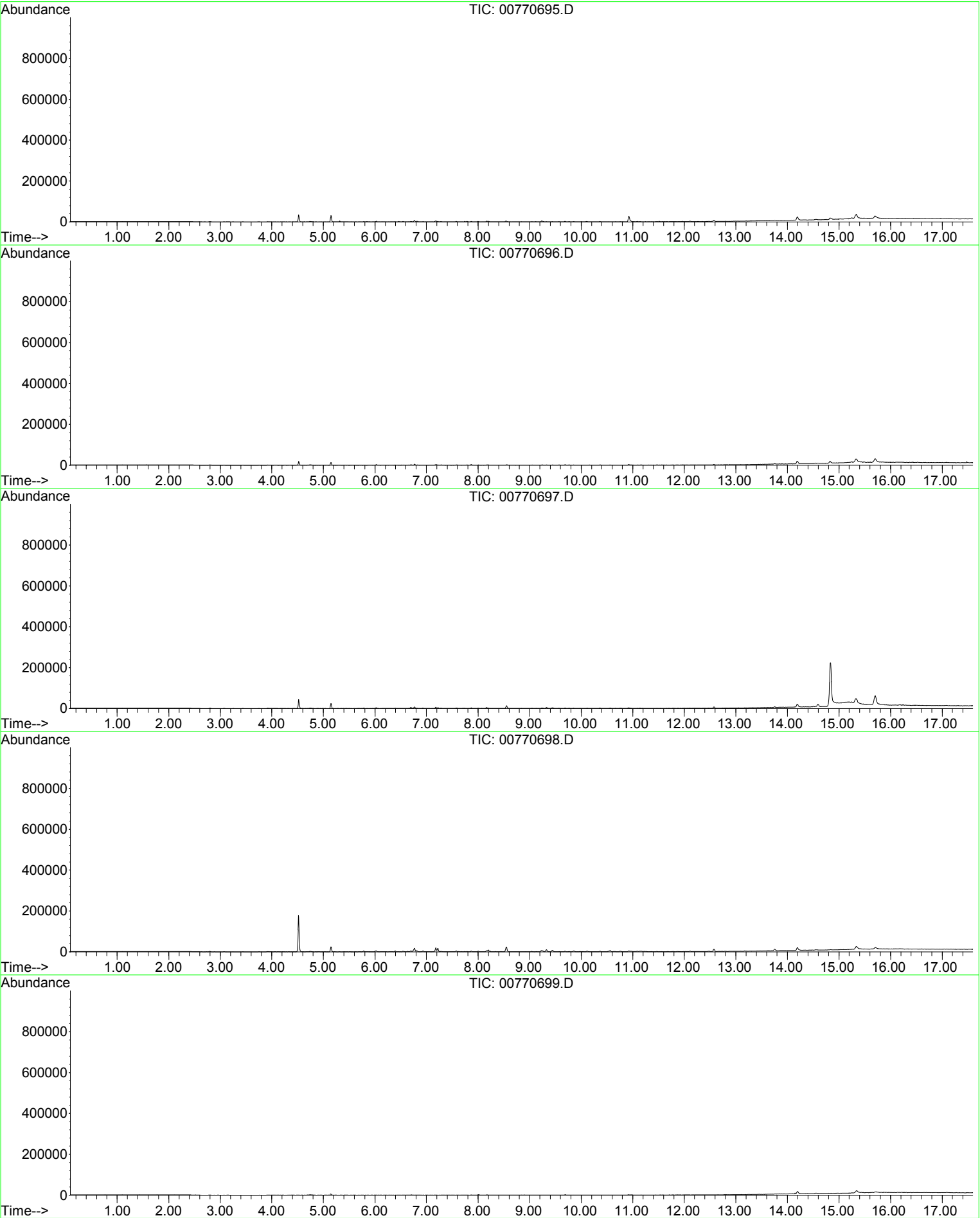
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www.agisurveys.net



Site: Comal & San Marcos Dec 2015
San Antonio, TX

Prepared for:

SWCA Environmental Consultants
6200 UTSA Boulevard
Suite 102
San Antonio, TX
UNITED STATES

Prepared on:
December 31, 2015

Project Summary and Objective

Amplified Geochemical Imaging, LLC. (AGI) provided the AGI Environmental Survey used at:

Comal & San Marcos Dec 2015

San Antonio, TX

The service provided by AGI included delivery of the required quantity of AGI Universal Samplers, analysis by the method described below for the requested organic compounds, reporting of the data, and contour mapping (as needed).

This report includes results for only the samples noted under the Laboratory Sample Report section. If contour maps are part of the project deliverable, the maps will be prepared and issued under a separate report cover, upon receipt of a usable sitemap (electronic) and compound choices for contouring.

Written/submitted by:

Kelly J Stringham

Project Manager

Reviewed/approved by:

Don D'Apolito

Project Manager

Analytical data approved by:

Ian McMullen

Chemist

Quality Assurance Statement

The AGI Laboratory, at Amplified Geochemical Imaging's facility in Newark, DE USA, operates under the guidelines of its ISO Standard 17025 DoD ELAP accreditation, and its Quality Assurance Manual, Operating Procedures, and Methods (SOP-QA-0462).

For this project, the analytical method, results, and observations reported do [] do not [✓] fall within the scope of AGI's ISO 17025 accreditation.

Screening/Concentration Method

The AGI Universal Samplers are analyzed at AGI's fixed laboratory using thermal desorption-gas chromatography/mass spectrometry (TD-GC/MS) instrumentation following modified U.S. EPA Method 8260 (SPG-WI-0292) which includes the following:

- **BFB Tuning Frequency:** A BFB tune is analyzed at the start of each analytical run and after every 30 samples.
- **Initial Calibration:** A minimum of a five point calibration curve is analyzed prior to the analysis of samples.
- **Initial Calibration Verification (ICV):** Following the calibration a second-source reference standard is analyzed to verify the accuracy of the calibration. Acceptance criteria for the ICV is +/- 30%.
- **Linearity of Target Compounds:** If the RSD of any target analyte is less than or equal to 25% then average response factor can be used for quantitation. If the RSD exceeds 25% for a target compound a regression equation can be used for quantitation.
- **Continuing Calibration Verification:** After every 10 samples, and at the end of each analytical batch, a mid-level second-source Reference Standard is analyzed. The acceptance criteria for all target analytes in the reference standards are +/- 50% of the true value.
- **Method Blank:** Analyzed prior to the analysis of field samples and every 30 samples.

Note: Analyte levels reported for the field-deployed AGI Universal Samplers that exceed trip and method blank levels, and/or the reporting limit, are more likely to have originated from on-site sources.

| | |
|----------------------------|---------------|
| Media Sampled: | WATER |
| Chemist - sample analysis: | Jasmine Smith |
| Chemist - data processor: | Jasmine Smith |
| Chemist - data review: | Ian McMullen |

Method deviations: None

Please note that data file names ending with R are rerun samples using the second pair of sorbers, in which the original results were not reported. Data file names ending in D are duplicate analysis results for the second set of sorbers from the same sampler, and are reported.

Additional Report Information

- Comments
- Laboratory Sample Report
- Chain of Custody
- Installation and Retrieval Log
- Analytical Results and Key
- Concentration Calculation Method Summary
- Total Ion Chromatograms

Project Specific Comments

None

Survey period ¹ Samplers were installed on December 1, 2015 and retrieved on December 15, 2015 for a sampling period of 14 days.

Tamper seal intact: Yes

Date received: 12/16/15 11:00 pm By: Clarence W Whigham

COC returned: Yes

Comments:

Shipment of samplers received by AGI included unused samplers provided to SWCA at an earlier date .

1 - Installation start to end of retrieval, as reported. See installation and retrieval log for individual deployment and retrieval dates and times (i.e., sampler exposure time).

General Comments

Analytical QA/QC

Laboratory instrumentation consists of gas chromatographs equipped with mass selective detectors, coupled with automated thermal desorption units. Sample preparation involves cutting the tip off the bottom of the AGI Universal Sampler, and transferring one or more "sorbents" to a thermal desorption tube for analysis. The insertion/retrieval cord prevents soil, water and other interferences from coming in contact with the adsorbent. No further sample preparation is required. Any replicate sorbents not consumed in the initial analysis will be discarded fifteen (15) days from the date of the laboratory report.

Data are archived and stored in a secure manner as per AGI's Quality Assurance program (SOP-QA-0462).

Total petroleum hydrocarbons (TPH), gasoline-range petroleum hydrocarbons (GRPH), and/or diesel range petroleum hydrocarbons (DRPH), when reported, are calculated using the area under the peaks observed in m/z 55 and 57 selected ion chromatograms. Quantitation of the mass values was performed using the response factor for a specific alkane (present in the calibration standards). TPH values include the entire chromatogram and provide estimates for aliphatic hydrocarbon ranges of C4 to C20. GRPH and DRPH include only the relevant regions of the chromatograms and provide estimates for C4 to C10 and C10 to C20 aliphatic hydrocarbons, respectively.

Trip blanks were provided to document potential exposures that were not part of the signal of interest (e.g., impact during sampler shipment, installation and/or retrieval, and storage). The trip blanks are identically manufactured and packaged AGI Universal Samplers to those samplers deployed in the field. The trip blanks remain unopened during all phases of the project. Levels reported on the trip blanks may indicate potential impact to the samplers other than the contaminant source of interest.

Unresolved peak envelopes (UPEs) are represented as a series of compound peaks clustered together around a central gas chromatograph elution time in the total ion chromatogram. UPEs may be indicative of complex fluid mixtures. UPEs observed early in the chromatograms are considered to indicate presence of more volatile fluids, while UPEs observed later in the chromatogram may indicate the presence of less volatile fluids. Multiple UPEs may indicate the presence of multiple complex fluids.

Total ion chromatograms (TICs) are included in the Attachments. The eight-digit serial number of each sampler is incorporated in the TIC identification (e.g., 12345678.D represents AGI Universal Sampler 12345678).

General Comments

Soil Gas Sampling

For soil gas sampling, the AGI Environmental Survey reports mass levels migrating through the open pore spaces of the soil and diffusing through the sampler membrane for sorption by the engineered, hydrophobic adsorbents, housed within the membrane tube. During the migration of the soil gas away from the source to the AGI Universal Sampler, the vapors are subject to a variety of attenuation factors. The soil gas masses reported on the samplers compare favorably with the concentrations reported in the soil or groundwater (e.g., where soil gas levels are reported at greater levels to other sampled locations on the site, the matrix data should reveal the same pattern, and vice versa). However, due to a variety of factors, a perfect comparison between matrix data and soil gas levels can rarely be achieved.

Soil gas concentrations ($\mu\text{g}/\text{m}^3$) are calculated following the method described in the Additional Report Information section.

Soil gas signals reported by this method cannot be correlated specifically to soil adsorbed, groundwater, and /or free-phase contamination. The soil gas signal reported from each AGI Universal Sampler can evolve from all of these sources. Differentiation between soil and groundwater contamination can only be achieved with prior knowledge of the site history (i.e., the site is known to have groundwater contamination only).

Air Sampling

For indoor, outdoor, and crawlspace air sampling, the AGI Environmental Survey reports mass levels present in the air and diffusing through the sampler membrane for sorption by the engineered adsorbents housed within the membrane tube.

Air concentrations ($\mu\text{g}/\text{m}^3$) are calculated following the method described in the Additional Report Information section.

Groundwater and Sediment Porewater Sampling

For groundwater and sediment porewater sampling, the AGI Environmental Survey reports the mass levels of compounds present in the water which, when coming in contact with the sampler membrane, partitions out of solution, and diffuses through the sampler membrane for sorption by the engineered adsorbents.

Water concentrations ($\mu\text{g}/\text{L}$) are calculated using the quantified mass, exposure period and the compound specific uptake rate. The rates were measured under controlled experimental conditions. The uptake rates are corrected for water pressure (depth of the AGI Universal Sampler below the water table), water temperature and the aquifer flow rate. For sediment porewater, the uptake rate is corrected for the reduced volume of water in the sediment, by multiplying the uptake rate by the pore water fraction.

LABORATORY SAMPLE REPORT

Project: ENV 01536

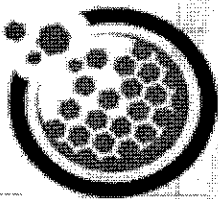
Site Name: Comal & San Marcos Dec 2015

Module Type: SPG0008

| Module ID | Sample Type | Field ID | |
|----------------------------|--------------------------|---------------------|-------------------|
| 00770871 | FIELD_SAMPLE | HCS410 | |
| 00770872 | FIELD_SAMPLE | HCS420 | |
| 00770873 | FIELD_SAMPLE | HCS430 | |
| 00770874 | FIELD_SAMPLE | HCS440 | |
| 00770875 | FIELD_SAMPLE | FDHCS440 | |
| 00770876 | LOST | HCS460 | |
| 00770877 | FIELD_SAMPLE | HSM410 | |
| 00770878 | FIELD_SAMPLE | HSM420 | |
| 00770879 | FIELD_SAMPLE | HSM430 | |
| 00770880 | FIELD_SAMPLE | FDHSM430 | |
| 00770881 | FIELD_SAMPLE | HSM440 | |
| 00770882 | FIELD_SAMPLE | HSM450 | |
| 00770883 | FIELD_SAMPLE | HSM460 | |
| 00770884 | FIELD_SAMPLE | HSM470 | |
| 00770885 | TRIP_BLANK | TB16 | |
| Total # "FIELD SAMPLES" | Total # "TRIP BLANKS" | Total # "UNUSED" | Total # "LOST" |
| 13 | 1 | 0 | 1* |
| Duplicate samples: | | | |

Duplicate samples: 0

***770876 was listed as "not retrieved" on the installation/retrieval log.**



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AGI Universal Passive Sampler Chain of Groundwater Sampling

Production Order #: **01536**

Customer Name: SWCA Environmental Consultants
Address: 6200 UTSA Boulevard
Suite 102

Site Name: Comal and San Marcos Rivers De
Site Address:

San Antonio, TX 78249
USA

Project Manager:

Serial # of Samplers Shipped
00770871 - 00770885

of Samplers for Installation 14.00
Total Samplers Shipped 15.00
Total Samplers Received 15
Total Samplers Installed 14

of Trip Blanks 1
Pieces
Pieces
Pieces

00770876 lost in field

00755239 - 00755241 > returned unused
00753380 - 00753383 from previous over stock

Serial # of Trip Blanks (Client Decides)

00770885

Prepared By:

Verified By:

Is Concurrent water sampling
planned this monitoring period?

YES

NO

Scheduled Sampling Date:

Installation Performed By:

Name: Jennifer Moreland Brittany Rios

Company: SWCA

Retrieval Performed By:

Name: Jennifer Moreland Guy Rubio

Company: SWCA

Installation Start Date / Time: 12/11/15 1013

Installation Complete Date / Time: 12/15/15 1412

Retrieval Start Date / Time: 12/15/15 1028

Retrieval Complete Date / Time: 12/15/15 1328

Total Samplers Retrieved:

13

Total Samplers Lost In Field:

1

Total Unused Samplers Returned:

7 unused samples returned from old 2014 supply or when we received 2 trip blanks

Relinquished By:

Company: AGI

Date/Time

Received By: Jennifer Moreland

Company: SWCA

Date/Time

11/19/15

12:30 PM

Relinquished By: Jennifer Moreland

Company: SWCA

Date/Time

12/15/15

1730

Received By: Clarence Whigham

Company: AGI

Date/Time

12/14/15

11:00



**AMPLIFIED
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IMAGING, LLC**

210 Executive Drive, Suite 1
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ph: 302-266-2428

AGI Project No.

ENV 01536

Site Name:

Comal and San Marcos Rivers De

Site Location:

New Braunfels and San Marcos, TX

AGI Soil Gas Sampling

Installation & Retrieval Log

Company Name:

SWCA Environmental Consultants

Location:

New Braunfels and San Marcos, TX

Samples collected by:

Jennifer Moreland, Brittany Rios, Guy Rubio

* Optional or as needed

| | | | | | WATER QUALITY M | | |
|--------------------|----------|---|--|---|--|--|---|
| SAMPLER SERIAL NO. | WELL ID | SAMPLE TYPE (Field Sample, Trip Blank, Field Blank, etc.) | INSTALLATION DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 13:00 | RETRIEVAL DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 15:00 | INSTALLATION DEPTH FROM TOP OF CASING (feet) | DEPTH TO WATER FROM TOP OF CASING (feet) | SCREENED INTERVAL FROM TOP OF CASING (feet) |
| 00770871 | HCS410 | FIELD_SAMPLE | 12/1/15 10:28 | 12/15/15 10:44 | 1.1 | | |
| 00770872 | HCS420 | FIELD_SAMPLE | 12/1/15 10:52 | 12/15/15 10:52 | 2.1 | | |
| 00770873 | HCS430 | FIELD_SAMPLE | 12/1/15 10:13 | 12/15/15 10:28 | 3.25 | | |
| 00770874 | HCS440 | FIELD_SAMPLE | 12/1/15 11:06 | 12/15/15 11:01 | 1.9 | | |
| 00770875 | FDHCS440 | FIELD_SAMPLE | 12/1/15 11:06 | 12/15/15 11:01 | 1.9 | | |
| 00770876 | HCS460 | FIELD_SAMPLE | 12/1/15 11:23 | 12/15/15 11:11 | 2.65 | | |
| 00770877 | HSM410 | FIELD_SAMPLE | 12/1/15 12:48 | 12/15/15 12:06 | 1.5 | | |
| 00770878 | HSM420 | FIELD_SAMPLE | 12/1/15 13:05 | 12/15/15 12:14 | 2.9 | | |
| 00770879 | HSM430 | FIELD_SAMPLE | 12/1/15 13:15 | 12/15/15 12:23 | 0.55 | | |
| 00770880 | FDHSM430 | FIELD_SAMPLE | 12/1/15 13:15 | 12/15/15 12:23 | 0.55 | | |
| 00770881 | HSM440 | FIELD_SAMPLE | 12/1/15 13:30 | 12/15/15 12:35 | 5 | | |
| 00770882 | HSM450 | FIELD_SAMPLE | 12/1/15 13:45 | 12/15/15 13:09 | 2.1 | | |
| 00770883 | HSM460 | FIELD_SAMPLE | 12/1/15 13:58 | 12/15/15 13:19 | 0.8 | | |
| 00770884 | HSM470 | FIELD_SAMPLE | 12/1/15 14:12 | 12/15/15 13:28 | 1.8 | | |
| 00770885 | TB16 | TRIP_BLANK | 12/1/15 10:13 | 12/15/15 13:28 | NA | | |
| | | | | | | | |
| | | | | | | | |



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ph: 302-266-2428

AGI Project No.

ENV 01536

Site Name:

Comal and San Marcos Rivers De

Site Location:

New Braunfels and San Marcos, TX

AGI Soil Gas Sampling

Installation & Retrieval Log

Company Name:

SWCA Environmental Consultants

Location:

New Braunfels and San Marcos, TX

Samples collected by:

Jennifer Moreland, Brittany Rios, Guy Rubio

* Optional or as needed

| | | | | | WATER QUALITY M | | |
|---|---------|---|--|---|--|--|---|
| SAMPLER SERIAL NO. | WELL ID | SAMPLE TYPE (Field Sample, Trip Blank, Field Blank, etc.) | INSTALLATION DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 13:00 | RETRIEVAL DATE & TIME MM/DD/YYYY HH:MM (24 Hour) ex. 12/30/2000 15:00 | INSTALLATION DEPTH FROM TOP OF CASING (feet) | DEPTH TO WATER FROM TOP OF CASING (feet) | SCREENED INTERVAL FROM TOP OF CASING (feet) |
| | | | | | | | |
| Old, unused samplers were also included in the shipment for return. There serial numbers were included on the chain of custody. | | | | | | | |



**AGI Soil Gas Sampling
Installation & Retrieval L**

* Optional or as needed

| MONITORING | | | OBSERVATIONS/COMMENTS* (e.g., observations, location description, missing, pulled from well, etc. - as needed) | YES / NO | |
|--------------------|--|--|--|--|--------|
| SAMPLER SERIAL NO. | WATER TEMPERATURE AT MODULE DEPTH (deg C) | FLOW RATE THROUGH SCREEN: HIGH OR LOW (HIGH = flow > than 10 m/day; LOW = flow < 10 m/day) | | EVIDENCE OF LIQUID PETROLEUM HYDROCARBONS? | ODOR ? |
| 00770871 | 22.6 | HIGH | | | |
| 00770872 | 22.6 | HIGH | All depths are water surface to top of sampler | | |
| 00770873 | 22.6 | HIGH | | | |
| 00770874 | 22.6 | HIGH | | | |
| 00770875 | 22.6 | HIGH | | | |
| 00770876 | 22.6 | HIGH | Not recovered | | |
| 00770877 | 21.4 | HIGH | | | |
| 00770878 | 21.4 | HIGH | | | |
| 00770879 | 21.4 | HIGH | | | |
| 00770880 | 21.4 | HIGH | | | |
| 00770881 | 21.4 | HIGH | | | |
| 00770882 | 21.4 | HIGH | | | |
| 00770883 | 21.4 | HIGH | | | |
| 00770884 | 21.4 | HIGH | | | |
| 00770885 | 21.4 | HIGH | | | |
| | | | | | |
| | | | | | |



**AGI Soil Gas Sampling
Installation & Retrieval L**

* Optional or as needed

| MONITORING | | | OBSERVATIONS/COMMENTS* (e.g., observations, location description, missing, pulled from well, etc. - as needed) | YES / NO | |
|---------------------------|--|--|--|--|--------|
| SAMPLER SERIAL NO. | WATER TEMPERATURE AT MODULE DEPTH (deg C) | FLOW RATE THROUGH SCREEN: HIGH OR LOW (HIGH = flow > than 10 m/day; LOW = flow < 10 m/day) | | EVIDENCE OF LIQUID PETROLEUM HYDROCARBONS? | ODOR ? |
| | | | | | |
| Old, unused samplers were | | | | | |



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PROJECT NUMBER: ENV 01536
SITE NAME: Comal & San Marcos Dec 2015
SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental
Consultants
San Antonio, TX 78249
USA

SAMPLER ID: 00770871 **FIELD_SAMPLE**

Dilution Factor: 1 Field ID: HCS410

Installation Date: 12/1/2015 10:28:00AM

Retrieval Date: 12/15/2015 10:44:00AM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Date Analyzed: 12/22/2015 5:47:00PM

Batch: ENV-151221-1

Reviewer:

Matrix: WATER

Product: SPG0008

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | 0.03 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.11 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770871 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS410

Installation Date: 12/1/2015 10:28:00AM

Retrieval Date: 12/15/2015 10:44:00AM

Date Analyzed: 12/22/2015 5:47:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770872 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HCS420

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 10:52:00AM

Retrieval Date: 12/15/2015 10:52:00AM

Date Analyzed: 12/22/2015 8:44:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | 0.11 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.38 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770872 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HCS420

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 10:52:00AM

Retrieval Date: 12/15/2015 10:52:00AM

Date Analyzed: 12/22/2015 8:44:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|-------------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | 0.11 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770873 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HCS430

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 10:13:00AM

Retrieval Date: 12/15/2015 10:28:00AM

Date Analyzed: 12/22/2015 5:17:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.58 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770873 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HCS430

Installation Date: 12/1/2015 10:13:00AM

Retrieval Date: 12/15/2015 10:28:00AM

Date Analyzed: 12/22/2015 5:17:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770874 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HCS440

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 11:06:00AM

Retrieval Date: 12/15/2015 11:01:00AM

Date Analyzed: 12/22/2015 2:49:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.38 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770874 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HCS440

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 11:06:00AM

Retrieval Date: 12/15/2015 11:01:00AM

Date Analyzed: 12/22/2015 2:49:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770875 FIELD_SAMPLE

Dilution Factor: 1

Field ID: FDHCS440

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 11:06:00AM

Retrieval Date: 12/15/2015 11:01:00AM

Date Analyzed: 12/22/2015 4:48:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.39 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770875 FIELD_SAMPLE

Dilution Factor: 1

Field ID: FDHCS440

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 11:06:00AM

Retrieval Date: 12/15/2015 11:01:00AM

Date Analyzed: 12/22/2015 4:48:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770877 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM410

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 12:48:00PM

Retrieval Date: 12/15/2015 12:06:00PM

Date Analyzed: 12/22/2015 3:19:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770877 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM410

Installation Date: 12/1/2015 12:48:00PM

Retrieval Date: 12/15/2015 12:06:00PM

Date Analyzed: 12/22/2015 3:19:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770878 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM420

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 1:05:00PM

Retrieval Date: 12/15/2015 12:14:00PM

Date Analyzed: 12/22/2015 2:19:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.10 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770878 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM420

Installation Date: 12/1/2015 1:05:00PM

Retrieval Date: 12/15/2015 12:14:00PM

Date Analyzed: 12/22/2015 2:19:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770879 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM430

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 1:15:00PM

Retrieval Date: 12/15/2015 12:23:00PM

Date Analyzed: 12/22/2015 6:46:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | 0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.81 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770879 FIELD_SAMPLE

Matrix: WATER

Product: SPG0008

Dilution Factor: 1

Field ID: HSM430

Installation Date: 12/1/2015 1:15:00PM

Retrieval Date: 12/15/2015 12:23:00PM

Date Analyzed: 12/22/2015 6:46:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770880 FIELD_SAMPLE

Dilution Factor: 1

Field ID: FDHSM430

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 1:15:00PM

Retrieval Date: 12/15/2015 12:23:00PM

Date Analyzed: 12/22/2015 8:15:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | 0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.76 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770880 FIELD_SAMPLE

Dilution Factor: 1

Field ID: FDHSM430

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 1:15:00PM

Retrieval Date: 12/15/2015 12:23:00PM

Date Analyzed: 12/22/2015 8:15:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770881 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM440

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 1:30:00PM

Retrieval Date: 12/15/2015 12:35:00PM

Date Analyzed: 12/22/2015 3:49:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.12 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770881 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM440

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 1:30:00PM

Retrieval Date: 12/15/2015 12:35:00PM

Date Analyzed: 12/22/2015 3:49:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |



PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770882 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM450

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 1:45:00PM

Retrieval Date: 12/15/2015 1:09:00PM

Date Analyzed: 12/22/2015 6:16:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.05 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770882 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM450

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 1:45:00PM

Retrieval Date: 12/15/2015 1:09:00PM

Date Analyzed: 12/22/2015 6:16:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



AMPLIFIED GEOCHEMICAL IMAGING, LLC

210 Executive Drive, Suite 1
Newark, DE 19702-3335 USA
ph: +1-302-266-2428
www.agisurveys.net

PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770883 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM460

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 1:58:00PM

Retrieval Date: 12/15/2015 1:19:00PM

Date Analyzed: 12/22/2015 4:18:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.10 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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www.agisurveys.net

PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770883 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM460

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 1:58:00PM

Retrieval Date: 12/15/2015 1:19:00PM

Date Analyzed: 12/22/2015 4:18:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770884 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM470

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 2:12:00PM

Retrieval Date: 12/15/2015 1:28:00PM

Date Analyzed: 12/22/2015 9:44:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|-------------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | 0.08 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770884 FIELD_SAMPLE

Dilution Factor: 1

Field ID: HSM470

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 2:12:00PM

Retrieval Date: 12/15/2015 1:28:00PM

Date Analyzed: 12/22/2015 9:44:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |



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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770885 TRIP_BLANK

Dilution Factor: 1

Field ID: TB16

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 10:13:00AM

Retrieval Date: 12/15/2015 1:28:00PM

Date Analyzed: 12/22/2015 9:14:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|---------------------------|-------------------|-------------|---------|
| Methyl tert-butyl ether | 1634-04-4 | <0.02 | 0.02 |
| trans-1,2-Dichloroethene | 156-60-5 | <0.02 | 0.02 |
| 1,1-Dichloroethane | 75-34-3 | <0.02 | 0.02 |
| cis-1,2-Dichloroethene | 156-59-2 | <0.02 | 0.02 |
| Chloroform | 67-66-3 | <0.02 | 0.02 |
| 1,1,1-Trichloroethane | 71-55-6 | <0.02 | 0.02 |
| 1,2-Dichloroethane | 107-06-2 | <0.02 | 0.02 |
| Benzene | 71-43-2 | <0.02 | 0.02 |
| Carbon Tetrachloride | 56-23-5 | <0.02 | 0.02 |
| Trichloroethene | 79-01-6 | <0.02 | 0.02 |
| 1,1,2-Trichloroethane | 79-00-5 | <0.02 | 0.02 |
| Toluene | 108-88-3 | <0.02 | 0.02 |
| Octane | 111-65-9 | <0.02 | 0.02 |
| Tetrachloroethene | 127-18-4 | <0.02 | 0.02 |
| Chlorobenzene | 108-90-7 | <0.02 | 0.02 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | <0.02 | 0.02 |
| Ethylbenzene | 100-41-4 | <0.02 | 0.02 |
| m,p-Xylene | 108-38-3/106-42-3 | <0.02 | 0.02 |
| o-Xylene | 95-47-6 | <0.02 | 0.02 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | <0.02 | 0.02 |
| 1,3,5-Trimethylbenzene | 108-67-8 | <0.02 | 0.02 |
| 1,2,4-Trimethylbenzene | 95-63-6 | <0.02 | 0.02 |
| 1,3-Dichlorobenzene | 541-73-1 | <0.02 | 0.02 |
| 1,4-Dichlorobenzene | 106-46-7 | <0.02 | 0.02 |
| 1,2-Dichlorobenzene | 95-50-1 | <0.02 | 0.02 |
| Undecane | 1120-21-4 | <0.05 | 0.05 |
| Naphthalene | 91-20-3 | <0.05 | 0.05 |
| Tridecane | 629-50-5 | <0.05 | 0.05 |
| 2-Methylnaphthalene | 91-57-6 | <0.05 | 0.05 |
| Acenaphthylene | 208-96-8 | <0.05 | 0.05 |
| Pentadecane | 629-62-9 | <0.05 | 0.05 |



AMPLIFIED GEOCHEMICAL IMAGING, LLC

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PROJECT NUMBER: ENV 01536

SITE NAME: Comal & San Marcos Dec 2015

SITE ADDRESS: San Antonio, TX

FOR: SWCA Environmental

Consultants

San Antonio, TX 78249

USA

SAMPLER ID: 00770885 TRIP_BLANK

Dilution Factor: 1

Field ID: TB16

Matrix: WATER

Product: SPG0008

Installation Date: 12/1/2015 10:13:00AM

Retrieval Date: 12/15/2015 1:28:00PM

Date Analyzed: 12/22/2015 9:14:00PM

Analyst: Jasmine Smith

Method: SPG-WI-0292

Batch: ENV-151221-1

Reviewer:

| Compound | CAS # | Result (ug) | RL (ug) |
|--------------------|------------|-------------|---------|
| Acenaphthene | 83-32-9 | <0.05 | 0.05 |
| Fluorene | 86-73-7 | <0.05 | 0.05 |
| TPH | | <0.50 | 0.50 |
| BTEX | | <0.02 | 0.02 |
| Phenanthrene | 85-01-8 | <0.05 | 0.05 |
| Anthracene | 120-12-7 | <0.05 | 0.05 |
| Fluoranthene | 206-44-0 | <0.05 | 0.05 |
| Pyrene | 129-00-0 | <0.05 | 0.05 |
| alpha-BHC | 319-84-6 | <0.05 | 0.05 |
| beta-BHC | 319-85-7 | <0.05 | 0.05 |
| gamma-BHC | 58-89-9 | <0.05 | 0.05 |
| delta-BHC | | <0.05 | 0.05 |
| Heptachlor | 76-44-8 | <0.05 | 0.05 |
| Aldrin | 309-00-2 | <0.05 | 0.05 |
| Heptachlor Epoxide | 1024-57-3 | <0.05 | 0.05 |
| Endosulfan I | 959-98-8 | <0.05 | 0.05 |
| 4,4-DDE | 72-55-9 | <0.05 | 0.05 |
| Dieldrin | 60-57-1 | <0.05 | 0.05 |
| Endrin | 72-20-8 | <0.05 | 0.05 |
| 4,4-DDD | 72-54-8 | <0.05 | 0.05 |
| Endosulfan II | 33213-65-9 | <0.05 | 0.05 |
| Endrin Aldehyde | 7421-93-4 | <0.05 | 0.05 |
| 4,4-DDT | 50-29-3 | <0.05 | 0.05 |
| Endosulfan Sulfate | 1031-07-8 | <0.05 | 0.05 |
| Endrin Ketone | 53494-70-5 | <0.05 | 0.05 |
| Methoxychlor | 72-43-5 | <0.05 | 0.05 |

AMPLIFIED GEOCHEMICAL IMAGING ANALYTICAL RESULTS
210 EXECUTIVE DRIVE, SUITE 1, NEWARK, DE
SWCA, SAN ANTONIO, TX
AGI STANDARD TARGET COMPOUNDS
ESTIMATED WATER CONCENTRATIONS
COMAL & SAN MARCOS RIVERS DECEMBER SAMPLING
ORDER #01536

| DATAFILE | FIELD | DATE/ TIME | DATE/ TIME | DATE/ TIME | DATE/ TIME | | estimated | | | | | | | | | | |
|----------------|----------|---------------|----------------|----------------|----------------|----|-----------|------------|--------------|-------------|--------------|-------------|--------------|-------------|------------|--|--|
| NAME | ID | INSTALLED | RETRIEVED | RECEIVED | ANALYZED | DF | TPH, ug/L | MTBE, ug/L | t12DCE, ug/L | 11DCA, ug/L | c12DCE, ug/L | CHCl3, ug/L | 111TCA, ug/L | 12DCA, ug/L | BENZ, ug/L | | |
| Average RL | | | | | | | 0.055 | 0.013 | 0.008 | 0.007 | 0.008 | 0.008 | 0.005 | 0.008 | 0.006 | | |
| 00770871 | HCS410 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.054 | <0.013 | <0.008 | <0.007 | <0.007 | 0.012 | <0.005 | <0.008 | <0.006 | | |
| 00770872 | HCS420 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.054 | <0.013 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | | |
| 00770873 | HCS430 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.054 | <0.013 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | | |
| 00770874 | HCS440 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.054 | <0.013 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | | |
| 00770875 | FDHCS440 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.054 | <0.013 | <0.008 | <0.007 | <0.007 | <0.007 | <0.005 | <0.008 | <0.006 | | |
| 00770877 | HSM410 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.056 | <0.014 | <0.008 | <0.008 | <0.008 | <0.008 | <0.005 | <0.008 | <0.007 | | |
| 00770878 | HSM420 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.056 | <0.014 | <0.008 | <0.008 | <0.008 | <0.008 | <0.005 | <0.008 | <0.007 | | |
| 00770879 | HSM430 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.056 | <0.014 | <0.008 | <0.008 | <0.008 | 0.009 | <0.005 | <0.008 | <0.007 | | |
| 00770880 | FDHSM430 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.056 | <0.014 | <0.008 | <0.008 | <0.008 | 0.009 | <0.005 | <0.008 | <0.007 | | |
| 00770881 | HSM440 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.056 | <0.014 | <0.008 | <0.008 | <0.008 | <0.008 | <0.005 | <0.008 | <0.007 | | |
| 00770882 | HSM450 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.056 | <0.014 | <0.008 | <0.008 | <0.008 | <0.008 | <0.005 | <0.008 | <0.007 | | |
| 00770883 | HSM460 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.056 | <0.014 | <0.008 | <0.008 | <0.008 | <0.008 | <0.005 | <0.008 | <0.007 | | |
| 00770884 | HSM470 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.056 | <0.014 | <0.008 | <0.008 | <0.008 | <0.008 | <0.005 | <0.008 | <0.007 | | |
| 00770885 | TB16 | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.055 | <0.013 | <0.008 | <0.007 | <0.008 | <0.008 | <0.005 | <0.008 | <0.006 | | |
| BLK ENV-277013 | | 12/1/2015:CST | 12/15/2015:CST | 12/16/2015:EST | 12/22/2015:EST | 1 | <0.055 | <0.013 | <0.008 | <0.007 | <0.008 | <0.008 | <0.005 | <0.008 | <0.006 | | |

AMPLIFIED GEOCHEMICAL IMAGING ANALYTICAL RESULTS
210 EXECUTIVE DRIVE, SUITE 1, NEWARK, DE
SWCA, SAN ANTONIO, TX
AGI STANDARD TARGET COMPOUNDS
ESTIMATED WATER CONCENTRATIONS
COMAL & SAN MARCOS RIVERS DECEMBER SAMPLING
ORDER #01536

| DATAFILE | FIELD | | | | | | | | | | | | | | | |
|----------------|----------|------------|-----------|--------------|-----------|-----------|-----------|--------------|-----------------|--------------|-------------|------------|-----------------|--------------|--------------|-------------|
| NAME | ID | CCl4, ug/L | TCE, ug/L | 112TCA, ug/L | TOL, ug/L | OCT, ug/L | PCE, ug/L | CIBENZ, ug/L | 1112TetCA, ug/L | ETBENZ, ug/L | mpXYL, ug/L | oXYL, ug/L | 1122TetCA, ug/L | 135TMB, ug/L | 124TMB, ug/L | 13DCB, ug/L |
| Average RL | | 0.005 | 0.006 | 0.010 | 0.006 | 0.005 | 0.006 | 0.006 | 0.007 | 0.005 | 0.005 | 0.006 | 0.011 | 0.005 | 0.005 | 0.006 |
| 00770871 | HCS410 | <0.005 | <0.006 | <0.009 | <0.006 | <0.005 | 0.023 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770872 | HCS420 | <0.005 | <0.006 | <0.010 | 0.027 | <0.005 | 0.071 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770873 | HCS430 | <0.005 | <0.006 | <0.009 | <0.006 | <0.005 | 0.10 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770874 | HCS440 | <0.005 | <0.006 | <0.010 | <0.006 | <0.005 | 0.070 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770875 | FDHCS440 | <0.005 | <0.006 | <0.010 | <0.006 | <0.005 | 0.072 | <0.006 | <0.007 | <0.005 | <0.005 | <0.005 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770877 | HSM410 | <0.005 | <0.006 | <0.010 | <0.006 | <0.005 | <0.006 | <0.007 | <0.007 | <0.006 | <0.006 | <0.006 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770878 | HSM420 | <0.005 | <0.006 | <0.010 | <0.006 | <0.005 | 0.022 | <0.007 | <0.007 | <0.006 | <0.006 | <0.006 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770879 | HSM430 | <0.005 | <0.006 | <0.010 | <0.006 | <0.005 | 0.15 | <0.007 | <0.007 | <0.006 | <0.006 | <0.006 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770880 | FDHSM430 | <0.005 | <0.006 | <0.010 | <0.006 | <0.005 | 0.14 | <0.007 | <0.007 | <0.006 | <0.006 | <0.006 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770881 | HSM440 | <0.005 | <0.006 | <0.010 | <0.006 | <0.005 | 0.028 | <0.007 | <0.007 | <0.006 | <0.006 | <0.006 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770882 | HSM450 | <0.005 | <0.006 | <0.010 | <0.006 | <0.005 | 0.013 | <0.007 | <0.007 | <0.006 | <0.006 | <0.006 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770883 | HSM460 | <0.005 | <0.006 | <0.010 | <0.006 | <0.005 | 0.023 | <0.007 | <0.007 | <0.006 | <0.006 | <0.006 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770884 | HSM470 | <0.005 | <0.006 | <0.010 | <0.006 | <0.005 | 0.019 | <0.007 | <0.007 | <0.006 | <0.006 | <0.006 | <0.011 | <0.005 | <0.005 | <0.006 |
| 00770885 | TB16 | <0.005 | <0.006 | <0.010 | <0.006 | <0.005 | <0.006 | <0.006 | <0.007 | <0.005 | <0.005 | <0.006 | <0.011 | <0.005 | <0.005 | <0.006 |
| BLK ENV-277013 | | <0.005 | <0.006 | <0.010 | <0.006 | <0.005 | <0.006 | <0.006 | <0.007 | <0.005 | <0.005 | <0.006 | <0.011 | <0.005 | <0.005 | <0.006 |

AMPLIFIED GEOCHEMICAL IMAGING ANALYTICAL RESULTS
 210 EXECUTIVE DRIVE, SUITE 1, NEWARK, DE
 SWCA, SAN ANTONIO, TX
 AGI STANDARD TARGET COMPOUNDS
 ESTIMATED WATER CONCENTRATIONS
 COMAL & SAN MARCOS RIVERS DECEMBER SAMPLING
 ORDER #01536

| DATAFILE | FIELD | | | | | estimated | | estimated | estimated | estimated | estimated |
|----------------|----------|-------------|-------------|-------------|------------|--------------|---------------|----------------------|----------------|--------------------|----------------|
| NAME | ID | 14DCB, ug/L | 12DCB, ug/L | UNDEC, ug/L | NAPH, ug/L | TRIDEC, ug/L | 2MeNAPH, ug/L | Acenaphthylene, ug/L | PENTADEC, ug/L | Acenaphthene, ug/L | Fluorene, ug/L |
| Average RL | | 0.006 | 0.006 | 0.021 | 0.017 | 0.021 | 0.014 | 0.015 | 0.021 | 0.015 | 0.015 |
| 00770871 | HCS410 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00770872 | HCS420 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00770873 | HCS430 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00770874 | HCS440 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00770875 | FDHCS440 | <0.006 | <0.006 | <0.020 | <0.016 | <0.020 | <0.013 | <0.014 | <0.020 | <0.014 | <0.014 |
| 00770877 | HSM410 | <0.006 | <0.007 | <0.021 | <0.017 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770878 | HSM420 | <0.006 | <0.007 | <0.021 | <0.017 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770879 | HSM430 | <0.006 | <0.007 | <0.021 | <0.017 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770880 | FDHSM430 | <0.006 | <0.007 | <0.021 | <0.017 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770881 | HSM440 | <0.006 | <0.007 | <0.021 | <0.017 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770882 | HSM450 | <0.006 | <0.007 | <0.021 | <0.017 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770883 | HSM460 | <0.006 | <0.007 | <0.021 | <0.017 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770884 | HSM470 | <0.006 | <0.007 | <0.021 | <0.017 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| 00770885 | TB16 | <0.006 | <0.006 | <0.021 | <0.017 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |
| BLK_ENV-277013 | | <0.006 | <0.006 | <0.021 | <0.017 | <0.021 | <0.014 | <0.015 | <0.021 | <0.015 | <0.015 |

KEY TO DATA TABLE

UNITS

| | |
|-------------------|--|
| µg | micrograms, relative mass value |
| µg/m ³ | micrograms per cubic meter; estimated soil gas concentration |
| µg/L | micrograms per Liter; calculated water concentration |

DATA QUALIFIERS

| | |
|---|--|
| > | greater than; value exceeds calibration range, estimated value |
| < | less than; compound value is below the LOD and RL |
| J | mass value below LOQ or RL, but above LOD, estimated mass value |
| E | mass value exceeds upper calibration level, estimated mass value |
| Q | one or more quality control parameters failed for the compound |

ABBREVIATIONS

| | |
|--------|--|
| AVG RL | average reporting limit; calculated based on individual field sample RLs |
| LOD | limit of detection |
| LOQ | limit of quantification |
| MDL | method detection limit |
| RL | reporting limit |

| | | | |
|-------------|---|----------|--|
| 1112TetCA | 1,1,1,2-tetrachloroethane | CIBENZ | chlorobenzene |
| 111TCA | 1,1,1-trichloroethane | ct12DCE | cis- & trans-1,2-dichloroethene |
| 1122TetCA | 1,1,2,2-tetrachloroethane | EtBENZ | ethylbenzene |
| 112TCA | 1,1,2-trichloroethane | mpXYL | m-, p-xylene |
| 11DCA | 1,1-dichloroethane | MTBE | methyl t-butyl ether |
| 11DCE | 1,1-dichloroethene | NAPH | naphthalene |
| 124TMB | 1,2,4-trimethylbenzene | OCT | octane |
| 12DCA | 1,2-dichloroethane | oXYL | o-xylene |
| 12DCB | 1,2-dichlorobenzene | PCE | tetrachloroethene |
| 135TMB | 1,3,5-trimethylbenzene | PENTADEC | pentadecane |
| 13DCB | 1,3-dichlorobenzene | PHEN | phenanthrene |
| 14DCB | 1,4-dichlorobenzene | t12DCE | trans-1,2-dichloroethene |
| 2MeNAPH | 2-methyl naphthalene | TCE | trichloroethene |
| BENZ | benzene | TMBs | combined masses of 1,3,5-trimethylbenzene and 1,2,4-trimethylbenzene |
| BTEX | combined masses of benzene, toluene, ethylbenzene, and total xylenes (Gasoline Range Aromatics) | TOL | toluene |
| C11,C13&C15 | combined masses of undecane, tridecane, and pentadecane (C11+C13+C15) (Diesel Range Alkanes) | TPH | total petroleum hydrocarbons |
| c12DCE | cis-1,2-dichloroethene | TRIDEC | tridecane |
| CCl4 | carbon tetrachloride | UNDEC | undecane |
| CHC13 | chloroform | VC | vinyl chloride |

SUMMARY OF SAMPLING RATE CALIBRATION FOR AGI UNIVERSAL SAMPLER IN AQUEOUS PHASES

INTRODUCTION:

The Amplified Geochemical Imaging, LLC (AGI) passive vapor sampler is designed to be used for soil gas, water, sediment pore water, and air sampling. This document describes the process used to calibrate the sampler's compound specific sampling or uptake rates in aqueous phases.

Sampling rates are measured following AGI's "Standard Practice for Determining the Sampling Rate of Passive Diffusion Samplers in Various Environmental Media": SPG-SOP-0493. Rates are used to calculate dissolved phase concentrations of volatile and semi-volatile contaminants in water. The calibration process is summarized in three parts: Part 1: shallow water, Part 2: deep water, and Part 3: sediment.

PURPOSE:

The purpose of this document is to:

1. Summarize the test protocol,
2. Summarize the methodology for analysis of data,
3. Present general results for generating concentration calibration of the AGI Universal Sampler

Principle of Operation of the AGI Sampler

The AGI Universal Sampler is designed with solid adsorbents enclosed inside a tubular microporous PTFE membrane. When placed in water, the pores and hydrophobic nature of the PTFE keep liquid water from entering the membrane until a water head of about 34 feet is reached. The membrane will not keep water vapor from entering but the adsorbents are very hydrophobic and through testing validated to be unaffected by this moisture vapor. In shallow water, <34', volatile and semi-volatile compounds will partition from the dissolved water into the air phase in the PTFE membrane according to Henry's Law. This partitioning is instantaneous and within seconds-minutes, the compound is adsorbed by the adsorbent inside the sealed tube. Because the diffusivity in air is about 10,000 times higher than the diffusivity in water, the sampling rate is controlled by the water contact area of the membrane that allows the Henry's Law effect to occur. This contact area is set by the membrane diameter and length of the sealed tube, which is fixed in AGI's manufacturing process.

Henry's law as well as diffusivity, which are fundamentally incorporated into the sampling rate, are affected by temperature, T , and follow an Arrhenius equation $H_T = H_r \times \exp\left(\frac{-E_a/R}{1/T_r - 1/T}\right)$. Because a 5°C temperature change can make a 15% change in sampling rate, the temperature of the sampled water should be known to get the most precise concentration.

The membrane pore size is also small enough that colloidal particles and microbes cannot pass through the membrane. This keeps the adsorbent from getting contaminated and eliminates any need to add preservative or chilling during storage or transportation.

When the water pressure exceeds the water entry pressure of the membrane, about 34 feet of water, the water becomes in contact with the solid adsorbent. Under this condition the compounds in the water will partition from the water into the solid. The partitioning coefficient, K_{AW} , can be approximated by the octanol-water coefficient, K_{OW} , but has been measured more precisely in the lab for AGI's specific solid adsorbent. The sampling rate is the product of the sampling rate at <34' of water and the K_{AW} .

In sediment, the sampler measures pore-water concentration, which is generally agreed to be the preferred measurement as it is more indicative of bioavailability. In sediment the volumetric

availability of water to the sampler is reduced by the volume fraction solids in the sediment, which typically varies from zero to 35%, but can be as high as 73% in well packed and broad particle size distribution sediments. As a result, sampling rates in sediment are multiplied by the fraction pore water in the sediment to determine concentration.

PART 1: Calibration in shallow water

Part 1 summarizes the work in shallow water generating calibration data, evaluating the physical and chemical factors affecting the sampling rate, and measurement of the actual sampling rates or regression calibration equations needed to determine concentrations.

Sample Generation in water

In this calibration work, solutions of analytes at known concentrations were formulated in clean 4 liter smoked glass jugs by injecting microliter measured amounts of environmental standards using a calibrated syringe into pure or deionized water and stirring for a minimum of 2 hours but generally overnight. Headspace in the jugs was minimized and generally less than 1% by volume during the tests. Jugs were temperature controlled by placing them in a water filled cooler, chilled via a copper tubing loop in the cooler. Temperature was measured with a certified digital temperature gauge and an average value used for each temperature experiment.

AGI samplers were weighted so they won't float and placed in the jugs at time zero. They were removed at various intervals to generate samples along with duplicates that showed mass increasing with exposure time. The sampler exposure time was selected to span minutes to hours and was generally reduced for high concentration tests to maintain uptake with time in roughly the linear dynamic range. Samplers were removed and dried with a paper towel and returned to their original container for analysis. They were analyzed by AGI's 8260C (SPG-WI-318 or SPG-WI-10028) method in duplicate, which is based on EPA SW846 Method 8260C.

Water samples were also taken and measured at an outside accredited lab using EPA SW846 Method 8260B. The concentrations agreed well with the calculated concentrations based on the standard certification, jug volume, and syringe injection. The variability of the outside lab 8260B values were found to be high, so for the sampling rate calculations we used the concentrations based on syringe dosing.

Calibrations were run at five concentrations, nominally at 6, 24, 118, 590, 1420 ug/L and five temperatures nominally at 5, 10, 15, 20, and 25 degrees centigrade. Samples were taken at 4 different exposure times. Samples were run in duplicate. A total of 176 data points were generated using 28 compounds from AGI's standard compounds list. Tridecane and pentadecane were not evaluated due to their very low solubility in water. In addition, another 23 compounds were tested using an 8260 liquid standard at nominal concentrations of 0.5, 1.0, 5.0, 15, 95, and 470 ug/L at a temperature typical of groundwater, 15°C. This is a living calibration and as additional data are generated, they may be qualified and added to this data set to improve the precision of the sampling rate calibration and broaden the compound list.

Key Variable Effects

As expected from theory, at short to moderate exposure times, mass will increase roughly linearly proportional to exposure time, as well as proportional to concentration, and exponentially with temperature following Arrhenius law. Temperature affects the Henry's law as well as diffusivity in water. Sampling rate is generally independent of concentration and time at mass values significantly below saturation. In the following sections we have characterized the sampling rate for each compound as affected by temperature and also developed calibrations using regression which account for the minor impact of time, and mass.

Concentration using Simple Sampling Rate Determination

A simple way to determine concentration is to measure mass on the AGI sampler, divide by exposure time, and divide by sampling rate, SR.

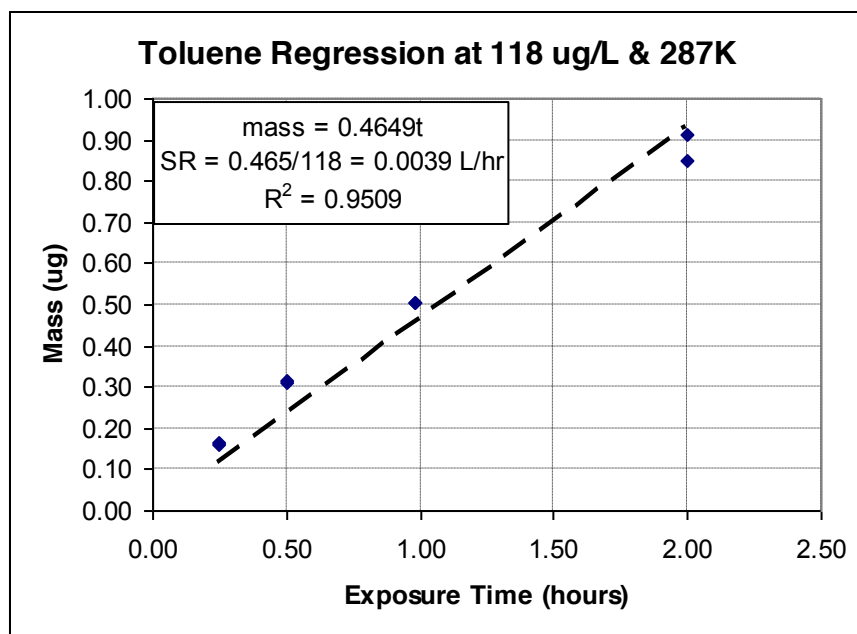
$$\text{Conc [ug/L]} = \text{mass/time/SR} \quad (1)$$

The sampling rate can be determined via measurements of mass versus time at a known concentration and temperature according to the following modification of equation (1).

$$\text{SR} = \text{mass/time/concentration} \quad (2)$$

Sampling rates in L/hr were determined by measuring the trend or regression mass uptake versus time and dividing by the concentration. A measurement like this will use 8 data points (4 times x 2 samples). Such a sampling rate can be measured at any concentration and temperature.

The chart to the right shows a plot of mass versus time for water at 118 ug/L and 287K (actual data from a single run). Slope of 0.465 ug/hr divided by the concentration of 118 ug/L yields a sampling rate, SR, of 0.0039 L/hr.



SR's typically range from about 0.004 to 0.007 L/hr at 15°C. Table A shows SR's measured for our standard compound list at 5 temperatures.

Rigorous Concentration using Regression

A preferred method for determining concentration that will yield improved accuracy over a wide range of concentrations, exposure times, and temperatures is to use all data in a regression analysis, which allows adjustments for the minor non-linear influences of mass and time as well as the effects of temperature. This step is done by regressing equation (1) or a universal version of equation (1):

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRo} * \exp(-\text{Ea}/\text{R}/\text{T})] \quad (3)$$

The subtle non-linear effects of mass and time will be evident in the deviation of coefficients b and d from 1.0. This regression generates four constants b, d, SRo, and $-\text{Ea}/\text{R}$ by regressing $\ln(\text{conc})$ versus $\ln(\text{mass})$, $\ln(\text{time})$, $1/\text{temp}$. These four constants can be used to determine concentration via the equation:

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRo} * \exp(-\text{Ea}/\text{R}(1/\text{T}))] \quad (4)$$

Where conc is in ug/L, mass is in ug, time in hours, T in degrees Kelvin.

Equation (4) can be also expressed at a reference temperature, Tr , such as 15°C by

$$\text{Conc} = (\text{mass})^b / (\text{time})^{-d} / [-\text{SRr} * \exp(-\text{Ea}/\text{R}(1/\text{Tr} - 1/\text{T}))] \quad (5)$$

This step allows sampling rates, SRr, at any reference temperature, Tr , and for any analyte to easily be compared. The values of SRr at 293.14K can be found in Table A.

When sampling times are between 0 and 4 hours, using the 4 constant equation (5) is preferred. For concentrations from about 5 to 1500 ug/L one hour exposure times generally give the lowest error, typically with average error of 6-20% and with total error range of 12%-32%. For low concentrations where sampling times are greater than 4 hours, it is preferred to use equation (1) to avoid unrealistic effects from the coefficient d or to set d to 1.0. In such a case SR in equation (1) can be substituted with $[\text{SRr} * \exp(-\text{Ea}/\text{R}(1/\text{Tr} - 1/\text{T}))]$ to use an SR representative of the well temperature, T.

The chart to the right is a plot of the calculated concentration from the 4 constant regression compared to the dosed concentration. Agreement is excellent for the 176 data points.

However, there does appear to be a slight high bias of 8.6% over the full range of this data, although it is well within acceptable limits of variability.

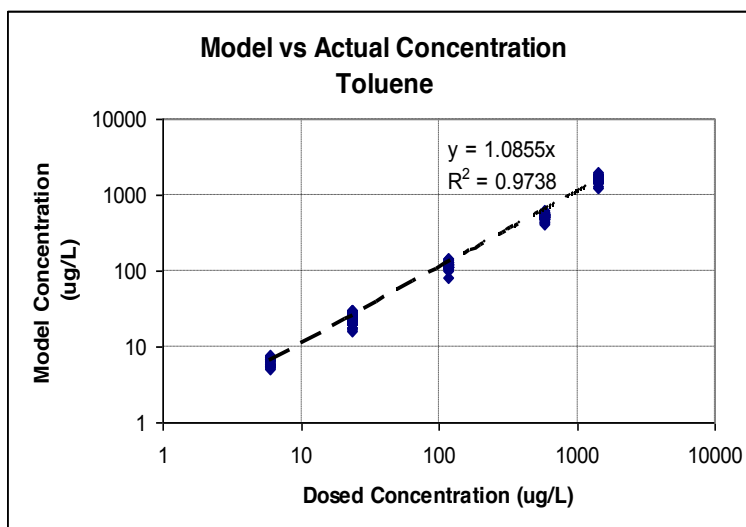


Table B shows the tabulated summary of the 4 constants regression with Rsq values and error estimates for the 4 constants for each analyte. Most regression Rsq values are 0.99 or greater for each analyte. In general, $-E_a/R$ is about 2400 \pm 400, b is about 0.9, d is about -0.75, and SR(15°C) ranges from .004 L/hr to 0.007 L/hr increasing with MW of the compound.

Error Estimates

The error in the water concentration values will depend on both the error in mass from the analytical method as well as the error in the concentration calibration. Table C shows the error in the mass values from the 8260C low sensitivity method.

The standard error of the regression and standard errors of the constants can be found in table B. For each compound we have measured the error between the derived concentration and the actual concentration. The error tends to be lowest at our recommended exposure time of one hour as shown by the example for Toluene to the right.

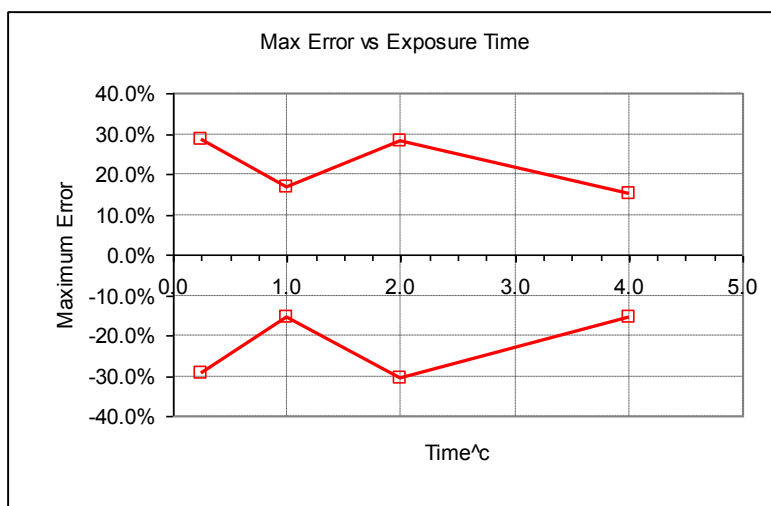
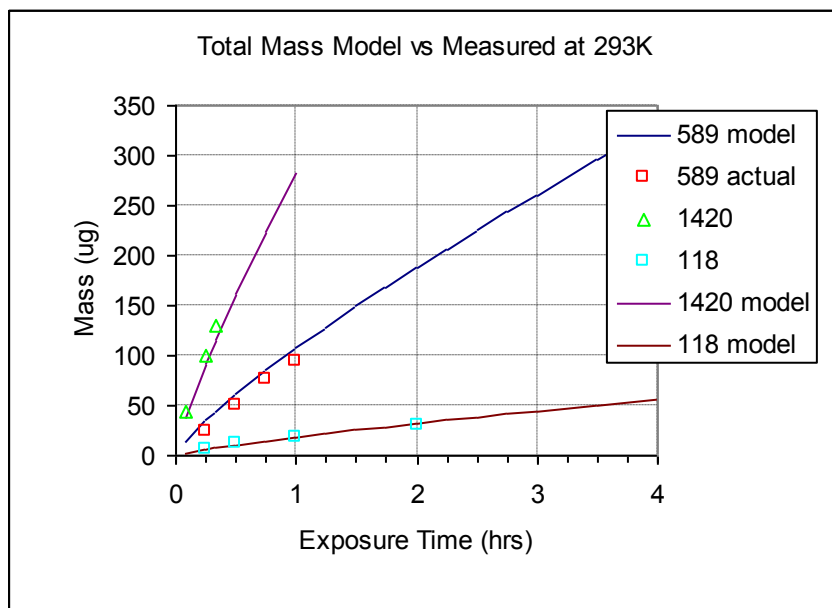


Table D shows the total average error in water concentration by compound as well as the low and high error. The average ranges from about 6% to 20%, which is similar to the analytical method errors. The low and high errors range from 12% to 32% and include contribution from measurement errors in both time and temperature.

Sorbent Saturation

As mass increases on a solid sorbent and approaches saturation, reverse diffusion can occur causing the sampling rate to drop. Eventually the mass level will reach a maximum steady state value at any concentration. A rate of mass uptake with time that deviates significantly from linear, indicates that sorbent saturation could be an issue. When using equation (1), staying in the linear range to avoid the effects of adsorbent saturation is important. We recommend keeping the total mass on the sampler below 50 ug or flagging when this is exceeded.

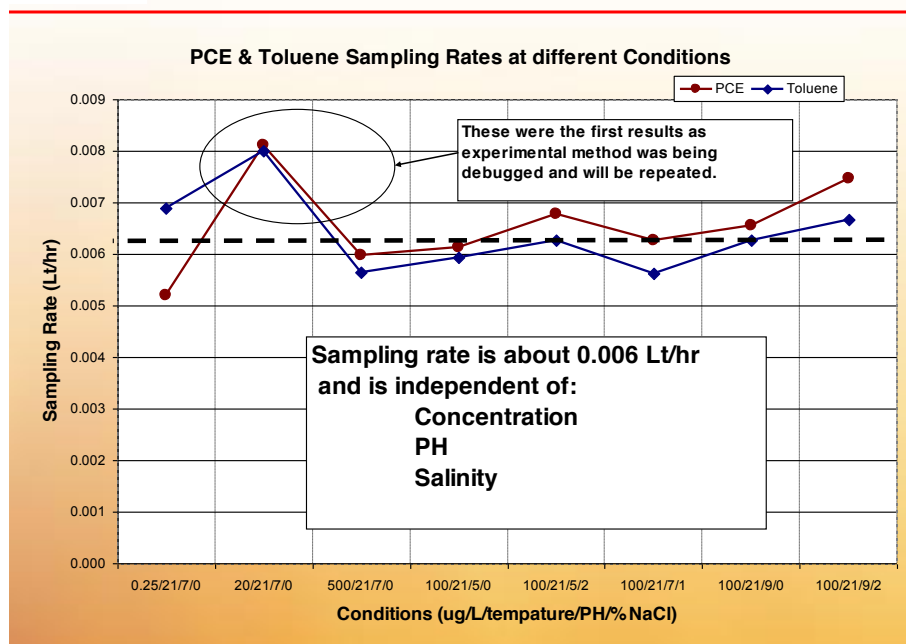
The 4 constant regression accounts for some of the non linearity allowing good accuracy at higher mass levels. From the experimental data we have found this safe range can be extended to 100 ug or higher as shown in the chart below. This chart compares total mass of all compounds (excluding heavy alkanes, which have solubility issues) versus time in comparison to that predicted from the 4-constant concentration equation.



Effect of PH and Salinity

Because neither PH nor salinity is known to have a significant impact on Henry's law or diffusivity in water, we did not expect them to have a significant impact on sampling rate. To confirm this, experiments were run varying PH from 5 to 9 and NaCl content from 0 to 2%. The chart below shows no significant impact for combinations of PH and NaCl content over this range on the sampling rate of toluene in water at 21°C.

Checked for Effects of PH & Salinity

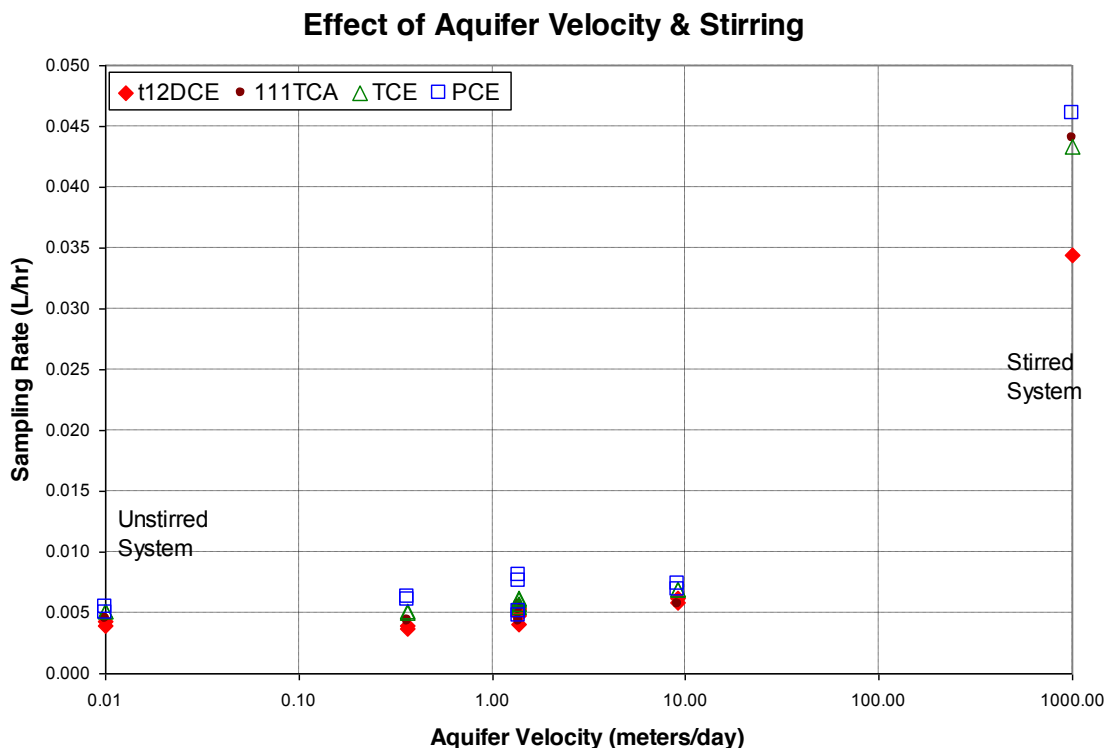


Impact of Aquifer Velocity

The velocity in most aquifers is quite slow, typically a meter/day or less. Occasionally water flow could be much higher such as encountered in karst aquifers, streams or rivers. Mass transfer coefficients are higher in high flow conditions, which will lead to higher sampling rates. We validated that a highly stirred system had sampling rates about 10 times higher than those that were non-stirred. We decided to evaluate the effect of aquifer velocity.

A test apparatus was built comprising a 3" PVC pipe tee filled with clean sand in each of the horizontal straight legs and screened to leave the center open. A test solution was run through this system using a variable flow pump and AGI samplers were placed into the simulated well through the vertical leg of the tee. Tests were run to examine the effect of velocity by varying the pumping rate and hence water velocity.

The chart below shows no significant effect of aquifer velocity up to a speed of about 10 meters/day. At velocities significantly above this, similar to a stirred system, sampling rates are about 10 times higher.

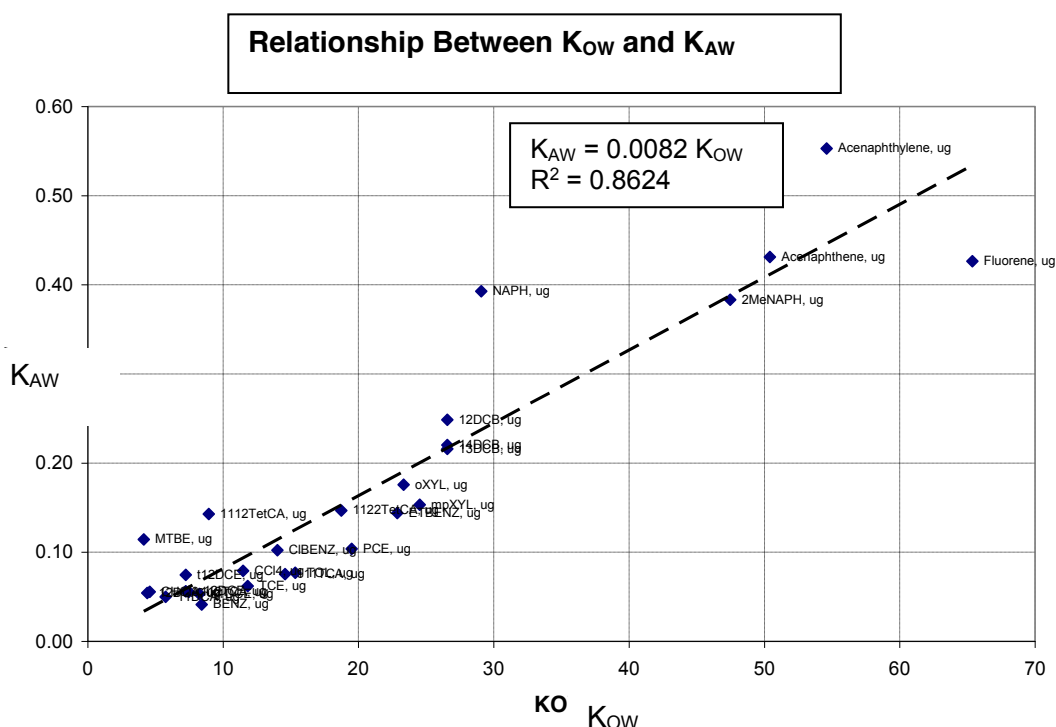


Part 2: Calibration in Deep (>34') water

Part 2 describes the effect of deep water on the AGI sampler and summarizes the effects on sampling rate and concentration measurement.

When the water pressure exceeds the water entry pressure of the membrane, about 34 feet of water, the water becomes in direct contact with the solid adsorbent. Under this condition the compounds in the water will partition from the water into the solid. The partitioning coefficient is closely related to the octanol-water coefficient, K_{OW} , but has been measured more precisely in the lab for AGI's specific solid adsorbent, K_{AW} . The sampling rate for deep water is the product of the sampling rate at <34' of water and the K_{AW} .

Measurement of the K_{AW} was done in a one liter stainless steel vessel pressurized with nitrogen to simulate water heads above 34' of water. Pressures of up to 465 psig or 200' of water head were used. The sampling rate change was the same at all pressures above 34' of water. The K_{AW} was determined as the ratio between the mass or sampling rate above 34' of head to the rate at <34' of head and is shown in the chart below.



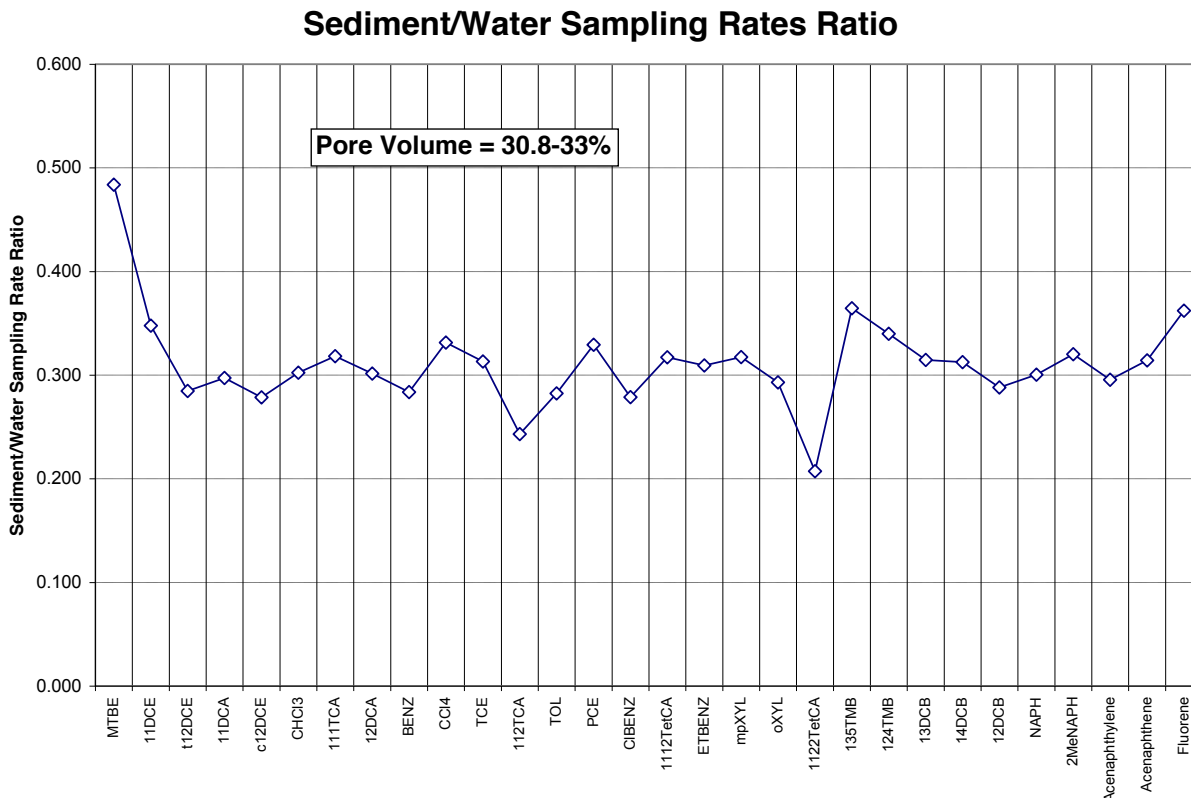
Part 3: Calibration in Sediment

Part 3 describes the effect of sediment solids or sediment pore volume on the sampling rate and concentration measurement.

In sediment, the sampler measures pore water concentration, which is generally agreed to be the preferred measurement as it is more indicative of bioavailability. In sediment the volumetric availability of water to the sampler is reduced by the volume fraction solids in the sediment. As a result sampling rates in sediment are multiplied by the fraction pore water to determine

concentration. Pore water fraction can range from 1.0 for water without sediment to as low as 0.25. Typically most sediments have pore fractions of 0.9 to 0.65.

A sampling rate study was done with water and with water added into a well-packed sorted sand. Pore water fraction in this test was measured between 30.8% and 33% by volume. Below is a plot of the ratio of sampling rates measured in the sediment to open water. The average ratio is equal to the pore water fraction confirming that sampling rate in sediment is on average equal to the product of pore water fraction times the sampling rate in water.



Summary

The AGI Sampler can be used to determine the concentration of volatile and semi-volatile compounds in a water phase. This requires knowing the exposure time and water temperature. It also requires knowing if the sample is above or below 34' of water head and if the water has a velocity above 10 meters/day. Regressions of large amounts of data were used to generate a four constant equation to generate concentration values in water. Potential error in the concentration values is excellent typically less than 25%.

TABLE A
WATER SAMPLING RATES STANDARD LIST

| | SRr 293.14 | SR @ 277.54 | SR @ 282.44 | SR @ 287.84 | SR @ 293.24 | SR @ 298.94 |
|-------------------|----------------------|-----------------------|-----------------------|-----------------------|-----------------------|-----------------------|
| MTBE | 0.0025 | 0.0014 | 0.0016 | 0.0018 | 0.0022 | 0.0029 |
| t12DCE | 0.0043 | 0.0028 | 0.0028 | 0.0027 | 0.0037 | 0.0048 |
| 11DCA | 0.0047 | 0.0031 | 0.0033 | 0.0033 | 0.0039 | 0.0052 |
| c12DCE | 0.0046 | 0.0030 | 0.0031 | 0.0031 | 0.0038 | 0.0051 |
| CHCl3 | 0.0046 | 0.0030 | 0.0031 | 0.0031 | 0.0038 | 0.0051 |
| 111TCA | 0.0066 | 0.0043 | 0.0047 | 0.0047 | 0.0056 | 0.0076 |
| 12DCA | 0.0045 | 0.0029 | 0.0029 | 0.0030 | 0.0036 | 0.0050 |
| BENZ | 0.0050 | 0.0031 | 0.0034 | 0.0035 | 0.0042 | 0.0056 |
| CCl4 | 0.0068 | 0.0044 | 0.0048 | 0.0047 | 0.0058 | 0.0080 |
| TCE | 0.0052 | 0.0030 | 0.0034 | 0.0034 | 0.0043 | 0.0058 |
| 112TCA | 0.0043 | 0.0027 | 0.0027 | 0.0028 | 0.0034 | 0.0048 |
| TOL | 0.0056 | 0.0034 | 0.0039 | 0.0039 | 0.0047 | 0.0062 |
| OCT | 0.0064 | 0.0046 | 0.0050 | 0.0040 | 0.0058 | 0.0089 |
| PCE | 0.0061 | 0.0036 | 0.0043 | 0.0043 | 0.0051 | 0.0069 |
| CIBENZ | 0.0054 | 0.0033 | 0.0039 | 0.0040 | 0.0045 | 0.0059 |
| 1112TetCA | 0.0061 | 0.0037 | 0.0042 | 0.0044 | 0.0050 | 0.0065 |
| EtBENZ | 0.0060 | 0.0037 | 0.0045 | 0.0044 | 0.0052 | 0.0069 |
| mpXYL | 0.0064 | 0.0039 | 0.0048 | 0.0046 | 0.0055 | 0.0072 |
| oXYL | 0.0066 | 0.0041 | 0.0050 | 0.0048 | 0.0057 | 0.0074 |
| 1122TetCA | 0.0044 | 0.0027 | 0.0029 | 0.0031 | 0.0036 | 0.0046 |
| 135TMB | 0.0079 | 0.0046 | 0.0059 | 0.0056 | 0.0071 | 0.0093 |
| 124TMB | 0.0078 | 0.0046 | 0.0060 | 0.0055 | 0.0071 | 0.0092 |
| 13DCB | 0.0072 | 0.0041 | 0.0055 | 0.0053 | 0.0063 | 0.0080 |
| 14DCB | 0.0071 | 0.0040 | 0.0054 | 0.0052 | 0.0062 | 0.0079 |
| 12DCB | 0.0070 | 0.0040 | 0.0053 | 0.0051 | 0.0060 | 0.0076 |
| UNDEC | | 0.0026 | 0.0024 | 0.0020 | 0.0031 | 0.0029 |
| NAPH | | 0.0041 | 0.0056 | 0.0054 | 0.0064 | 0.0081 |
| TRIDEC | | | | | | |
| 2MeNAPH | | 0.0043 | 0.0066 | 0.0066 | 0.0080 | 0.0108 |
| PENTADEC | | | | | | |
| Total mass | 0.1177 | 0.0822 | 0.1339 | 0.1334 | 0.1773 | 0.1981 |

Notes:

Values in L/hr

Total mass does not include UNDEC, TRIDEC, PENTADEC (28 compounds)

TABLE B
4 CONSTANT REGRESSION OUTPUT

| | Adjusted Rsq | Standard Error | ln(SR0) | b | -Ea/R | d | Std Error ln(SR0) | Std Error b | Std Error - Ea/R | Std Error d |
|------------|-------------------------|---------------------------|----------------|----------|--------------|----------|----------------------------------|----------------------------|-------------------------------------|----------------------------|
| MTBE | 0.997 | 0.0960 | -3.217 | 0.981 | 2704 | -0.709 | 0.2881 | 0.0062 | 83 | 0.0082 |
| t12DCE | 0.992 | 0.1659 | -1.877 | 0.905 | 2147 | -0.760 | 0.4971 | 0.0100 | 144 | 0.0138 |
| 11DCA | 0.995 | 0.1272 | -1.346 | 0.916 | 1965 | -0.746 | 0.3809 | 0.0077 | 110 | 0.0106 |
| c12DCE | 0.995 | 0.1299 | -1.905 | 0.911 | 2137 | -0.751 | 0.3892 | 0.0078 | 112 | 0.0109 |
| CHCl3 | 0.996 | 0.1260 | -1.841 | 0.912 | 2118 | -0.748 | 0.3776 | 0.0076 | 109 | 0.0105 |
| 111TCA | 0.995 | 0.1279 | -2.684 | 0.902 | 2259 | -0.761 | 0.3836 | 0.0076 | 111 | 0.0106 |
| 12DCA | 0.995 | 0.1263 | -2.161 | 0.908 | 2218 | -0.746 | 0.3786 | 0.0076 | 109 | 0.0106 |
| BENZ | 0.995 | 0.1323 | -2.207 | 0.920 | 2198 | -0.754 | 0.3965 | 0.0080 | 114 | 0.0110 |
| CCl4 | 0.994 | 0.1405 | -3.121 | 0.889 | 2379 | -0.776 | 0.4220 | 0.0083 | 122 | 0.0116 |
| TCE | 0.992 | 0.1655 | -3.338 | 0.900 | 2522 | -0.772 | 0.4969 | 0.0099 | 144 | 0.0137 |
| 112TCA | 0.995 | 0.1264 | -2.412 | 0.896 | 2302 | -0.724 | 0.3790 | 0.0075 | 109 | 0.0107 |
| TOL | 0.994 | 0.1426 | -2.873 | 0.916 | 2364 | -0.756 | 0.4281 | 0.0087 | 124 | 0.0119 |
| OCT | 0.938 | 0.4698 | -5.984 | 0.822 | 3235 | -0.827 | 1.4231 | 0.0277 | 412 | 0.0388 |
| PCE | 0.991 | 0.1773 | -3.780 | 0.877 | 2601 | -0.775 | 0.5329 | 0.0103 | 154 | 0.0147 |
| CIBENZ | 0.994 | 0.1457 | -2.601 | 0.911 | 2292 | -0.747 | 0.4370 | 0.0088 | 126 | 0.0122 |
| 1112TetCA | 0.996 | 0.1235 | -2.676 | 0.898 | 2281 | -0.725 | 0.3705 | 0.0073 | 107 | 0.0104 |
| EtBENZ | 0.993 | 0.1597 | -2.930 | 0.918 | 2357 | -0.752 | 0.4794 | 0.0097 | 138 | 0.0134 |
| mpXYL | 0.992 | 0.1678 | -3.036 | 0.909 | 2372 | -0.749 | 0.5037 | 0.0101 | 145 | 0.0140 |
| oXYL | 0.993 | 0.1555 | -2.862 | 0.911 | 2312 | -0.740 | 0.4667 | 0.0094 | 135 | 0.0131 |
| 1122TetCA | 0.996 | 0.1118 | -1.971 | 0.913 | 2167 | -0.691 | 0.3351 | 0.0067 | 97 | 0.0096 |
| 135TMB | 0.988 | 0.2024 | -4.435 | 0.897 | 2720 | -0.738 | 0.6093 | 0.0121 | 176 | 0.0170 |
| 124TMB | 0.989 | 0.1997 | -4.126 | 0.890 | 2631 | -0.731 | 0.6009 | 0.0118 | 173 | 0.0169 |
| 13DCB | 0.991 | 0.1832 | -3.422 | 0.888 | 2449 | -0.730 | 0.5503 | 0.0108 | 159 | 0.0155 |
| 14DCB | 0.991 | 0.1802 | -3.263 | 0.892 | 2408 | -0.724 | 0.5413 | 0.0107 | 156 | 0.0153 |
| 12DCB | 0.992 | 0.1697 | -2.970 | 0.894 | 2327 | -0.716 | 0.5092 | 0.0101 | 147 | 0.0144 |
| UNDEC | 0.694 | 0.374 | -1.406 | 0.426 | 1708 | -0.806 | 1.792 | 0.028 | 517 | 0.053 |
| NAPH | 0.992 | 0.166 | -3.374 | 0.915 | 2430 | -0.671 | 0.497 | 0.010 | 144 | 0.014 |
| TRIDEC | | | | | | | | | | |
| 2MeNAPH | 0.984 | 0.238 | -5.498 | 0.869 | 2990 | -0.689 | 0.72 | 0.014 | 208 | 0.021 |
| PENTADEC | | | | | | | | | | |
| Total mass | 0.993 | 0.1543 | -6.111 | 0.907 | 2419 | -0.732 | 0.4666 | 0.0093 | 134 | 0.0130 |

TABLE C
8260C MASS UNCERTAINTY

**AGI 8260C Method for Mass using SPG-0008
Samplers**

| | 99% Uncertainty Range +/- | 95% Uncertainty Range +/- |
|-----------|---------------------------------|---------------------------------|
| MTBE | 20% | 14% |
| t12DCE | 22% | 15% |
| 11DCA | 18% | 12% |
| c12DCE | 18% | 12% |
| CHCl3 | 16% | 11% |
| 111TCA | 18% | 12% |
| 12DCA | 20% | 13% |
| BENZ | 16% | 10% |
| CCl4 | 19% | 12% |
| TCE | 15% | 10% |
| 112TCA | 18% | 12% |
| TOL | 15% | 10% |
| OCT | 20% | 13% |
| PCE | 16% | 11% |
| CIBENZ | 18% | 12% |
| 1112TetCA | 19% | 13% |
| EtBENZ | 18% | 12% |
| mpXYL | 18% | 12% |
| oXYL | 18% | 12% |
| 1122TetCA | 23% | 15% |
| 135TMB | 21% | 14% |
| 124TMB | 20% | 14% |
| 13DCB | 19% | 13% |
| 14DCB | 19% | 13% |
| 12DCB | 20% | 14% |
| NAPH | 21% | 14% |
| 2MeNAPH | 25% | 17% |

TABLE D
4 CONSTANT WATER CONCENTRATION UNCERTAINTY
ERROR IN CONCENTRATION REPORTING (1)

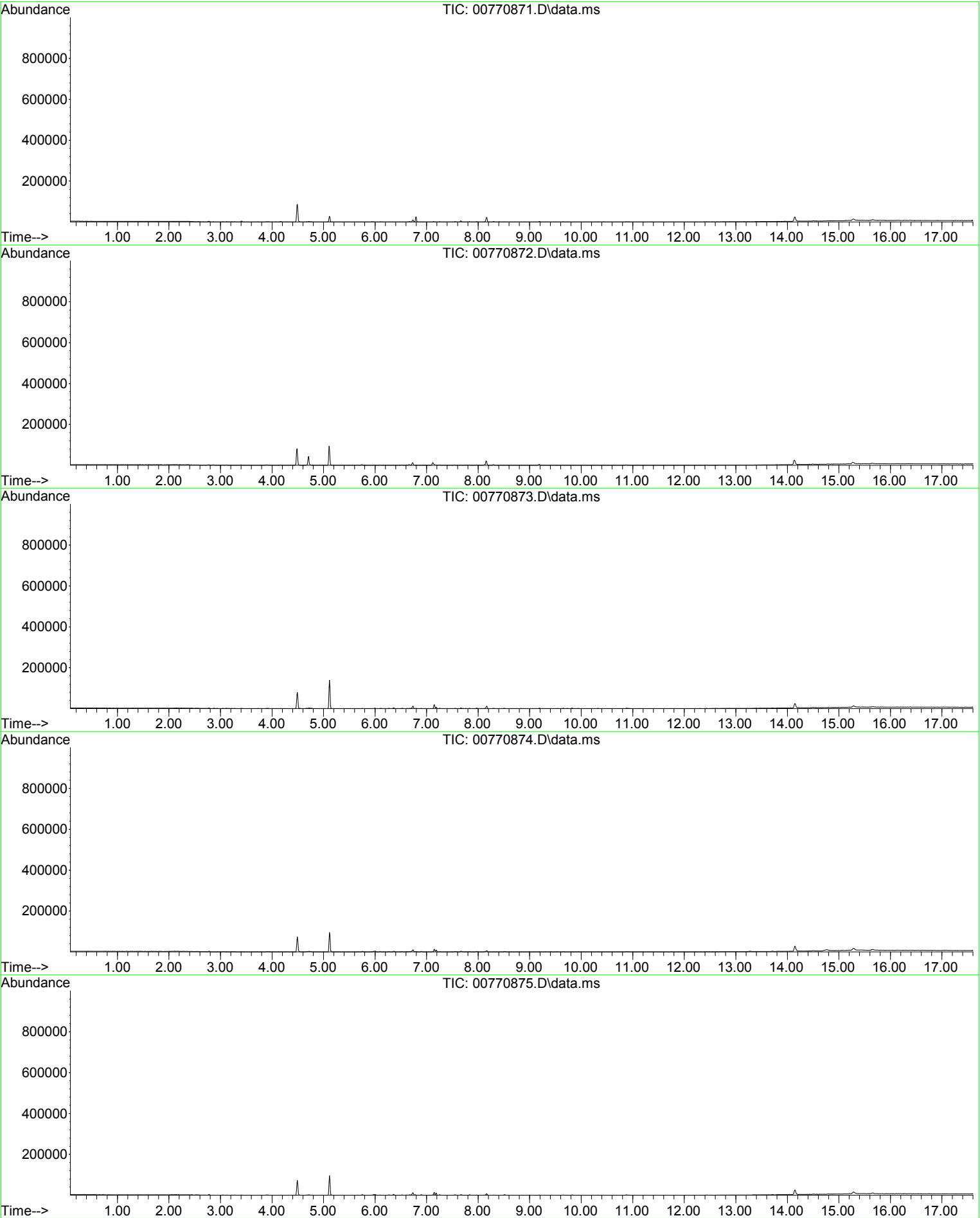
| | Average Error | Minimum Error | Maximum Error |
|-----------|--------------------------|--------------------------|--------------------------|
| MTBE | 6% | -12% | 12% |
| t12DCE | 11% | -26% | 21% |
| 11DCA | 8% | -19% | 13% |
| c12DCE | 9% | -19% | 15% |
| CHCl3 | 9% | -20% | 14% |
| 111TCA | 9% | -19% | 23% |
| 12DCA | 10% | -19% | 17% |
| BENZ | 8% | -18% | 13% |
| CCl4 | 10% | -23% | 22% |
| TCE | 10% | -21% | 14% |
| 112TCA | 11% | -21% | 21% |
| TOL | 7% | -17% | 14% |
| OCT | 20% | -41% | 42% |
| PCE | 10% | -24% | 15% |
| CIBENZ | 7% | -16% | 14% |
| 1112TetCA | 8% | -17% | 18% |
| EtBENZ | 6% | -19% | 14% |
| mpXYL | 7% | -22% | 13% |
| oXYL | 7% | -19% | 13% |
| 1122TetCA | 8% | -16% | 17% |
| 135TMB | 9% | -23% | 17% |
| 124TMB | 10% | -28% | 19% |
| 13DCB | 10% | -22% | 17% |
| 14DCB | 10% | -22% | 17% |
| 12DCB | 9% | -23% | 17% |
| NAPH | 10% | -24% | 21% |
| 2MeNAPH | 13% | -32% | 30% |

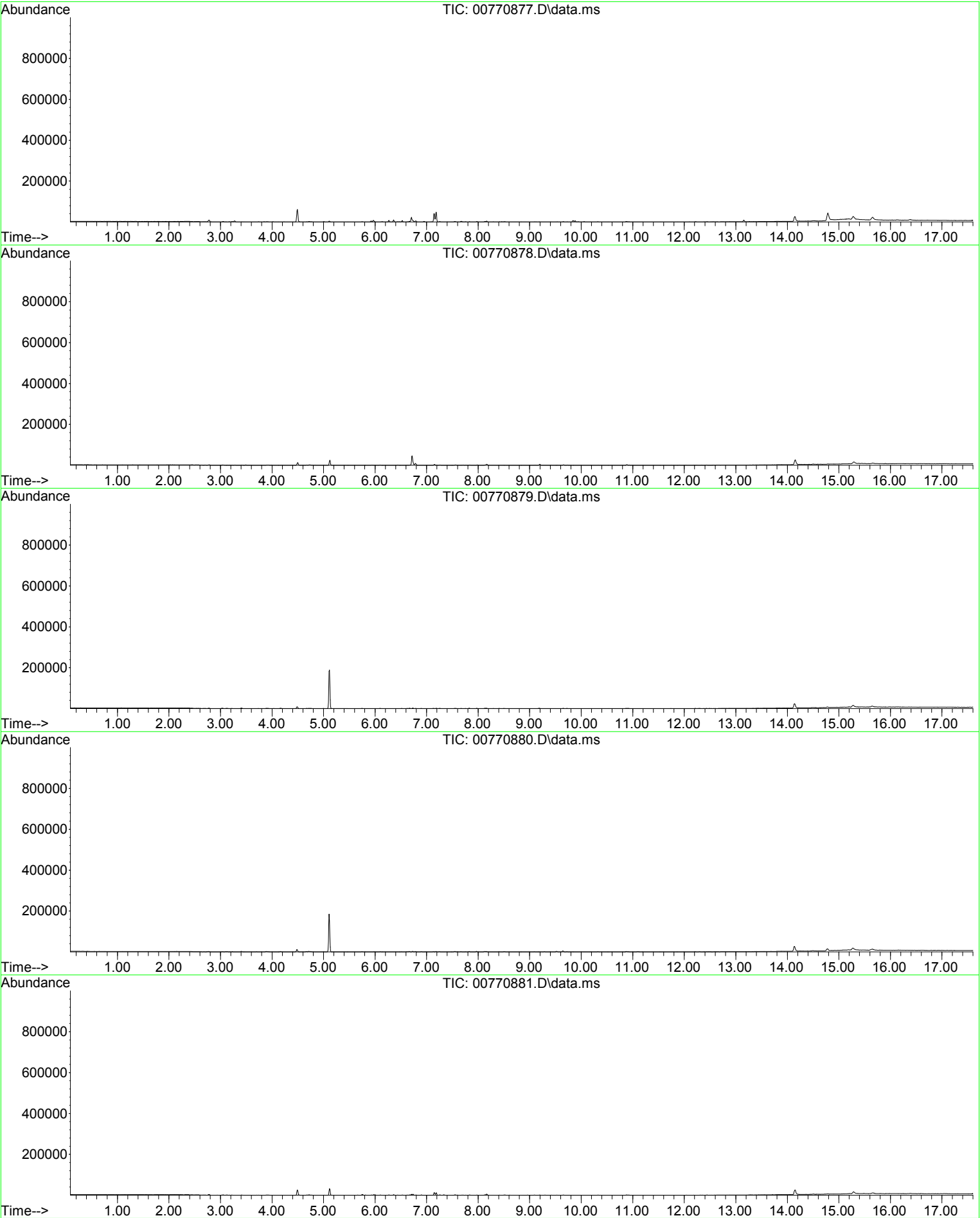
(1) For 1 hour exposure, includes error related to mass value from AGI analytical method 8260C

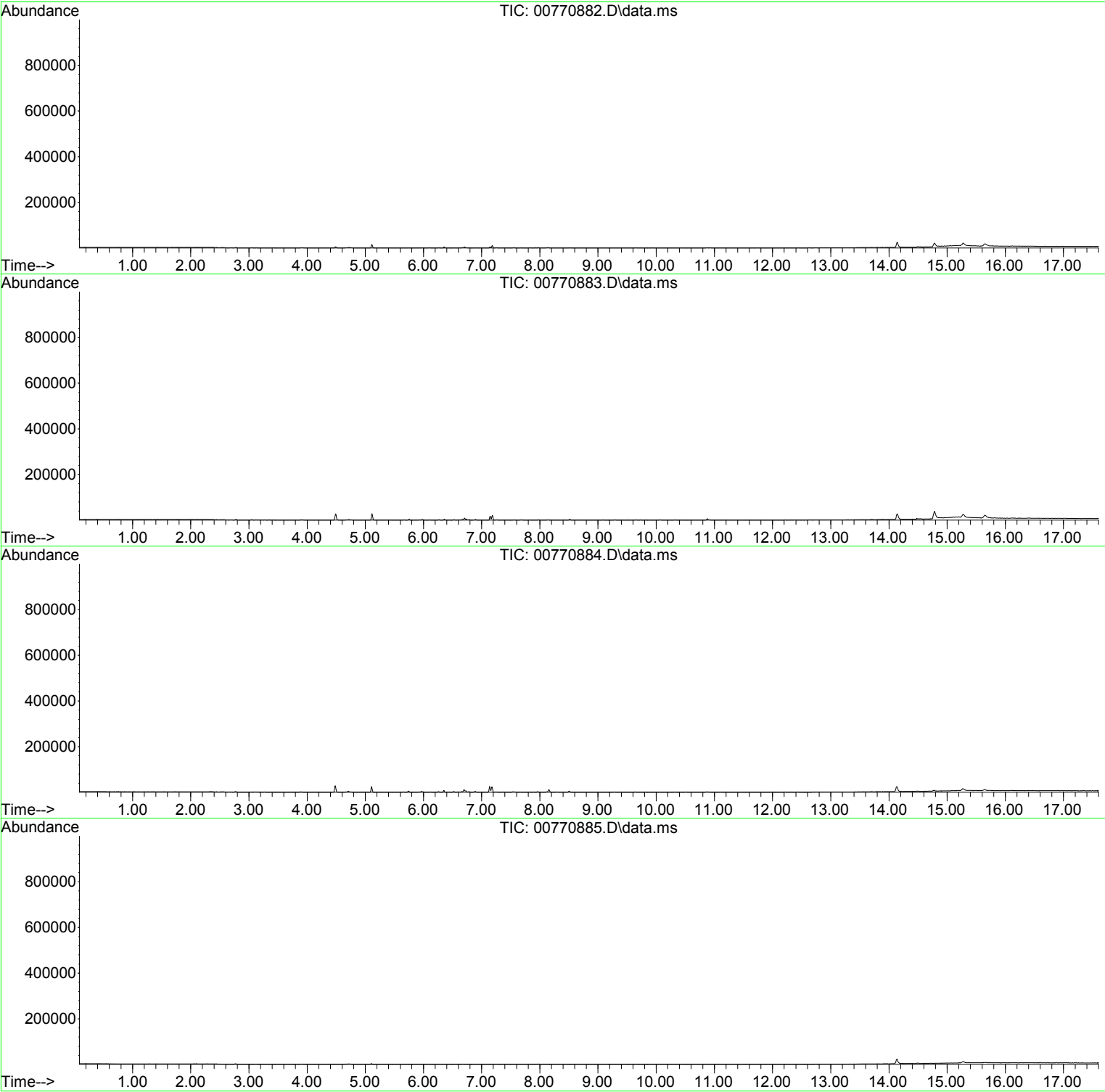
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APPENDIX H

ANALYTICAL DATA VALIDATION DISCUSSION

Introduction

This appendix provides an overview of the SWCA's post analyses review of the Eurofins/Calscience (Calscience), San Antonio River Authority (SARA), and Amplified Geochemical Imaging, LLC (AGI) analytical data set. In general, the data are considered valid for the intended purpose of assessing the baseline of sediment quality, stormwater runoff quality, surface water (under base flow conditions) and passive diffusion samplers (PDS) at a screening level for Comal and San Marcos Springs. Analyses with any associated laboratory issues are listed herein.

Worth noting in this assessment are the detection of plasticizer compounds such as bis(2-ethylhexyl) phthalate (DEHP) and other phthalate compounds. While it is possible these compounds were introduced post-sample collection from either sampling equipment, or laboratory equipment, they should not be completely dismissed.

Based on analysis of 2013 laboratory data, the EAA concluded that three compounds detected in sediment samples may have been laboratory artifacts. The compounds were DEHP, di-n-octyl phthalate, and di-n-butyl phthalate. The EAA noted in the 2013 *Edwards Aquifer Habitat Conservation Plan Expanded Water Quality Report* that as the data set grows, additional conclusions could be drawn. The 2014 laboratory analyses of sediment samples did not detect di-n-octyl phthalate or di-n-butyl phthalate. However, DEHP was detected in three of the sediment samples in 2014 (HSM320, HSM330, and HSM350). DEHP was also detected in the three of the 2015 sediment samples (HSM340, HSM350, and HSM360).

DEHP was not detected in the laboratory method blank or trip blank, and a different laboratory and sampling equipment was used by SWCA. Therefore, it appears DEHP is present within sediment samples in the middle reaches of the San Marcos Spring complex. Future analyses may help better understand the possible presence of DEHP within sediments in both Comal and San Marcos Spring complexes.

Analytical results are discussed by analytical laboratory sample data group number, and by sample event type and date. Each event (surface water/base flow, stormwater, PDS, or sediment) is discussed by sample data group with sample names and date outlined for each event in the beginning of the discussion.

A key to sample names is provided below:

Key to Sample Names

H CS 1 10

H=HCP

CS=Comal Springs (**SM**=San Marcos Springs)

1=Sample Type (1=Surface Water, 2=Storm, 3=Sediment, 4=PDS)

10=Sample Location

Field Duplicates are identified with the prefix FD followed by the sample identification described above. Trip Blank samples are denoted with the prefix "TB" followed by a sequential number. Equipment Blank samples are denoted with the prefix "EB" followed by a sequential number.

Corrective Action Comments

The two following bulleted items describe corrective actions from 2014 that were effective in at reducing problems:

- TPH was consistently detected in PDS Trip Blanks from AGI. AGI had provided an annual supply of PDS. SWCA returned unused PDS and requested separate, new batches of PDS for the December 2014 and each individual 2015 sampling event.
- Calscience packaged three VOA vials within a tight zip-lock bag within a bubble wrap pouch. Frequently, some of the VOA vials broke during return shipping. During the last two sampling events of 2014 and all events of 2015, SWCA individually wrapped the VOA vials as opposed to having the three vials packaged together.

Analytical Data Review Summary for HCP samples collected in 2015.

Data Group Numbers (HCP surface water/base flow samples collected March 16, 2015, and March 25, 2015, Comal and San Marcos springs):

| | |
|-------------------------------------|-------------------------------------|
| <i>15-03-1281-1</i> (HCS 110) | <i>15-03-2020-1</i> (HSM 110) |
| <i>15-03-1281-2</i> (HCS 120) | <i>15-03-2020-2</i> (HSM 120) |
| <i>15-03-1281-3</i> (HCS 130) | <i>15-03-2020-3</i> (HSM 130) |
| <i>15-03-1281-4</i> (HCS 140) | <i>15-03-2020-4</i> (HSM 140) |
| <i>15-03-1281-5</i> (HCS 160) | <i>15-03-2020-5</i> (HSM 150) |
| <i>15-03-1281-6</i> (FDHCS 120) | <i>15-03-2020-6</i> (HSM 160) |
| <i>15-03-1281-7</i> (Trip Blank 02) | <i>15-03-2020-7</i> (HSM 170) |
| | <i>15-03-2020-8</i> (FDHSM110) |
| | <i>15-03-2020-9</i> (Trip Blank 04) |

General Comments

Although some analytical issues are noted for the data group, unless otherwise noted in the detailed discussion, the data are considered valid for the purposes of the investigation. pH values are collected in the field at the time of sample collection and are listed in the field parameters for each sample event. As such, the laboratory pH values which are flagged for hold time exceedances are not used for sample assessment purposes.

Trip Blanks

There were no detections in the trip blanks associated with this sample set.

QA/QC Discussion – Comal and San Marcos Springs Surface/Base Flow Samples (Sampled March 16, 2015 and March 25, 2015)

Issues associated with work order 15-03-1281

Magnesium was detected in the Method Blank associated with all of the Comal surface samples at a level of 0.00760 J mg/L. The detected concentration is below the Reporting Limit, but above the Method Detection Limit. It was therefore “J” flagged. Due to the low concentration and lack of concerns with magnesium concentrations in the subject samples, the results are accepted.

Method EPA 365.1 (Batch No. 150323LO1) – Recovery of matrix spike (MS) and matrix spike duplicates (MSD) were outside of control limits due to suspected matrix interference for total phosphorous. The associated Laboratory Control Sample (LCS) was in control; therefore, the results are accepted.

Method SM 5310 B (Batch No. F0326DOCS1) – Recovery of the MSD compound was outside of control limits due to suspected matrix interference for dissolved organic carbon (DOC). The associated Laboratory Control Sample (LCS) was in control; therefore, the results are accepted.

Method EPA 6010B (Batch No. 150330SA2) - For analytes calcium, magnesium, sodium, and silicon, spike recovery and relative percent difference (RPD) control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method EPA 6020 (Batch No. 150319SA1) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for aluminum. The associated LCS was in control; therefore, the results are accepted. The recovery of the pulse digestion spike was also out of controls for aluminum due to suspected matrix interference. The result was biased high, the result is accepted.

Issues associated with work order 15-03-2020

Magnesium was detected in the Method Blank at a level of 0.0449 J mg/L. The detected concentration is below the Reporting Limit, but above the Method Detection Limit. It was therefore “J” flagged. Due to the low concentration and lack of concerns with magnesium concentrations in the subject samples, the results are accepted.

Method EPA 6010B (Batch No. 150327SA4A) - For analytes calcium, magnesium, and silicon, spike recovery and RPD control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Data Group Numbers (HCP surface water/base flow samples collected September 9 and 17, 2015, Comal and San Marcos springs):

| | |
|-------------------------------------|-------------------------------------|
| <i>15-09-0733-1</i> (HCS 110) | <i>15-09-1440-1</i> (HSM 110) |
| <i>15-09-0733-2</i> (HCS 120) | <i>15-09-1440-2</i> (FDHSM 110) |
| <i>15-09-0733-3</i> (FDHCS 120) | <i>15-09-1440-3</i> (HSM 120) |
| <i>15-09-0733-4</i> (HCS 130) | <i>15-09-1440-4</i> (HSM 130) |
| <i>15-09-0733-5</i> (HCS 140) | <i>15-09-1440-5</i> (HSM 140) |
| <i>15-09-0733-6</i> (HCS 160) | <i>15-09-1440-6</i> (HCS 150) |
| <i>15-09-0733-7</i> (Trip Blank 11) | <i>15-09-1440-7</i> (HSM 160) |
| | <i>15-09-1440-8</i> (HSM 170) |
| | <i>15-09-1440-9</i> (Trip Blank 12) |

General Comments

Although some analytical issues are noted for the data group, unless otherwise noted in the detailed discussion, the data are considered valid for the purposes of the investigation. pH values are collected in the field at the time of sample collection and are listed in the field parameters for each sample event. As such, the laboratory pH values which are flagged for hold time exceedances are not used for sample assessment purposes.

Trip Blank (Comal and San Marcos)

Trip Blank (TB11) associated with this sample set had a detection of 0.25 J ug/L for toluene. This constituent was not detected in the subject samples.

**QA/QC Discussion – Comal and San Marcos Springs Surface/Base Flow Samples
(Sampled September 9 and 17, 2015)**

Issues associated with work order 15-09-0733

Calcium, magnesium, potassium and sodium were detected in the Method Blank associated with all of the Comal surface samples. The detected concentrations are below the Reporting Limits, but above the Method Detection Limits. They were therefore “J” flagged. Due to the low concentration and lack of concerns with these analyte concentrations in the subject samples, the results are accepted.

Method EPA 300.0 (Batch No. 150910S01) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for chloride and sulfate. The associated LCSs were in control; therefore, the results are accepted.

Method EPA 365.1 (Batch No. 150915S01) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for total phosphorous. The associated LCS was in control; therefore, the results are accepted.

Method EPA 6010B (Batch No. 150911SA5) - For analytes calcium, magnesium, and silicon, spike recovery and relative percent difference (RPD) control limits do not apply—resulting from the parameter

concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method EPA 8141A (Batch No. 150911L03) – LCS and LCSD percent recoveries were out of control for merphos. This compound was not detected in the subjected samples. The results could be biased high; therefore, the results are accepted.

Method 1694 - Caffeine analysis was subcontracted to ALS Environmental by Calscience. The subcontracted laboratory received less than optimal sample volume which resulted in elevated detection limits. In the future, adequate sample volume will be provided to ALS Environmental by Calscience.

Issues associated with work order 15-09-1440

Magnesium was detected in the Method Blank associated with all of the San Marcos surface samples. The detected concentration, 0.00711 mg/L, is below the Reporting Limit, but above the Method Detection Limit. The result is therefore “J” flagged. Due to the low concentration and lack of concerns with this analyte in the subject samples, the results are accepted.

Method EPA 300.0 (Batch No. 150918S01) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for nitrate and sulfate. The associated LCSs were in control; therefore, the results are accepted.

Method EPA 365.1 (Batch No. 151002S01) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for nitrate and sulfate. The associated LCSs were in control; therefore, the results are accepted.

Method EPA 6010B (Batch No. 150922SA6A) - For analytes calcium, magnesium, and silicon, spike recovery and relative percent difference (RPD) control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method EPA 6020 (Batch No. 150922SA3A) - For manganese, spike recovery and relative percent difference (RPD) control limits do not apply to the MS, MSD or pulse digestion spike samples—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method EPA 6010B (Batch No. 150922LA6F) – The LCS percent recovery was out of range for silicon. Results may be biased low but are accepted.

Method EPA 8141A (Batch No. 150124L03) – LCS and LCSD percent recoveries were out of control for merphos. This compound was not detected in the subjected samples. The results could be biased high; therefore, the results are accepted.

Data Group Numbers (HCP stormwater samples collected January 22-23, 2014, at Comal Springs):

| | |
|------------------------------------|--|
| <i>15-01-1404-1</i> (HCS 210 Lead) | <i>15-01-1511-5</i> (HCS 270 Peak) |
| <i>15-01-1404-2</i> (HCS 240 Lead) | <i>15-01-1511-6</i> (HCS 210 Trail) |
| <i>15-01-1404-3</i> (HCS 250 Lead) | <i>15-01-1511-7</i> (HCS 240 Trail) |
| <i>15-01-1404-4</i> (HCS 260 Lead) | <i>15-01-1511-8</i> (HCS 250 Trail) |
| <i>15-01-1404-5</i> (HCS 270 Lead) | <i>15-01-1511-9</i> (HCS 260 Trail) |
| <i>15-01-1511-1</i> (HCS 210 Peak) | <i>15-01-1511-10</i> (HCS 270 Trail) |
| <i>15-01-1511-2</i> (HCS 240 Peak) | <i>15-01-1511-11</i> (FDHCS 260 Trail) |
| <i>15-01-1511-3</i> (HCS 250 Peak) | <i>15-01-1511-12</i> (FDHCS 270 Trail) |
| <i>15-01-1511-4</i> (HCS 260 Peak) | <i>15-01-1511-13</i> (Trip Blank 1) |

General Comments

Although some analytical issues are noted for the data group, unless otherwise noted in the detailed discussion, the data are considered valid for the purposes of the investigation. pH values are collected in the field at the time of sample collection and are listed in the field parameters for each sample event. As such, the laboratory pH values which are flagged for hold time exceedances are not used for sample assessment purposes.

All samples were collected outside of laboratory operating hours. Therefore all *E. coli* samples were analyzed outside of their hold times. The initial blank for *E. coli* was also analyzed outside of hold times for batch E_COLI_QUANTITRAY-41482 associated with samples HCS 250 Trail, HCS 260 Trail, HCS 270 Trail, FDHCS 260 Trail and FDHCS 270 Trail.

Trip Blank

There were no detections in the trip blank associated with these samples.

**QA/QC Discussion – Comal Springs Stormwater Samples
(Sampled January 22-23, 2015)**

Issues associated with all January 2015 stormwater samples

Magnesium was detected at a concentration of 0.0522 J mg/L in batch 150126L6FF within the Method Blank, which is below the Reporting Limit, but above the Method Detection Limit. It was therefore “J” flagged. Due to the low concentration and lack of issues with magnesium concentrations in the subject samples, the results are accepted.

Method EPA 6010B (Batch No. 150126SA6) - For analytes calcium and magnesium, spike recovery and RPD control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method EPA 6010B (Batch No. 150126L6FF) – Percent recovery of silicon in the LCS was outside of the control limit and results may be biased low. Due to lack of issues with silicon concentrations in the subject samples, the results are accepted.

Issues specific to work order 15-01-1404

Method EPA 365.1 (Batch No. 150202SO1) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for total phosphorous. The associated LCS was in control; therefore, the results are accepted.

Method EPA 8141A (Batch No. 150124L03) – LCS and LCSD percent recoveries were out of control for merphos. This compound was not detected in the subjected samples. The results could be biased high; therefore, the results are accepted.

Issues specific to individual samples

HCS 260 Trail Method EPA 8082 – Recovery of the surrogate compounds was outside of control limits. Results may be biased high; however, no compounds were detected in the sample and the method blank surrogate recovery was in control. The results are accepted.

Data Group Numbers (HCP stormwater samples collected May 5-6, 2015, at San Marcos springs):

| | |
|-------------------------------------|--|
| <i>15-05-0402-1</i> (HSM 210 Lead) | <i>15-05-0402-14</i> (HSM 270 Peak) |
| <i>15-05-0402-2</i> (HSM 230 Lead) | <i>15-05-0402-15</i> (Trip Blank 06) |
| <i>15-05-0402-3</i> (HSM 231 Lead) | <i>15-05-0402-16</i> (HSM 210 Trail) |
| <i>15-05-0402-4</i> (HSM 240 Lead) | <i>15-05-0402-17</i> (HSM 230 Trail) |
| <i>15-05-0402-5</i> (HSM 250 Lead) | <i>15-05-0402-18</i> (HSM 231 Trail) |
| <i>15-05-0402-6</i> (HSM 260 Lead) | <i>15-05-0402-19</i> (HSM 240 Trail) |
| <i>15-05-0402-7</i> (HSM 270 Lead) | <i>15-05-0402-20</i> (HSM 250 Trail) |
| <i>15-05-0402-8</i> (HSM 210 Peak) | <i>15-05-0402-21</i> (HSM 260 Trail) |
| <i>15-05-0402-9</i> (HSM 230 Peak) | <i>15-05-0402-22</i> (HSM 270 Trail) |
| <i>15-05-0402-10</i> (HSM 231 Peak) | <i>15-05-0402-23</i> (FDHSM 210 Trail) |
| <i>15-05-0402-11</i> (HSM 240 Peak) | <i>15-05-0402-24</i> (FDHSM 230 Trail) |
| <i>15-05-0402-12</i> (HSM 250 Peak) | <i>15-05-0402-25</i> (FDHSM 231 Trail) |
| <i>15-05-0402-13</i> (HSM 260 Peak) | |

General Comments

Although some analytical issues are noted for the data group, unless otherwise noted in the detailed discussion, the data are considered valid for the purposes of the investigation. pH values are collected in the field at the time of sample collection and are listed in the field parameters for each sample event. As such, the laboratory pH values which are flagged for hold time exceedances are not used for sample assessment purposes. All samples were collected outside of laboratory operating hours. Therefore all *E. coli* samples were analyzed outside of their hold times.

Trip Blanks

There were no detections in the trip blank associated with these samples.

**QA/QC Discussion – San Marcos Springs Stormwater Samples
(Sampled May 5 and 6, 2015)**

Issues associated with specific batches of stormwater samples

Method EPA 6010B (Batch No. 150507LA6A) - Magnesium was detected in the Method Blank at a level of 0.00476 J mg/L. The detected concentration is below the Reporting Limit, but above the Method Detection Limit. It was therefore “J” flagged. Due to the low concentration and lack of concerns with magnesium concentrations in the subject samples, the results are accepted.

Method EPA 365.1 (Batch No. 150520S01) – Recovery of the MS and MSD compound was outside of control limits due to suspected matrix interference for total phosphorous. The associated LCS was in control; therefore, the results are accepted.

Method EPA 6010B (Batch No. 150507SA5) - For analytes calcium, magnesium, and silicon, spike recovery and RPD control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method EPA 6010B (Batch No. 150507SA6) - For analytes calcium, magnesium, and silicon, spike recovery and RPD control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted. Recovery of the MS and MSD compound was out of control due to suspected matrix interference for strontium. The results may be biased high; however, the LCS was in control and the results are accepted.

Method EPA 6020 (Batch No. 150507SA4) – The MS/MSD percent difference was out of control due to suspected matrix interference for antimony, arsenic, beryllium, cadmium, chromium, copper, lead, nickel, selenium, thallium, and zinc. Recovery of the MS compound was outside of control limits due to suspected matrix interference for cadmium. The associated LCS was in control; therefore, the results are accepted.

Samples HSM 270 Trail, FDHSM 210 Trail, FDHSM 230 Trail, and FDHSM 231 Trail Method EPA 6020 (Batch No. 150507LA6F) - Arsenic was detected in the Method Blank at a level of 0.000469 J mg/L. The detected concentration is below the Reporting Limit, but above the Method Detection Limit. It was therefore “J” flagged. The detections in these results are similar to the levels detected in the Lead and Peak samples. However, none of the other trail samples had arsenic detections. These results may be biased high but are within the range of results in the unaffected samples. Results are accepted.

Method SM 6020 (Batch No. 150507SA6) – Recovery of the MSD compound was outside of control limits due to suspected matrix interference for antimony, arsenic, and cadmium. The associated LCS was in control; therefore, the results are accepted.

Method EPA 8260B (Batch No. 150515S020) – The MS/MSD percent difference was out of control due to suspected matrix interference for methyl-t-butyl ether (MTBE). MS and MSD percent recoveries both in acceptable range, therefore the results are accepted.

Method EPA 8260B (Batch No. 150518S007) – Recovery of the MSD compound was outside of control limits due to suspected matrix interference for carbon tetrachloride and 1,1-dichloroethene. The associated LCS was in control; therefore, the results are accepted.

Issues specific to individual samples

HSM 240 Peak Method EPA 8081A – Recovery of one surrogate compound was outside of control limits. Results may be biased low. The method blank surrogate recovery was in control. The results are accepted.

HSM 230 Lead Method EPA 8141A – Recovery of the surrogate compound was outside of control limits. Results may be biased high; however, no compounds were detected in the sample and the method blank surrogate recovery was in control. The results are accepted.

Data Group Numbers (HCP stormwater samples collected October 23, 2015, at Comal Springs and San Marcos Springs):

| | |
|---------------------------------------|--|
| <i>15-10-1856-1</i> (HCS 210 Lead) | <i>15-10-1995-21</i> (HSM 210 Lead) |
| <i>15-10-1856-2</i> (HCS 240 Lead) | <i>15-10-1995-22</i> (HSM 230 Lead) |
| <i>15-10-1856-3</i> (HCS 250 Lead) | <i>15-10-1995-23</i> (HSM 231 Lead) |
| <i>15-10-1856-4</i> (HCS 260 Lead) | <i>15-10-1995-24</i> (HSM 240 Lead) |
| <i>15-10-1856-5</i> (HCS 270 Lead) | <i>15-10-1995-25</i> (HSM 250 Lead) |
| <i>15-10-1856-7</i> (HCS 210 Peak) | <i>15-10-1995-26</i> (HSM 260 Lead) |
| <i>15-10-1856-8</i> (HCS 240 Peak) | <i>15-10-1995-27</i> (HSM 270 Lead) |
| <i>15-10-1856-9</i> (HCS 250 Peak) | <i>15-10-1995-28</i> (HSM 210 Peak) |
| <i>15-10-1995-1</i> (HCS 260 Peak) | <i>15-10-1995-29</i> (HSM 230 Peak) |
| <i>15-10-1995-2</i> (HCS 270 Peak) | <i>15-10-1995-30</i> (HSM 231 Peak) |
| <i>15-10-1995-3</i> (HCS 210 Trail) | <i>15-10-1995-31</i> (HSM 240 Peak) |
| <i>15-10-1995-4</i> (HCS 240 Trail) | <i>15-10-1995-32</i> (HSM 250 Peak) |
| <i>15-10-1995-5</i> (HCS 250 Trail) | <i>15-10-1995-33</i> (HSM 260 Peak) |
| <i>15-10-1995-6</i> (HCS 260 Trail) | <i>15-10-1995-34</i> (HSM 270 Peak) |
| <i>15-10-1995-7</i> (HCS 270 Trail) | <i>15-10-1995-11</i> (HSM 210 Trail) |
| <i>15-10-1995-8</i> (FDHCS 260 Trail) | <i>15-10-1995-12</i> (HSM 230 Trail) |
| <i>15-10-1995-9</i> (FDHCS 270 Trail) | <i>15-10-1995-13</i> (HSM 231 Trail) |
| | <i>15-10-1995-14</i> (HSM 240 Trail) |
| <i>15-10-1856-6</i> (Trip Blank 14) | <i>15-10-1995-15</i> (HSM 250 Trail) |
| <i>15-10-1995-10</i> (Trip Blank 15) | <i>15-10-1995-16</i> (HSM 260 Trail) |
| | <i>15-10-1995-17</i> (HSM 270 Trail) |
| | <i>15-10-1995-18</i> (FDHSM 210 Trail) |
| | <i>15-10-1995-19</i> (FDHSM 230 Trail) |
| | <i>15-10-1995-20</i> (FDHSM 231 Trail) |

General Comments

Although some analytical issues are noted for the data group, unless otherwise noted in the detailed discussion, the data are considered valid for the purposes of the investigation. pH values are collected in the field at the time of sample collection and are listed in the field parameters for each sample event. As such, the laboratory pH values which are flagged for hold time exceedances are not used for sample assessment purposes. Because sample collection occurred late in the day on a Friday, samples could not be packaged and delivered to FedEx in time for a Saturday delivery. Therefore, samples associated with work order 15-10-1995 were held on ice at SWCA over the weekend and shipped to the laboratory Monday, October 26, 2015. This resulted in the nitrate samples from this work order being received outside of hold times.

The majority of samples were collected outside of laboratory operating hours. Therefore many *E. coli* samples were analyzed outside of their hold times. In addition, *E. coli* sample HSM230 Lead contained limited air space and samples HCS260 Peak, HSM231 Lead and FDHSM230 Trail contained limited sample volume.

High sample concentrations of caffeine required that samples HCS210 Lead and HCS250 Lead be diluted for effective analysis.

Trip Blank

There were no detections in the trip blanks associated with these samples.

QA/QC Discussion – Comal and San Marcos Springs Stormwater Samples (Sampled October 23, 2015)

Issues specific to work order 15-10-1856

Method EPA 365.1 (Batch No. 151109S02) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for total phosphorous. The associated LCS was in control; therefore, the results are accepted.

Method SM 5310 B (Batch No. F1105TOCS1) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for total organic carbon. The associated LCS was in control; therefore, the results are accepted.

Method EPA 6010B (Batch No. 151026SA4A) - For analytes calcium, magnesium, and silicon spike recovery and RPD control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method EPA 6020 (Batch No. 151026SA1) – The MS/MSD RPD was out of control for nickel due to suspected matrix interference. The MS and MSD were each within control. Recovery of the MS for aluminum was below the control due to suspected matrix interference. Spike recovery and RPD control limits do not apply for manganese for the MS, MSD or pulse digestion spike samples due to the sample concentration exceeding the spike concentration by a factor of four or greater. The results are accepted.

Issues specific to work order 15-10-1995

Method 6010B (Batch No. 151028LA8F)- Calcium was detected in the Method Blank at a level of 0.0244 J mg/L. The detected concentration is below the Reporting Limit, but above the Method Detection Limit. It was therefore “J” flagged. Due to the low concentration and lack of concerns with calcium concentrations in the subject samples, the results are accepted.

Method 6020 (Batch No. 151028LA5F)- Zinc was detected in the Method Blank at a level of 0.00185 J mg/L. The detected concentration is below the Reporting Limit, but above the Method Detection Limit. It was therefore “J” flagged. Due to the low concentration and lack of concerns with zinc concentrations in the subject samples, the results are accepted.

Method EPA 300.0 (Batch No. 151031S01) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for sulfate. The associated LCS was in control; therefore, the results are accepted.

Method EPA 365.1 (Batch No. 151110S01) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for total phosphorous. The associated LCS was in control; therefore, the results are accepted.

Method EPA 365.1 (Batch No. 151110S02) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for total phosphorous. The associated LCS was in control; therefore, the results are accepted.

Method SM 5310 B (Batch No. F1109TOCS2) – Recovery of the MS and the MS/MSD RPD were outside of control limits due to suspected matrix interference for total organic carbon. The associated LCS was in control; therefore, the results are accepted.

Method EPA 6010B (Batch No. 151028SA8) – For analytes calcium, magnesium, and silicon spike recovery and RPD control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method EPA 6010B (Batch No. 151028SA9) - For analytes calcium, magnesium, and silicon spike recovery and RPD control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method EPA 6020 (Batch No. 151028SA4) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for zinc. The associated LCS was in control; therefore, the results are accepted.

Method EPA 6020 (Batch No. 151026SA1) – The MS/MSD RPD was out of control for selenium due to suspected matrix interference. The MS and MSD were each within control. The results are accepted.

Method EPA 8260B (Batch No. 151030S009) – Recovery of the MS and MSD were outside of control limits due to suspected matrix interference for vinyl chloride. The associated LCS was in control; therefore, the results are accepted.

Method EPA 6020 (Batch No. 151028SA4) – Recovery of the pulse digestion spike was outside of control limits due to suspected matrix interference for selenium. All detections were below the reporting limit and two orders of magnitude lower than the regulatory standard used for result comparison (MCL); therefore, the results are accepted.

Method EPA 6020 (Batch No. 151028SA5) – Recovery of the pulse digestion spike was outside of control limits due to suspected matrix interference for selenium. All detections were below the reporting limit and two orders of magnitude lower than the regulatory standard used for result comparison (MCL); therefore, the results are accepted.

Issues specific to individual samples

HCS 260 Trail Method EPA 8082 – Recovery of a surrogate compound was above the acceptance limit. Results may be biased high; however, no compounds were detected in the sample and the method blank surrogate recovery was in control. The results are accepted.

HSM 260 Trail Method EPA 8151A – Recovery of the surrogate compound was above the acceptance limit. Results may be biased high; however, no compounds were detected in the sample and the method blank surrogate recovery was in control. The results are accepted.

FDHSM 210 Trail Method EPA 8151A – Recovery of the surrogate compound was above the acceptance limit. Results may be biased high; however, no compounds were detected in the sample and the method blank surrogate recovery was in control. The results are accepted.

HSM 270 Peak Method EPA 8151A – Recovery of the surrogate compound was above the acceptance limit. Results may be biased high; however, no compounds were detected in the sample and the method blank surrogate recovery was in control. The results are accepted.

Data Group Numbers (HCP sediment samples collected June 4 and 5, 2015, at Comal springs):

| | |
|--|---|
| <i>15-06-0463-1</i> (HCS 310) | <i>15-06-0567-1</i> (HSM 310) |
| <i>15-06-0463-2</i> (HCS 320) | <i>15-06-0567-2</i> (HSM 320) |
| <i>15-06-0463-3</i> (HCS 330) | <i>15-06-0567-3</i> (HSM 330) |
| <i>15-06-0463-4</i> (HCS 340) | <i>15-06-0567-4</i> (HSM 340) |
| <i>15-06-0463-5</i> (HCS 360) | <i>15-06-0567-5</i> (HSM 350) |
| <i>15-06-0463-6</i> (FDHCS 360) | <i>15-06-0567-6</i> (HSM 360) |
| <i>15-06-0463-7</i> (Trip Blank 07) | <i>15-06-0567-7</i> (HSM 370) |
| | <i>15-06-0567-8</i> (FDHSM370) |
| | <i>15-06-0567-9</i> (Trip Blank 08) |
| <i>15-06-0567-1</i> (Equipment Blank 02) | <i>15-06-0567-10</i> (Equipment Blank 01) |

General Comments

Although some analytical issues are noted for the data group, unless otherwise noted in the detailed discussion, the data are considered valid for the purposes of the investigation. Samples were collected in glass jars using stainless steel trowels. The water depth and sediment availability at HCS 330 made collection with a trowel impossible. The HCS 330 sample was collected in a plastic sleeve using a hand core sampler and extruded into the glass sample jars.

Trip Blanks

There were no detections in the trip blanks associated with these samples.

Equipment Blanks

Two equipment blanks were collected. One sample was collected by pouring ASTM Type II Reagent Grade water over a decontaminated trowel. The second sample was collected by pouring ASTM Type II Reagent Grade water through a new sample tube. There were several compounds detected in both of the two equipment blanks. The detections are similar between each blank despite the differences in sampling equipment. SWCA suspects that the detected compounds were present in the ASTM Type II Reagent Grade water used. The plastic sample tube was a new, single-use tube from the box and was therefore not decontaminated, eliminating decontamination procedures as the source of interference. Many of the detections were “J” flagged and substantially lower than the levels detected in the sediment samples. Therefore, the equipment blank analyses results are not anticipated to have affected the usability of results of the sediment sample analyses. The equipment blank detections are summarized in Table 1.

Table 1. Detections in sediment equipment blanks June 2015

| Sample | Associated Equipment | Date Collected | Calcium mg/L | Magnesium mg/L | Sodium mg/L | Barium mg/L | Copper mg/L | Nickel mg/L | Zinc mg/L | Manganese mg/L | Chloroform µg/L | Toluene µg/L | Carbon, Total Organic mg/L | Phosphorous, Total mg/L |
|--------|----------------------|----------------|-----------------|-------------------|----------------|----------------|----------------|----------------|--------------|-------------------|--------------------|-----------------|-------------------------------|----------------------------|
| EB01 | Trowel | 6/5/2015 | 0.0548J | 0.0451B,J | 0.150J | <0.0000986 | 0.00104 | 0.000299J | 0.000853J | 0.00158J | 23 | 0.27J | 0.55 | <0.020 |
| EB02 | Sample tube | 6/10/2015 | <0.0118 | 0.0131B,J | <0.103 | 0.000130 | 0.00155 | 0.000425 | 0.00129 | 0.000253 | 26 | 0.45J | 0.78 | 0.020J |

QA/QC Discussion – Comal and San Marcos Springs Sediment Samples (Sampled June 4, and 5, 2015)

Issues associated with work order 15-06-0463

Method EPA 300.0 (Batch No. 150610S01P) – Recovery of the MS and MSD compound was outside of control limits due to suspected matrix interference for fluoride. The associated LCS was in control; therefore, the results are accepted.

Method EPA 6010B (Batch No. 150610S04) - For analytes calcium, magnesium, potassium and silicon, spike recovery and RPD control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method EPA 6020 (Batch No. 150608S01) - For analytes barium, aluminum, iron and manganese, spike recovery and RPD control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method SM 8270C (Batch No. 150613S07) – Recovery of the MS and MSD compound was outside of control limits due to suspected matrix interference for 1,4-dichlorobenzene and 1,2,4-trichlorobenzene. The associated LCS was in control; therefore, the results are accepted.

Method SM 8260B (Batch No. 150613S007) – Recovery of the MS and/or MSD compound was outside of control limits due to suspected matrix interference for benzene, bromodichloromethane, carbon disulfide, chloroform, 1,1-dichloroethene, c-1,2-dichloroethene, t-1,2-dichloroethene, 1,2-dichloropropane, c-1,3-dichloropropene, methylene chloride, toluene, 1,1,1-trichloroethane and trichloroethene. The associated LCS was in control; therefore, the results are accepted.

Method EPA 6020 (Batch No. 150608L01) – LCS recovery percentage was in the Marginal Exceedance (ME) control limit range for iron. This method allows for one ME result; therefore, results are accepted.

Method EPA 8260B (Batch No. 150607L024) – LCS recovery percentage was in the Marginal Exceedance (ME) control limit range for 1,2-dichloroethane. This method allows for one ME result; therefore, results are accepted.

Issues associated with work order 15-06-0463

Method EPA365.1 (Batch No. 150613S01) - Recovery of the MS and MSD compound was out of control due to suspected matrix interference for total phosphorous. The results may be biased high; however, the LCS was in control, and the results are accepted.

Method EPA 6010B (Batch No. 150611S09A) - For analytes calcium, magnesium, and silicon, spike recovery and RPD control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Recovery of the MS and/or MSD compound was out of control due to suspected matrix interference for potassium, sodium, and strontium. The results may be biased high; however, the LCS was in control, and the results are accepted.

Method EPA 6010B (Batch No. 150615SA4) - For analytes calcium, magnesium, sodium and silicon, spike recovery and RPD control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method EPA 6020 (Batch No. 150612S01) - For analytes barium, aluminum, iron, and manganese spike recovery and RPD control limits do not apply—resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. Results are accepted.

Method EPA 8260B (Batch No. 150608S026) – Recovery of the MS and MSD compound was outside of control limits due to suspected matrix interference for 1,2-dibromoethane, 1,2-dichloroethane, p/m-xylene, and o-xylene. The associated LCS was in control; therefore, the results are accepted.

Method EPA 8260B (Batch No. 150608S015) – Recovery of the MS and/or MSD compound was outside of control limits due to suspected matrix interference for benzene, chlorobenzene, 1,2-dibromoethane, 1,2-dichloroethane, toluene, p/m-xylene, and o-xylene. The associated LCS was in control; therefore, the results are accepted.

Issues specific to individual samples

EB01 Method 6010B - Magnesium was detected in the Method Blank at a level of 0.0372 J mg/L. The detected concentration is below the Reporting Limit, but above the Method Detection Limit. It was therefore “J” flagged. Due to the low concentration and lack of concerns with magnesium concentrations in the subject samples, the results are accepted.

HSM 350 Method EPA 8081A – Recovery of the surrogate compounds were outside of control limits. Results may be biased high; however, no compounds were detected in the sample and the method blank surrogate recovery was in control. The results are accepted.

Data Group Numbers (HCP PDS deployed February 3 – February 17, 2015, at Comal and San Marcos springs):

HCS 410 **00755224**

HCS 420 **00755225**

HCS 430 **00755226**

HCS 440 **00755227**

FDHCS 440 **00755228**

HCS 460 **00755228**

Trip Blank 02 **00755228**

HSM 410 **00755230**

HSM 420 **00755231**

HSM 430 **00755232**

FDHSM 430 **00755233**

HSM 440 **00755234**

HSM 450 **00755235**

HSM 460 **00755236**

HSM 470 **00755237**

General Comments

Although some analytical issues are noted for the data group, unless otherwise noted in the detailed discussion, the data are considered valid for the purposes of the investigation. PDSs were deployed from February 3, 2015, through February 17, 2015. The initial method blank data was accidentally overwritten by the laboratory and not available for inclusion in the laboratory report. However, there was no evidence of system contamination observed in any other quality control sample; the results are accepted.

Trip Blanks

There were no detections in the trip blank associated with these samples.

Equipment Blanks

Equipment blanks were not practicable due to sampler type.

Data Group Numbers (HCP PDS deployed April 7 – April 21, 2015, at Comal and San Marcos springs):

| | |
|-------------------------------|---------------------------|
| HCS 410 <i>00757309</i> | HSM 410 <i>00757315</i> |
| HCS 420 <i>00757310</i> | HSM 420 <i>00757316</i> |
| HCS 430 <i>00757311</i> | HSM 430 <i>00757317</i> |
| HCS 440 <i>00757312</i> | FDHSM 430 <i>00757318</i> |
| FDHCS 440 <i>00757313</i> | HSM 440 <i>00757319</i> |
| HCS 460 <i>00757314</i> | HSM 450 <i>00757320</i> |
| | HSM 460 <i>00757321</i> |
| Trip Blank 05 <i>00757323</i> | HSM 470 <i>00757322</i> |

General Comments

Although some analytical issues are noted for the data group, unless otherwise noted in the detailed discussion, the data are considered valid for the purposes of the investigation. PDSs were deployed from April 7, 2015, through April 21, 2015.

Trip Blanks

There were no detections in the trip blank associated with these samples.

Equipment Blanks

Equipment blanks were not practicable due to sampler type.

Data Group Numbers (HCP PDS deployed June 1 – June 15, 2015, at Comal and San Marcos springs):

| | |
|-------------------------------|---------------------------|
| HCS 410 <i>00759936</i> | HSM 410 <i>00759941</i> |
| HCS 420 <i>00759937</i> | HSM 420 <i>00759942</i> |
| HCS 430 <i>00759935</i> | HSM 430 <i>00759943</i> |
| HCS 440 <i>00759938</i> | FDHSM 430 <i>00759944</i> |
| FDHCS 440 <i>00759939</i> | HSM 440 <i>00759945</i> |
| HCS 460 <i>00759940</i> | HSM 450 <i>00759946</i> |
| | HSM 460 <i>00759947</i> |
| Trip Blank 09 <i>00759949</i> | HSM 470 <i>00759948</i> |

General Comments

Although some analytical issues are noted for the data group, unless otherwise noted in the detailed discussion, the data are considered valid for the purposes of the investigation. PDSs were deployed from June 1, 2015, through June 15, 2015.

Trip Blanks

There were no detections in the trip blank associated with these samples.

Equipment Blanks

Equipment blanks were not practicable due to sampler type.

Data Group Numbers (HCP PDS deployed August 13 – August 27, 2015, at Comal and San Marcos springs):

| | |
|-------------------------------|---------------------------|
| HCS 410 <i>00767201</i> | HSM 410 <i>00767206</i> |
| HCS 420 <i>00767202</i> | HSM 420 <i>00767207</i> |
| HCS 430 <i>00767200</i> | HSM 430 <i>00767208</i> |
| HCS 440 <i>00767203</i> | FDHSM 430 <i>00767209</i> |
| FDHCS 440 <i>00767204</i> | HSM 440 <i>00767210</i> |
| HCS 460 <i>00767205</i> | HSM 450 <i>00767211</i> |
| | HSM 460 <i>00767212</i> |
| Trip Blank 10 <i>00767214</i> | HSM 470 <i>00767213</i> |

General Comments

Although some analytical issues are noted for the data group, unless otherwise noted in the detailed discussion, the data are considered valid for the purposes of the investigation. PDSs were deployed from August 13, 2015, through August 27, 2015.

Samples were analyzed on two different instruments, one of which had an elevated TPH background. The TPH detections in samples HCS 440, FDHCS 440, HCS 460, and HSM 440 may be biased high. An unused sampler analyzed on this instrument was found to have 0.65 µg of TPH.

Trip Blanks

There were no detections in the trip blank associated with these samples.

Equipment Blanks

Equipment blanks were not practicable due to sampler type.

Data Group Numbers (HCP PDS deployed October 6 – October 20, 2015, at Comal and San Marcos springs):

| | |
|-------------------------------|---------------------------|
| HCS 410 <i>00770685</i> | HSM 410 <i>00770691</i> |
| HCS 420 <i>00770686</i> | HSM 420 <i>00770692</i> |
| HCS 430 <i>00770687</i> | HSM 430 <i>00770693</i> |
| HCS 440 <i>00770688</i> | FDHSM 430 <i>00770694</i> |
| FDHCS 440 <i>00770689</i> | HSM 440 <i>00770695</i> |
| HCS 460 <i>00770690</i> | HSM 450 <i>00770696</i> |
| | HSM 460 <i>00770697</i> |
| Trip Blank 13 <i>00770699</i> | HSM 470 <i>00770698</i> |

General Comments

Although some analytical issues are noted for the data group, unless otherwise noted in the detailed discussion, the data are considered valid for the purposes of the investigation. PDSs were deployed from October 6, 2015, through October 20, 2015.

Trip Blanks

There were no detections in the trip blank associated with these samples.

Equipment Blanks

Equipment blanks were not practicable due to sampler type.

Data Group Numbers (HCP PDS deployed December 1 – December 15, 2015, at Comal and San Marcos springs):

| | |
|-------------------------------|---------------------------|
| HCS 410 <i>00770871</i> | HSM 410 <i>00770877</i> |
| HCS 420 <i>00770872</i> | HSM 420 <i>00770878</i> |
| HCS 430 <i>00770873</i> | HSM 430 <i>00770879</i> |
| HCS 440 <i>00770874</i> | FDHSM 430 <i>00770880</i> |
| FDHCS 440 <i>00770875</i> | HSM 440 <i>00770881</i> |
| HCS 460 <i>00770876</i> | HSM 450 <i>00770882</i> |
| | HSM 460 <i>00770883</i> |
| Trip Blank 16 <i>00770885</i> | HSM 470 <i>00770884</i> |

General Comments

Although some analytical issues are noted for the data group, unless otherwise noted in the detailed discussion, the data are considered valid for the purposes of the investigation. PDSs were deployed from December 1, 2015, through December 15, 2015. The PDS from HCS460 was not recovered.

Trip Blanks

There were no detections in the trip blank associated with these samples.

Equipment Blanks

Equipment blanks were not practicable due to sampler type.

Relative Percent Differences between Field Samples and their duplicates

Relative Percent Differences (RPD) values for parent samples and associated duplicate samples are provided in Table 2 below. In general, the RPD are less than 20% indicating parent and duplicate sample constituent concentrations are similar. Some differences observed in water samples may be the result of concentrations that vary naturally due to stream flow. Because the streams sampled are constantly flowing, constituent concentrations can change or fluctuate during the time period in which multiple sample bottles, which are all part of the same sample, are filled. Although sediment sample cores were collected in close proximity to one another, soil concentrations may vary spatially, which can cause differences between parent and duplicate sample constituent concentrations.

The differences observed do not show wide variations where a parent sample concentration exceeds a regulatory threshold or comparison value and a duplicate does not, or vice versa.

It should be noted that the RPDs between parent and duplicate field samples not only show differences between the parent and duplicate samples but also include differences inherent to laboratory procedures when the two separate samples are analyzed. Therefore, the laboratory RPDs contribute to the parent and field duplicate constituent concentration RPDs.

Table 2. Relative Percent Differences between Field Samples and their duplicates

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|----------------------|----------------|-------------------------|-------------|--------------|------------|-----------|------------|---------|
| Surface Water | | | | | | | | |
| HCS 120 | 3/16/2015 | Chloride | mg/L | 17 | | 17 | | 0.00% |
| | | Fluoride | mg/L | 0.23 | | 0.22 | | 4.44% |
| | | Nitrate | mg/L | 1.7 | | 1.7 | | 0.00% |
| | | Sulfate | mg/L | 29 | | 28 | | 3.51% |
| | | Calcium | mg/L | 87.2 | | 85.6 | | 1.85% |
| | | Magnesium | mg/L | 16.3 | B | 16.2 | B | 0.62% |
| | | Potassium | mg/L | 1.65 | | 1.67 | | 1.20% |
| | | Silicon | mg/L | 4.44 | | 4.44 | | 0.00% |
| | | Sodium | mg/L | 13.4 | | 13 | | 3.03% |
| | | Strontium | mg/L | 0.69 | | 0.668 | | 3.24% |
| | | Aluminum | mg/L | <0.00331 | | 0.00459 | J | 32.41% |
| | | Antimony | mg/L | <0.0000995 | | 0.000166 | J | 50.09% |
| | | Barium | mg/L | 0.0555 | | 0.0562 | | 1.25% |
| | | Chromium | mg/L | <0.000402 | | 0.000426 | J | 5.80% |
| | | Copper | mg/L | 0.000538 | J | 0.000783 | J | 37.09% |
| | | Iron | mg/L | 0.046 | J | 0.0555 | | 18.72% |
| | | Lead | mg/L | <0.0000898 | | 0.000117 | J | 26.31% |
| | | Manganese | mg/L | 0.00178 | | 0.00607 | | 109.30% |
| | | Nickel | mg/L | 0.00201 | | 0.00225 | | 11.27% |
| | | Selenium | mg/L | 0.000334 | J | 0.000414 | J | 21.39% |
| | | Zinc | mg/L | 0.0239 | | 0.0391 | | 48.25% |
| | | Bicarbonate | mg/L | 235 | | 235 | | 0.00% |
| | | Total Alkalinity | mg/L | 235 | | 235 | | 0.00% |
| | | Total dissolved solids | mg/L | 350 | | 335 | | 4.38% |
| | | pH | pH units | 7.36 | BV,BU | 7.41 | BV,BU | 0.68% |
| | | Total Kjeldahl Nitrogen | mg/L | 0.42 | J | 0.35 | J | 18.18% |
| | | DOC | mg/L | 2.9 | | 2.1 | | 32.00% |
| | | TOC | mg/L | 11 | | 10 | | 9.52% |
| | | E. coli | MPN/ 100 mL | 47 | | 51 | | 8.16% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|------------------------|-------------|--------------|------------|-----------|------------|---------|
| HCS 120 | 9/9/2015 | Caffeine | ng/L | <4 | | 16 | | 120.00% |
| | | Bromide | mg/L | 0.1 | | 0.096 | J | 4.08% |
| | | Chloride | mg/L | 17 | | 17 | | 0.00% |
| | | Fluoride | mg/L | 0.23 | | 0.22 | | 4.44% |
| | | Nitrate | mg/L | 1.8 | | 1.8 | | 0.00% |
| | | Sulfate | mg/L | 25 | | 25 | | 0.00% |
| | | Calcium | mg/L | 82.2 | B | 81.3 | B | 1.10% |
| | | Magnesium | mg/L | 16.3 | B | 15.4 | B | 5.68% |
| | | Potassium | mg/L | 2.3 | B | 2.04 | B | 11.98% |
| | | Silicon | mg/L | 5.55 | | 5.2 | | 6.51% |
| | | Sodium | mg/L | 12.7 | B | 12.2 | B | 4.02% |
| | | Strontium | mg/L | 0.67 | | 0.665 | | 0.75% |
| | | Aluminum | mg/L | 0.00699 | J | 0.00954 | J | 30.85% |
| | | Arsenic | mg/L | 0.000388 | J | <0.000386 | | 0.52% |
| | | Barium | mg/L | 0.0549 | | 0.0537 | | 2.21% |
| | | Chromium | mg/L | 0.000524 | J | 0.000482 | J | 8.35% |
| | | Copper | mg/L | 0.00504 | | 0.00078 | J | 146.39% |
| | | Iron | mg/L | 0.0364 | J | 0.0406 | J | 10.91% |
| | | Manganese | mg/L | 0.00082 | J | 0.00064 | J | 24.66% |
| | | Nickel | mg/L | 0.00182 | | 0.00187 | | 2.71% |
| | | Selenium | mg/L | 0.000521 | J | 0.000932 | J | 56.57% |
| | | Thallium | mg/L | 0.000121 | J | <0.000101 | | 18.02% |
| | | Zinc | mg/L | 0.00984 | | 0.00773 | | 24.02% |
| | | Bicarbonate | mg/L | 221 | | 212 | | 4.16% |
| | | Total Alkalinity | mg/L | 221 | | 212 | | 4.16% |
| | | Total dissolved solids | mg/L | 355 | | 340 | | 4.32% |
| | | pH | pH units | 7.23 | BV,BU | 7.25 | BV,BU | 0.28% |
| | | DOC | mg/L | 12 | | 11 | | 8.70% |
| | | TOC | mg/L | 8 | | 5.7 | | 33.58% |
| | | E. coli | MPN/ 100 mL | 67 | | 54 | | 21.49% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|-------------------------|-------------|--------------|------------|-----------|------------|---------|
| HSM 110 | 3/25/2015 | Caffeine | ng/L | 16 | | 10 | | 46.15% |
| | | Bromide | mg/L | 0.16 | | 0.15 | | 6.45% |
| | | Chloride | mg/L | 24 | | 24 | | 0.00% |
| | | Fluoride | mg/L | 0.23 | | 0.23 | | 0.00% |
| | | Nitrate | mg/L | 0.36 | | 0.36 | | 0.00% |
| | | Sulfate | mg/L | 33 | | 33 | | 0.00% |
| | | Phosphorous, total | mg/L | <0.02 | | 0.022 | J | 9.52% |
| | | Calcium | mg/L | 93.9 | | 87.4 | | 7.17% |
| | | Magnesium | mg/L | 18 | B | 20.1 | B | 11.02% |
| | | Potassium | mg/L | 1.31 | | 1.75 | | 28.76% |
| | | Silicon | mg/L | 3.63 | | 3.59 | | 1.11% |
| | | Sodium | mg/L | 12.7 | | 15.3 | | 18.57% |
| | | Strontium | mg/L | 0.568 | | 0.688 | | 19.11% |
| | | Barium | mg/L | 0.0409 | | 0.0429 | | 4.77% |
| | | Copper | mg/L | 0.000788 | J | 0.000673 | J | 15.74% |
| | | Iron | mg/L | 0.0411 | J | 0.0485 | J | 16.52% |
| | | Manganese | mg/L | 0.0222 | | 0.022 | | 0.90% |
| | | Nickel | mg/L | 0.00193 | | 0.00203 | | 5.05% |
| | | Selenium | mg/L | 0.00031 | J | <0.000168 | | 59.41% |
| | | Zinc | mg/L | 0.00287 | J | <0.000479 | | 142.79% |
| | | Bicarbonate | mg/L | 244 | | 253 | | 3.62% |
| | | Total Alkalinity | mg/L | 244 | | 253 | | 3.62% |
| | | Total dissolved solids | mg/L | 365 | | 385 | | 5.33% |
| | | Total suspended solids | mg/L | 5.9 | | 9 | | 41.61% |
| | | pH | pH units | 7.29 | BV,BU | 7.57 | BV,BU | 3.77% |
| | | Total Kjeldahl Nitrogen | mg/L | 0.42 | J | 0.28 | J | 40.00% |
| | | DOC | mg/L | 1.9 | | 3.1 | | 48.00% |
| | | TOC | mg/L | 1.4 | | 4.6 | | 106.67% |
| | | E. coli | MPN/ 100 mL | 1 | | 9 | | 160.00% |
| | | E. coli | MPN/ 100 mL | 45 | | 75 | | 50.00% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|-------------------------|-------------|--------------|------------|-----------|------------|--------|
| HSM110 | 9/17/2015 | Caffeine | ng/L | 19 | | 13 | | 37.50% |
| | | Bromide | mg/L | 0.17 | | 0.18 | | 5.71% |
| | | Chloride | mg/L | 29 | | 29 | | 0.00% |
| | | Fluoride | mg/L | 0.28 | | 0.27 | | 3.64% |
| | | Nitrate | mg/L | 0.56 | | 0.59 | | 5.22% |
| | | Sulfate | mg/L | 36 | | 36 | | 0.00% |
| | | Phosphorous, total | mg/L | 0.038 | J | 0.039 | J | 2.60% |
| | | Calcium | mg/L | 94.8 | | 91.2 | | 3.87% |
| | | Magnesium | mg/L | 20.1 | B | 19.4 | B | 3.54% |
| | | Potassium | mg/L | 2.18 | | 2.08 | | 4.69% |
| | | Silicon | mg/L | 5.53 | | 5.33 | | 3.68% |
| | | Sodium | mg/L | 17.8 | | 17.1 | | 4.01% |
| | | Strontium | mg/L | 0.756 | | 0.725 | | 4.19% |
| | | Barium | mg/L | 0.0407 | | 0.0394 | | 3.25% |
| | | Copper | mg/L | 0.00175 | J | 0.00216 | J | 20.97% |
| | | Manganese | mg/L | 0.283 | | 0.282 | | 0.35% |
| | | Nickel | mg/L | 0.00176 | J | 0.00168 | J | 4.65% |
| | | Bicarbonate | mg/L | 266 | | 271 | | 1.86% |
| | | Total Alkalinity | mg/L | 266 | | 271 | | 1.86% |
| | | Total dissolved solids | mg/L | 420 | | 400 | | 4.88% |
| | | Total suspended solids | mg/L | 4.5 | | 4.6 | | 2.20% |
| | | pH | pH units | 7.03 | BV,BU | 7.1 | BV,BU | 0.99% |
| | | Total Kjeldahl Nitrogen | mg/L | 0.28 | J | 0.35 | J | 22.22% |
| | | DOC | mg/L | 1.1 | | 1.2 | | 8.70% |
| | | TOC | mg/L | 1.4 | | 1.4 | | 0.00% |
| | | E. coli | MPN/ 100 mL | 45 | | 75 | | 50.00% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|----------------------|----------------|-------------------------|------------|--------------|------------|-----------|------------|---------|
| Storm Water | | | | | | | | |
| HCS 260 Trail | 1/23/2015 | Caffeine | ng/L | 110 | | 67 | | 48.59% |
| | | Bromide | mg/L | 0.077 | J | 0.073 | J | 5.33% |
| | | Chloride | mg/L | 12 | | 12 | | 0.00% |
| | | Fluoride | mg/L | 0.2 | | 0.19 | | 5.13% |
| | | Nitrate | mg/L | 1.2 | | 1.2 | | 0.00% |
| | | Sulfate | mg/L | 25 | | 25 | | 0.00% |
| | | Phosphorous, total | mg/L | 0.11 | | 0.1 | | 9.52% |
| | | Calcium | mg/L | 57.6 | | 56 | | 2.82% |
| | | Magnesium | mg/L | 8.87 | B | 8.74 | B | 1.48% |
| | | Potassium | mg/L | 3.91 | | 3.82 | | 2.33% |
| | | Silicon | mg/L | 3.52 | | 3.68 | | 4.44% |
| | | Sodium | mg/L | 10.2 | | 10 | | 1.98% |
| | | Strontium | mg/L | 0.409 | | 0.401 | | 1.98% |
| | | Aluminum | mg/L | 0.0224 | J | 0.188 | | 157.41% |
| | | Antimony | mg/L | 0.000118 | J | 0.000104 | J | 12.61% |
| | | Arsenic | mg/L | 0.0014 | | 0.00134 | | 4.38% |
| | | Barium | mg/L | 0.0471 | | 0.0466 | | 1.07% |
| | | Chromium | mg/L | <0.000402 | | 0.000525 | J | 26.54% |
| | | Copper | mg/L | 0.00142 | | 0.00141 | | 0.71% |
| | | Iron | mg/L | 0.0678 | | 0.129 | | 62.20% |
| | | Lead | mg/L | <0.0000898 | | 0.00011 | J | 20.22% |
| | | Manganese | mg/L | 0.00767 | | 0.00662 | | 14.70% |
| | | Nickel | mg/L | 0.00191 | | 0.00171 | | 11.05% |
| | | Selenium | mg/L | 0.000207 | J | 0.00018 | J | 13.95% |
| | | Zinc | mg/L | 0.00978 | | 0.0135 | | 31.96% |
| | | Bicarbonate | mg/L | 155 | | 157 | | 1.28% |
| | | Total Alkalinity | mg/L | 155 | | 157 | | 1.28% |
| | | Total dissolved solids | mg/L | 280 | | 400 | | 35.29% |
| | | Total suspended solids | mg/L | 182 | | 89 | | 68.63% |
| | | pH | pH units | 7.75 | BV,BU | 7.71 | BV,BU | 0.52% |
| | | Total Kjeldahl Nitrogen | mg/L | 1 | | 0.77 | | 25.99% |
| | | DOC | mg/L | 4.7 | | 5.6 | | 17.48% |
| | | TOC | mg/L | 11 | | 9.5 | | 14.63% |
| | | E. coli | MPN/100 mL | 2300 | H | 3400 | H | 38.60% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|-------------------------|------------|--------------|------------|-----------|------------|---------|
| HCS 260 Trail | 10/23/2015 | Aluminum | mg/L | 0.00338 | J | 0.00574 | J | 51.75% |
| | | Barium | mg/L | 0.0535 | | 0.0529 | | 1.13% |
| | | Bicarbonate | mg/L | 219 | | 220 | | 0.46% |
| | | Bromide | mg/L | 0.099 | J | 0.1 | | 1.01% |
| | | Caffeine | ng/L | 80 | | 56 | | 35.29% |
| | | Calcium | mg/L | 79.4 | B | 78 | B | 1.78% |
| | | Chloride | mg/L | 17 | | 17 | | 0.00% |
| | | Copper | mg/L | 0.00056 | J | 0.000267 | J | 70.86% |
| | | E. coli | MPN/100 mL | 20000 | H | 24000 | H | 18.18% |
| | | Fluoride | mg/L | 0.19 | | 0.2 | | 5.13% |
| | | Iron | mg/L | 0.0502 | | 0.0559 | | 10.74% |
| | | Magnesium | mg/L | 16.6 | | 16.5 | | 0.60% |
| | | Manganese | mg/L | 0.00299 | | <0.000139 | | 182.23% |
| | | Nickel | mg/L | 0.00152 | | 0.00152 | | 0.00% |
| | | Nitrate | mg/L | 1.7 | BU | 1.7 | BU | 0.00% |
| | | pH | pH units | 7.66 | BV,BU | 7.68 | BV,BU | 0.26% |
| | | Phosphorous, total | mg/L | 0.04 | J | 0.039 | J | 2.53% |
| | | Potassium | mg/L | 1.87 | | 1.83 | | 2.16% |
| | | Selenium | mg/L | 0.00027 | J | 0.000327 | J | 19.10% |
| | | Silicon | mg/L | 5.9 | | 5.83 | | 1.19% |
| | | Sodium | mg/L | 12.7 | | 12.7 | | 0.00% |
| | | Strontium | mg/L | 0.651 | | 0.647 | | 0.62% |
| | | Sulfate | mg/L | 30 | | 30 | | 0.00% |
| | | Total Alkalinity | mg/L | 219 | | 220 | | 0.46% |
| | | Total dissolved solids | mg/L | 275 | | 285 | | 3.57% |
| | | Total Kjeldahl Nitrogen | mg/L | 0.84 | | 0.56 | | 40.00% |
| | | Total suspended solids | mg/L | 5.7 | | 2.8 | | 68.24% |
| | | Zinc | mg/L | 0.00363 | J | 0.014 | | 117.64% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|-------------------------|------------|--------------|------------|-----------|------------|---------|
| HCS 270 Trail | 1/23/2015 | Caffeine | ng/L | 100 | | 65 | | 42.42% |
| | | Bromide | mg/L | 0.082 | J | 0.089 | J | 8.19% |
| | | Chloride | mg/L | 12 | | 12 | | 0.00% |
| | | Fluoride | mg/L | 0.2 | | 0.18 | | 10.53% |
| | | Nitrate | mg/L | 1.2 | | 1.2 | | 0.00% |
| | | Sulfate | mg/L | 24 | | 23 | | 4.26% |
| | | Phosphorous, total | mg/L | 0.096 | | 0.12 | | 22.22% |
| | | Calcium | mg/L | 57.8 | | 55.2 | | 4.60% |
| | | Magnesium | mg/L | 8.83 | B | 8.93 | B | 1.13% |
| | | Potassium | mg/L | 4.06 | | 3.86 | | 5.05% |
| | | Silicon | mg/L | 3.48 | | 3.64 | | 4.49% |
| | | Sodium | mg/L | 10.1 | | 9.77 | | 3.32% |
| | | Strontium | mg/L | 0.414 | | 0.399 | | 3.69% |
| | | Aluminum | mg/L | <0.00331 | | 0.0847 | | 184.96% |
| | | Antimony | mg/L | 0.00011 | J | 0.000108 | J | 1.83% |
| | | Arsenic | mg/L | 0.00109 | | 0.00121 | | 10.43% |
| | | Barium | mg/L | 0.0498 | | 0.0467 | | 6.42% |
| | | Chromium | mg/L | <0.000402 | | 0.000417 | J | 3.66% |
| | | Copper | mg/L | 0.00142 | | 0.00186 | | 26.83% |
| | | Iron | mg/L | 0.0535 | | 0.0808 | | 40.66% |
| | | Manganese | mg/L | 0.00408 | | 0.00446 | | 8.90% |
| | | Nickel | mg/L | 0.0018 | | 0.00204 | | 12.50% |
| | | Selenium | mg/L | 0.000171 | J | 0.000258 | J | 40.56% |
| | | Zinc | mg/L | 0.00605 | | 0.0105 | | 53.78% |
| | | Bicarbonate | mg/L | 152 | | 154 | | 1.31% |
| | | Total Alkalinity | mg/L | 152 | | 154 | | 1.31% |
| | | Total dissolved solids | mg/L | 265 | | 275 | | 3.70% |
| | | Total suspended solids | mg/L | 114 | | 114 | | 0.00% |
| | | pH | pH units | 7.78 | BV,BU | 7.77 | BV,BU | 0.13% |
| | | Total Kjeldahl Nitrogen | mg/L | <0.28 | | 0.56 | | 66.67% |
| | | DOC | mg/L | 7 | | 7.8 | | 10.81% |
| | | TOC | mg/L | 7.4 | | 8.2 | | 10.26% |
| | | E. coli | MPN/100 mL | 4400 | H | 2500 | H | 55.07% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|-------------------------|------------|--------------|------------|-----------|------------|---------|
| HSM 270 Trail | 10/23/2015 | Aluminum | mg/L | 0.00724 | | 0.00581 | J | 21.92% |
| | | Barium | mg/L | 0.0512 | | 0.049 | | 4.39% |
| | | Bicarbonate | mg/L | 206 | | 206 | | 0.00% |
| | | Bromide | mg/L | 0.088 | | 0.089 | J | 1.13% |
| | | Caffeine | ng/L | 450 | | 130 | | 110.34% |
| | | Calcium | mg/L | 73.8 | | 71.5 | B | 3.17% |
| | | Chloride | mg/L | 18 | | 18 | | 0.00% |
| | | Copper | mg/L | 0.000721 | | 0.000496 | J | 36.98% |
| | | E. coli | MPN/100 mL | 20000 | | 17000 | H | 16.22% |
| | | Fluoride | mg/L | 0.2 | | 0.19 | | 5.13% |
| | | Iron | mg/L | 0.0553 | | 0.0502 | | 9.67% |
| | | Magnesium | mg/L | 15.3 | | 14.8 | | 3.32% |
| | | Manganese | mg/L | 0.000998 | | 0.000298 | J | 108.02% |
| | | Nickel | mg/L | 0.00145 | | 0.00139 | | 4.23% |
| | | Nitrate | mg/L | 1.6 | | 1.6 | BU | 0.00% |
| | | pH | pH units | 7.68 | | 7.69 | BV,BU | 0.13% |
| | | Phosphorous, total | mg/L | 0.05 | | 0.038 | J | 27.27% |
| | | Potassium | mg/L | 2.04 | | 1.99 | | 2.48% |
| | | Selenium | mg/L | 0.00035 | | 0.000318 | J | 9.58% |
| | | Silicon | mg/L | 5.52 | | 5.29 | | 4.26% |
| | | Sodium | mg/L | 13.1 | | 12.8 | | 2.32% |
| | | Strontium | mg/L | 0.605 | | 0.593 | | 2.00% |
| | | Sulfate | mg/L | 28 | | 28 | | 0.00% |
| | | Total Alkalinity | mg/L | 206 | | 206 | | 0.00% |
| | | Total dissolved solids | mg/L | 315 | | 295 | | 6.56% |
| | | Total Kjeldahl Nitrogen | mg/L | 0.98 | | 0.84 | | 15.38% |
| | | Total suspended solids | mg/L | 2.4 | | 4.4 | | 58.82% |
| | | Zinc | mg/L | 0.00559 | | 0.0101 | | 57.49% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|-------------------------|------------|--------------|------------|-----------|------------|---------|
| HSM 210 Trail | 5/6/2015 | Caffeine | ng/L | 13 | | 14 | | 7.41% |
| | | Bromide | mg/L | 0.14 | | 0.16 | | 13.33% |
| | | Chloride | mg/L | 26 | | 25 | | 3.92% |
| | | Fluoride | mg/L | 0.22 | | 0.21 | | 4.65% |
| | | Nitrate | mg/L | 0.41 | | 0.41 | | 0.00% |
| | | Sulfate | mg/L | 35 | | 35 | | 0.00% |
| | | Phosphorous, total | mg/L | 0.054 | | <0.02 | | 91.89% |
| | | Calcium | mg/L | 90.1 | | 91.7 | | 1.76% |
| | | Magnesium | mg/L | 20.3 | | 20.1 | B | 0.99% |
| | | Potassium | mg/L | 2.29 | | 2.06 | | 10.57% |
| | | Silicon | mg/L | 4.45 | | 4.41 | | 0.90% |
| | | Sodium | mg/L | 16.5 | | 16.5 | | 0.00% |
| | | Strontium | mg/L | 0.723 | | 0.715 | | 1.11% |
| | | Aluminum | mg/L | 0.00574 | J | 0.00347 | J | 49.29% |
| | | Arsenic | mg/L | <0.000386 | | 0.00124 | B | 105.04% |
| | | Barium | mg/L | 0.0362 | | 0.0397 | | 9.22% |
| | | Copper | mg/L | 0.000804 | J | 0.000945 | J | 16.12% |
| | | Iron | mg/L | 0.0724 | | 0.0708 | | 2.23% |
| | | Manganese | mg/L | 0.0184 | | 0.0203 | | 9.82% |
| | | Nickel | mg/L | 0.00239 | | 0.00207 | | 14.35% |
| | | Selenium | mg/L | 0.000394 | J | 0.000263 | J | 39.88% |
| | | Zinc | mg/L | 0.0106 | | 0.011 | | 3.70% |
| | | p-Isopropyltoluene | ug/L | 0.18 | J | <0.16 | | 11.76% |
| | | Bicarbonate | mg/L | 240 | | 256 | | 6.45% |
| | | Total Alkalinity | mg/L | 240 | | 256 | | 6.45% |
| | | Total dissolved solids | mg/L | 405 | | 340 | | 17.45% |
| | | Total suspended solids | mg/L | 2.3 | | 3.8 | | 49.18% |
| | | pH | pH units | 7.48 | BV,BU | 7.47 | BV,BU | 0.13% |
| | | Total Kjeldahl Nitrogen | mg/L | 0.56 | | 0.56 | | 0.00% |
| | | DOC | mg/L | 2.4 | J | 3 | | 22.22% |
| | | TOC | mg/L | 0.46 | J | 3 | | 146.82% |
| | | E. coli | MPN/100 mL | 270 | H | 330 | H | 20.00% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|-------------------------|------------|--------------|------------|-----------|------------|--------|
| HSM 210 Trail | 10/23/2015 | Aluminum | mg/L | 0.0057 | J | 0.00569 | J | 0.18% |
| | | Barium | mg/L | 0.0341 | | 0.0347 | | 1.74% |
| | | Bicarbonate | mg/L | 228 | | 233 | | 2.17% |
| | | Bromide | mg/L | 0.17 | | 0.17 | | 0.00% |
| | | Caffeine | ng/L | 2.4 | | 5.7 | | 81.48% |
| | | Calcium | mg/L | 78.6 | B | 80.4 | B | 2.26% |
| | | Chloride | mg/L | 27 | | 27 | | 0.00% |
| | | Copper | mg/L | <0.000140 | | 0.000179 | J | 24.45% |
| | | E. coli | MPN/100 mL | 450 | H | 180 | H | 85.71% |
| | | Fluoride | mg/L | 0.21 | | 0.2 | | 4.88% |
| | | Iron | mg/L | 0.0519 | | 0.0609 | | 15.96% |
| | | Magnesium | mg/L | 20 | | 20.7 | | 3.44% |
| | | Manganese | mg/L | 0.0865 | | 0.0846 | | 2.22% |
| | | Nickel | mg/L | 0.00157 | | 0.00165 | | 4.97% |
| | | Nitrate | mg/L | 0.61 | BU | 0.62 | BU | 1.63% |
| | | pH | pH units | 7.43 | BV,BU | 7.39 | BV,BU | 0.54% |
| | | Phosphorous, total | mg/L | 0.026 | J | 0.036 | J | 32.26% |
| | | Potassium | mg/L | 2.01 | | 2.08 | | 3.42% |
| | | Selenium | mg/L | 0.000261 | J | <0.000168 | | 43.36% |
| | | Silicon | mg/L | 5.31 | | 5.45 | | 2.60% |
| | | Sodium | mg/L | 15.4 | | 15.8 | | 2.56% |
| | | Strontium | mg/L | 0.658 | | 0.675 | | 2.55% |
| | | Sulfate | mg/L | 34 | | 34 | | 0.00% |
| | | Total Alkalinity | mg/L | 228 | | 233 | | 2.17% |
| | | Total dissolved solids | mg/L | 345 | | 370 | | 6.99% |
| | | Total Kjeldahl Nitrogen | mg/L | 0.56 | | 0.42 | J | 28.57% |
| | | Total suspended solids | mg/L | 2 | | 4.7 | | 80.60% |
| | | Zinc | mg/L | 0.0193 | | 0.00791 | | 83.72% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|-------------------------|------------|--------------|------------|-----------|------------|--------|
| HSM 230 Trail | 5/6/2015 | Caffeine | ng/L | 290 | | 340 | | 15.87% |
| | | Bromide | mg/L | 0.076 | J | 0.079 | J | 3.87% |
| | | Chloride | mg/L | 17 | | 17 | | 0.00% |
| | | Fluoride | mg/L | 0.16 | | 0.16 | | 0.00% |
| | | Nitrate | mg/L | 1.5 | | 1.5 | | 0.00% |
| | | Sulfate | mg/L | 27 | | 27 | | 0.00% |
| | | Phosphorous, total | mg/L | 0.045 | J | 0.029 | J | 43.24% |
| | | Calcium | mg/L | 83.8 | | 86.8 | | 3.52% |
| | | Magnesium | mg/L | 12.1 | | 12.5 | B | 3.25% |
| | | Potassium | mg/L | 2.52 | | 2.05 | | 20.57% |
| | | Silicon | mg/L | 4.01 | | 4.11 | | 2.46% |
| | | Sodium | mg/L | 12.4 | | 12.9 | | 3.95% |
| | | Strontium | mg/L | 0.404 | | 0.411 | | 1.72% |
| | | Aluminum | mg/L | 0.0119 | J | 0.0117 | J | 1.69% |
| | | Antimony | mg/L | 0.000248 | J | 0.000244 | J | 1.63% |
| | | Arsenic | mg/L | <0.000386 | | 0.000733 | B,J | 62.02% |
| | | Barium | mg/L | 0.0359 | | 0.0387 | | 7.51% |
| | | Chromium | mg/L | 0.00056 | J | 0.000634 | J | 12.40% |
| | | Copper | mg/L | 0.00229 | | 0.00275 | | 18.25% |
| | | Iron | mg/L | 0.0929 | | 0.0758 | | 20.27% |
| | | Lead | mg/L | 0.000122 | J | 0.000124 | J | 1.63% |
| | | Manganese | mg/L | 0.00719 | | 0.00682 | | 5.28% |
| | | Nickel | mg/L | 0.00234 | | 0.00235 | | 0.43% |
| | | Selenium | mg/L | 0.000308 | J | 0.000317 | J | 2.88% |
| | | Zinc | mg/L | 0.0308 | | 0.0278 | | 10.24% |
| | | 4,4-DDD | ug/L | 0.038 | J | <0.027 | | 33.85% |
| | | Bicarbonate | mg/L | 199 | | 198 | | 0.50% |
| | | Total Alkalinity | mg/L | 199 | | 198 | | 0.50% |
| | | Total dissolved solids | mg/L | 325 | | 200 | | 47.62% |
| | | Total suspended solids | mg/L | 2.4 | | 3.3 | | 31.58% |
| | | pH | pH units | 7.35 | BV,BU | 7.48 | BV,BU | 1.75% |
| | | Total Kjeldahl Nitrogen | mg/L | 0.56 | | 0.56 | | 0.00% |
| | | DOC | mg/L | 2.6 | | 6.6 | | 86.96% |
| | | TOC | mg/L | 3.2 | | 4 | | 22.22% |
| | | E. coli | MPN/100 mL | 5200 | H | 4600 | H | 12.24% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|-------------------------|------------|--------------|------------|-----------|------------|---------|
| HSM 230 Trail | 10/23/2015 | Aluminum | mg/L | 0.00775 | J | 0.00881 | J | 12.80% |
| | | Antimony | mg/L | 0.000173 | J | 0.000143 | J | 18.99% |
| | | Barium | mg/L | 0.0357 | | 0.0369 | | 3.31% |
| | | Bicarbonate | mg/L | 228 | | 222 | | 2.67% |
| | | Bromide | mg/L | 0.1 | | 0.11 | | 9.52% |
| | | Caffeine | ng/L | 470 | | 450 | | 4.35% |
| | | Calcium | mg/L | 83.2 | B | 80.9 | B | 2.80% |
| | | Chloride | mg/L | 19 | | 19 | | 0.00% |
| | | Copper | mg/L | 0.00118 | | 0.00151 | | 24.54% |
| | | E. coli | MPN/100 mL | 73000 | H | 65000 | H | 11.59% |
| | | Fluoride | mg/L | 0.18 | | 0.2 | | 10.53% |
| | | Iron | mg/L | 0.0568 | | 0.0529 | | 7.11% |
| | | Magnesium | mg/L | 15.1 | | 15.1 | | 0.00% |
| | | Manganese | mg/L | 0.00236 | | 0.00354 | | 40.00% |
| | | Nickel | mg/L | 0.00192 | | 0.00188 | | 2.11% |
| | | Nitrate | mg/L | 1.5 | BU | 1.6 | BU | 6.45% |
| | | pH | pH units | 7.36 | BV,BU | 7.1 | BV,BU | 3.60% |
| | | Phosphorous, total | mg/L | 0.039 | J | 0.058 | | 39.18% |
| | | Potassium | mg/L | 2.25 | | 2.22 | | 1.34% |
| | | Selenium | mg/L | 0.000374 | J | 0.00027 | J | 32.30% |
| | | Silicon | mg/L | 5.22 | | 5.18 | | 0.77% |
| | | Sodium | mg/L | 12.3 | | 12.2 | | 0.82% |
| | | Strontium | mg/L | 0.471 | | 0.465 | | 1.28% |
| | | Sulfate | mg/L | 30 | | 30 | | 0.00% |
| | | Total Alkalinity | mg/L | 228 | | 222 | | 2.67% |
| | | Total dissolved solids | mg/L | 290 | | 310 | | 6.67% |
| | | Total Kjeldahl Nitrogen | mg/L | 0.84 | | 0.84 | | 0.00% |
| | | Total suspended solids | mg/L | <0.83 | | 4 | | 131.26% |
| | | Zinc | mg/L | 0.0401 | | 0.00928 | | 124.83% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|-------------------------|------------|--------------|------------|-----------|------------|--------|
| HSM 231 Trail | 5/6/2015 | Caffeine | ng/L | 25 | | 25 | | 0.00% |
| | | Bromide | mg/L | 0.094 | J | 0.097 | J | 3.14% |
| | | Chloride | mg/L | 19 | | 19 | | 0.00% |
| | | Fluoride | mg/L | 0.19 | | 0.19 | | 0.00% |
| | | Nitrate | mg/L | 1.2 | | 1.2 | | 0.00% |
| | | Sulfate | mg/L | 28 | | 28 | | 0.00% |
| | | Phosphorous, total | mg/L | <0.02 | | 0.042 | J | 70.97% |
| | | Calcium | mg/L | 98.7 | | 97.2 | | 1.53% |
| | | Magnesium | mg/L | 18.4 | | 18.4 | B | 0.00% |
| | | Potassium | mg/L | 1.53 | | 1.48 | | 3.32% |
| | | Silicon | mg/L | 4.64 | | 4.65 | | 0.22% |
| | | Sodium | mg/L | 14 | | 13.5 | | 3.64% |
| | | Strontium | mg/L | 0.608 | | 0.597 | | 1.83% |
| | | Aluminum | mg/L | 0.00844 | J | <0.00331 | | 87.32% |
| | | Arsenic | mg/L | <0.000386 | | 0.000941 | B,J | 83.65% |
| | | Barium | mg/L | 0.0382 | | 0.043 | | 11.82% |
| | | Chromium | mg/L | 0.000402 | J | <0.000402 | | 0.00% |
| | | Copper | mg/L | 0.000596 | J | 0.00105 | | 55.16% |
| | | Iron | mg/L | 0.0695 | | 0.086 | | 21.22% |
| | | Lead | mg/L | <0.0000898 | | 0.000146 | J | 47.67% |
| | | Manganese | mg/L | 0.00172 | | 0.00175 | | 1.73% |
| | | Nickel | mg/L | 0.00256 | | 0.00225 | | 12.89% |
| | | Selenium | mg/L | 0.000386 | J | 0.000319 | J | 19.01% |
| | | Zinc | mg/L | 0.018 | | 0.0361 | | 66.91% |
| | | Bicarbonate | mg/L | 252 | | 255 | | 1.18% |
| | | Total Alkalinity | mg/L | 252 | | 255 | | 1.18% |
| | | Total dissolved solids | mg/L | 385 | | 255 | | 40.63% |
| | | Total suspended solids | mg/L | 1.2 | | <0.83 | | 36.45% |
| | | pH | pH units | 7.48 | BV,BU | 7.52 | BV,BU | 0.53% |
| | | Total Kjeldahl Nitrogen | mg/L | 0.84 | | 0.56 | | 40.00% |
| | | DOC | mg/L | 5 | | 3 | | 50.00% |
| | | TOC | mg/L | 4 | | 2.5 | | 46.15% |
| | | E. coli | MPN/100 mL | 280 | H | 330 | H | 16.39% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|-------------------------|------------|--------------|------------|-----------|------------|---------|
| HSM 231 Trail | 10/23/2015 | Aluminum | mg/L | 0.00682 | J | 0.00514 | J | 28.09% |
| | | Barium | mg/L | 0.0367 | | 0.0374 | | 1.89% |
| | | Bicarbonate | mg/L | 245 | | 245 | | 0.00% |
| | | Bromide | mg/L | 0.1 | | 0.1 | | 0.00% |
| | | Caffeine | ng/L | 19 | | 15 | | 23.53% |
| | | Calcium | mg/L | 87.7 | B | 86.2 | B | 1.73% |
| | | Chloride | mg/L | 18 | | 18 | | 0.00% |
| | | Copper | mg/L | 0.000193 | J | <0.000140 | | 31.83% |
| | | E. coli | MPN/100 mL | 1700 | H | 1700 | H | 0.00% |
| | | Fluoride | mg/L | 0.18 | | 0.18 | | 0.00% |
| | | Iron | mg/L | 0.055 | | 0.0707 | | 24.98% |
| | | Magnesium | mg/L | 18.6 | | 18.2 | | 2.17% |
| | | Manganese | mg/L | 0.000472 | J | <0.000139 | | 109.00% |
| | | Nickel | mg/L | 0.0017 | | 0.00183 | | 7.37% |
| | | Nitrate | mg/L | 1.3 | BU | 1.2 | BU | 8.00% |
| | | pH | pH units | 7.48 | BV,BU | 7.28 | BV,BU | 2.71% |
| | | Phosphorous, total | mg/L | 0.034 | J | 0.031 | J | 9.23% |
| | | Potassium | mg/L | 1.51 | | 1.55 | | 2.61% |
| | | Selenium | mg/L | 0.000309 | J | 0.000292 | J | 5.66% |
| | | Silicon | mg/L | 5.74 | | 5.62 | | 2.11% |
| | | Sodium | mg/L | 11.4 | | 11.4 | | 0.00% |
| | | Strontium | mg/L | 0.529 | | 0.524 | | 0.95% |
| | | Sulfate | mg/L | 25 | | 25 | | 0.00% |
| | | TOC | mg/L | <0.24 | | 7.4 | | 187.43% |
| | | Total Alkalinity | mg/L | 245 | | 245 | | 0.00% |
| | | Total dissolved solids | mg/L | 235 | | 295 | | 22.64% |
| | | Total Kjeldahl Nitrogen | mg/L | 0.84 | | 0.28 | J | 100.00% |
| | | Total suspended solids | mg/L | <0.83 | | 1.1 | | 27.98% |
| | | Zinc | mg/L | 0.0136 | | 0.0502 | | 114.73% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|------------------------|----------|--------------|------------|-----------|------------|--------|
| Sediment | | | | | | | | |
| HCS 360 | 6/4/2015 | Chloride | mg/kg | 22 | | 22 | | 0.00% |
| | | Fluoride | mg/kg | 3.1 | | 3.2 | | 3.17% |
| | | Sulfate | mg/kg | 49 | | 68 | | 32.48% |
| | | Calcium | mg/kg | 146000 | | 165000 | | 12.22% |
| | | Magnesium | mg/kg | 3060 | | 3560 | | 15.11% |
| | | Potassium | mg/kg | 2300 | | 2580 | | 11.48% |
| | | Silicon | mg/kg | 1410 | | 1690 | | 18.06% |
| | | Sodium | mg/kg | 207 | | 193 | | 7.00% |
| | | Strontium | mg/kg | 199 | | 222 | | 10.93% |
| | | Aluminum | mg/kg | 10800 | | 11900 | | 9.69% |
| | | Arsenic | mg/kg | 3.83 | | 4.15 | | 8.02% |
| | | Barium | mg/kg | 68.9 | | 72.9 | | 5.64% |
| | | Chromium | mg/kg | 16.1 | | 18 | | 11.14% |
| | | Copper | mg/kg | 14.3 | | 16 | | 11.22% |
| | | Iron | mg/kg | 9750 | | 10700 | | 9.29% |
| | | Lead | mg/kg | 23.9 | | 26.2 | | 9.18% |
| | | Manganese | mg/kg | 216 | | 227 | | 4.97% |
| | | Nickel | mg/kg | 17 | | 18.4 | | 7.91% |
| | | Zinc | mg/kg | 142 | | 160 | | 11.92% |
| | | pH | pH units | 7.35 | | 7.41 | | 0.81% |
| | | TOC | mg/kg | 53000 | | 50000 | | 5.83% |
| | | Bicarbonate | mg/kg | 580 | | 590 | | 1.71% |
| | | Total Alkalinity | mg/kg | 580 | | 590 | | 1.71% |
| | | Total dissolved solids | mg/kg | 3460 | | 4320 | | 22.11% |
| | | Phosphorous, total | mg/kg | 5.6 | | 5.8 | | 3.51% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------|----------------|------------------------|----------|--------------|------------|-----------|------------|--------|
| HSM370 | 6/5/2015 | Chloride | mg/kg | 17 | J | 16 | J | 6.06% |
| | | Fluoride | mg/kg | 1.7 | | 2 | | 16.22% |
| | | Nitrate | mg/kg | 0.69 | J | 0.61 | J | 12.31% |
| | | Sulfate | mg/kg | 30 | | 38 | | 23.53% |
| | | Calcium | mg/kg | 339000 | | 312000 | | 8.29% |
| | | Magnesium | mg/kg | 4920 | | 5780 | | 16.07% |
| | | Potassium | mg/kg | 1860 | | 2230 | | 18.09% |
| | | Silicon | mg/kg | 1860 | | 2050 | | 9.72% |
| | | Strontium | mg/kg | 293 | | 286 | | 2.42% |
| | | Aluminum | mg/kg | 5760 | | 8180 | | 34.72% |
| | | Arsenic | mg/kg | 11 | | 8.3 | | 27.98% |
| | | Barium | mg/kg | 42.7 | | 55.7 | | 26.42% |
| | | Beryllium | mg/kg | 0.436 | J | 0.54 | J | 21.31% |
| | | Cadmium | mg/kg | 0.612 | J | 0.56 | J | 8.87% |
| | | Chromium | mg/kg | 11.6 | | 14 | | 18.75% |
| | | Copper | mg/kg | 8.3 | | 8.41 | | 1.32% |
| | | Iron | mg/kg | 15300 | | 13800 | | 10.31% |
| | | Lead | mg/kg | 17.9 | | 19.6 | | 9.07% |
| | | Manganese | mg/kg | 439 | | 471 | | 7.03% |
| | | Nickel | mg/kg | 18 | | 14.9 | | 18.84% |
| | | Selenium | mg/kg | <0.451 | | 0.67 | J | 39.07% |
| | | Silver | mg/kg | 0.282 | J | <0.262 | | 7.35% |
| | | Zinc | mg/kg | 42.6 | | 43.5 | | 2.09% |
| | | Mercury | mg/kg | 0.0166 | J | 0.0268 | J | 47.00% |
| | | Acetone | ug/kg | 14 | J | <12 | | 15.38% |
| | | p-Isopropyltoluene | ug/kg | <1.1 | | 1.4 | J | 24.00% |
| | | Toluene | ug/kg | <0.89 | | 2.2 | J | 84.79% |
| | | 2-Methylnaphthalene | mg/kg | 0.15 | J | <0.11 | | 30.77% |
| | | Di-n-Octyl Phthalate | mg/kg | 0.48 | J | <0.19 | | 86.57% |
| | | Fluoranthene | mg/kg | 0.11 | J | <0.11 | | 0.00% |
| | | pH | pH units | 7.46 | | 7.49 | | 0.40% |
| | | TOC | mg/kg | 25000 | | 22000 | | 12.77% |
| | | Bicarbonate | mg/kg | 440 | | 360 | | 20.00% |
| | | Total Alkalinity | mg/kg | 440 | | 360 | | 20.00% |
| | | Total dissolved solids | mg/kg | 5500 | | 7660 | | 32.83% |
| | | Phosphorous, total | mg/kg | 3.1 | | 3.5 | | 12.12% |

| Sample Location | Date Collected | Analyte | Units | Field Sample | Qualifiers | Duplicate | Qualifiers | RPD |
|-----------------------------------|----------------|-------------------|-------|--------------|------------|-----------|------------|--------|
| Passive Diffusion Sampling | | | | | | | | |
| HCS 440 | 2/17/2015 | Tetrachloroethene | µg | 0.28 | | 0.28 | | 0.00% |
| | | TPH | µg | 0.61 | | 0.76 | | 21.90% |
| | 4/21/2015 | Tetrachloroethene | µg | 0.41 | | 0.37 | | 10.26% |
| | | TPH | µg | 0.65 | | 1.05 | | 47.06% |
| | 6/15/2015 | Tetrachloroethene | µg | 0.4 | | 0.42 | | 4.88% |
| | | TPH | µg | 0.72 | | <0.5 | | 36.07% |
| | 8/27/2015 | Tetrachloroethene | µg | 0.39 | | 0.44 | | 12.05% |
| | | TPH | µg | 1.23 | | 1.44 | | 15.73% |
| | 10/20/2015 | Tetrachloroethene | µg | 0.47 | | 0.46 | | 2.15% |
| | 12/15/2015 | Tetrachloroethene | µg | 0.38 | | 0.39 | | 2.63% |
| HSM 430 | 2/17/2015 | Tetrachloroethene | µg | 0.13 | | 0.16 | | 20.69% |
| | | TPH | µg | 1.27 | | 0.99 | | 24.78% |
| | 4/21/2015 | Tetrachloroethene | µg | 0.8 | | 0.72 | | 10.53% |
| | | Toluene | µg | <0.02 | | <0.02 | | 0.00% |
| | | TPH | µg | <0.5 | | 0.54 | | 7.69% |
| | 6/15/2015 | Tetrachloroethene | µg | 1.98 | | 2.35 | | 17.09% |
| | | TPH | µg | 0.96 | | 0.54 | | 56.00% |
| | 8/27/2015 | Tetrachloroethene | µg | 0.68 | | 0.89 | | 26.75% |
| | 10/20/2015 | Tetrachloroethene | µg | 0.94 | | 1.01 | | 7.18% |
| | 12/15/2015 | Chloroform | µg | 0.02 | | 0.02 | | 0.00% |
| | | Tetrachloroethene | µg | 0.81 | | 0.76 | | 6.37% |

B - Analyte detected in the laboratory method blank at the quantity listed

BV,BU - Sample exceeded hold time

H - Sample exceeded hold time

J - Result above the method detection limit but below the reporting limit

Method detection limits or reporting limits were used to calculate RPD for results not detected above these limits.

APPENDIX I

SAMPLE RECORD

COMAL SURFACE WATER

| Location / Sample Name | Date Sampled | Time Sampled | Latitude (dd) | Longitude (dd) | County | Location Generic Name |
|------------------------|--------------|--------------|---------------|----------------|------------|-----------------------|
| HCS110 | 3/16/2015 | 12:44 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS120 | 3/16/2015 | 13:20 | 29.718084 | -98.131644 | Comal | Upper Landa Lake |
| FDHCS120 | 3/16/2015 | 13:20 | 29.709566 | -98.133749 | Comal | Lower Landa Lake |
| HCS130 | 3/16/2015 | 11:29 | 29.709566 | -98.133749 | Comal | Lower Landa Lake |
| HCS140 | 3/16/2015 | 13:52 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS160 | 3/16/2015 | 14:13 | 29.707454 | -98.122762 | Comal | USGS Gauge |
| TB02 | 3/16/2015 | NA | NA | NA | Comal | VOC Blank |
| HCS110 | 9/9/2015 | 10:11 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS120 | 9/9/2015 | 10:44 | 29.718084 | -98.131644 | Comal | Upper Landa Lake |
| FDHCS120 | 9/9/2015 | 10:44 | 29.709566 | -98.133749 | Comal | Lower Landa Lake |
| HCS130 | 9/9/2015 | 9:30 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS140 | 9/9/2015 | 11:12 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS160 | 9/9/2015 | 11:39 | 29.707454 | -98.122762 | Comal | USGS Gauge |
| TB11 | 9/9/2015 | NA | NA | NA | Comal/Hays | VOC Blank |

COMAL STORM WATER

| Location / Sample Name | Date Sampled | Time Sampled | Latitude (dd) | Longitude (dd) | County | Location Generic Name |
|------------------------|--------------|--------------|---------------|----------------|--------|------------------------------|
| HCS210 Lead | 1/22/2015 | 0:36 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS240 Lead | 1/22/2015 | 0:51 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS250 Lead | 1/22/2015 | 0:35 | 29.709491 | -98.122578 | Comal | Lower Old Channel |
| HCS260 Lead | 1/22/2015 | 1:10 | 29.708007 | -98.127301 | Comal | New Channel |
| HCS270 Lead | 1/22/2015 | 0:54 | 29.704014 | -98.115791 | Comal | Comal River above confluence |
| HCS210 Peak | 1/22/2015 | 16:28 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS240 Peak | 1/22/2015 | 16:51 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS250 Peak | 1/22/2015 | 16:40 | 29.709491 | -98.122578 | Comal | Lower Old Channel |
| HCS260 Peak | 1/22/2015 | 17:03 | 29.708007 | -98.127301 | Comal | New Channel |
| HCS270 Peak | 1/22/2015 | 16:55 | 29.704014 | -98.115791 | Comal | Comal River above confluence |
| HCS210 Trail | 1/23/2015 | 3:01 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS240 Trail | 1/23/2015 | 3:21 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS250 Trail | 1/23/2015 | 3:03 | 29.709491 | -98.122578 | Comal | Lower Old Channel |
| HCS260 Trail | 1/23/2015 | 3:36 | 29.708007 | -98.127301 | Comal | New Channel |
| FDHCS260 Trail | 1/23/2015 | 3:36 | 29.708007 | -98.127301 | Comal | New Channel |
| HCS270 Trail | 1/23/2015 | 3:30 | 29.704014 | -98.115791 | Comal | Comal River above confluence |

| | | | | | | |
|----------------|------------|-------|-----------|------------|-------|------------------------------|
| FDHCS270 Trail | 1/23/2015 | 3:30 | 29.704014 | -98.115791 | Comal | Comal River above confluence |
| TB01 | 1/23/2015 | NA | NA | NA | Comal | VOC Blank |
| HCS210 Lead | 10/23/2015 | 11:46 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS240 Lead | 10/23/2015 | 12:00 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS250 Lead | 10/23/2015 | 12:03 | 29.709491 | -98.122578 | Comal | Lower Old Channel |
| HCS260 Lead | 10/23/2015 | 12:24 | 29.708007 | -98.127301 | Comal | New Channel |
| HCS270 Lead | 10/23/2015 | 12:20 | 29.704014 | -98.115791 | Comal | Comal River above confluence |
| HCS210 Peak | 10/23/2015 | 14:19 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS240 Peak | 10/23/2015 | 14:36 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS250 Peak | 10/23/2015 | 14:59 | 29.709491 | -98.122578 | Comal | Lower Old Channel |
| HCS260 Peak | 10/23/2015 | 15:01 | 29.708007 | -98.127301 | Comal | New Channel |
| HCS270 Peak | 10/23/2015 | 15:16 | 29.704014 | -98.115791 | Comal | Comal River above confluence |
| HCS210 Trail | 10/23/2015 | 16:14 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS240 Trail | 10/23/2015 | 16:29 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS250 Trail | 10/23/2015 | 17:09 | 29.709491 | -98.122578 | Comal | Lower Old Channel |
| HCS260 Trail | 10/23/2015 | 16:50 | 29.708007 | -98.127301 | Comal | New Channel |
| FDHCS260 Trail | 10/23/2015 | 16:50 | 29.708007 | -98.127301 | Comal | New Channel |
| HCS270 Trail | 10/23/2015 | 17:34 | 29.704014 | -98.115791 | Comal | Comal River above confluence |
| FDHCS270 Trail | 10/23/2015 | 17:34 | 29.704014 | -98.115791 | Comal | Comal River above confluence |
| TB14 | 10/23/2015 | NA | NA | NA | Comal | VOC Blank |

COMAL SEDIMENT

| Location / Sample Name | Date Sampled | Time Sampled | Latitude (dd) | Longitude (dd) | County | Location Generic Name |
|-------------------------------|---------------------|---------------------|----------------------|-----------------------|---------------|------------------------------|
| HCS310 | 6/4/2015 | 11:22 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS320 | 6/4/2015 | 11:43 | 29.718084 | -98.131644 | Comal | Upper Landa Lake |
| HCS330 | 6/4/2015 | 12:20 | 29.709566 | -98.133749 | Comal | Lower Landa Lake |
| HCS340 | 6/4/2015 | 13:57 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS360 | 6/4/2015 | 14:24 | 29.707454 | -98.122762 | Comal | USGS Gauge |
| FDHCS360 | 6/4/2015 | 14:24 | 29.707454 | -98.122762 | Comal | USGS Gauge |
| EB01 | 6/5/2015 | 15:41 | NA | NA | Comal/Hays | Equipment Blank |
| EB02 | 6/10/2015 | 13:26 | NA | NA | Comal/Hays | Equipment Blank |
| TB07 | 6/5/2015 | NA | NA | NA | Comal | VOC Blank |

COMAL PDS

| Location / Sample Name | Installed | Retrieved | Latitude (dd) | Longitude (dd) | County | Location Generic Name |
|-----------------------------------|--------------------|--------------------|--------------------------|---------------------------|----------------|----------------------------------|
| HCS410 | 2/3/2015 10:02 | 2/17/2015 11:56 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS420 | 2/3/2015 10:13 | 2/17/2015 12:01 | 29.718084 | -98.131644 | Comal | Upper Landa Lake |
| HCS430 | 2/3/2015 9:11 | 2/17/2015 11:25 | 29.709566 | -98.133749 | Comal | Lower Landa Lake |
| HCS440 | 2/3/2015 10:27 | 2/17/2015 12:12 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| FDHCS440 | 2/3/2015 10:27 | 2/17/2015 12:12 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS460 | 2/3/2015 10:38 | 2/17/2015 12:24 | 29.707454 | -98.122762 | Comal | USGS Gauge |
| TB02 | 2/3/2015 9:11 | 2/17/2015 12:24 | NA | NA | Comal/ Hays | Trip Blank |
| HCS410 | 4/7/2015 10:00 | 4/21/2015 11:34 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS420 | 4/7/2015 10:11 | 4/21/2015 11:50 | 29.718084 | -98.131644 | Comal | Upper Landa Lake |
| HCS430 | 4/7/2015 9:41 | 4/21/2015 11:10 | 29.709566 | -98.133749 | Comal | Lower Landa Lake |
| HCS440 | 4/7/2015 10:23 | 4/21/2015 12:05 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| FDHCS440 | 4/7/2015 10:23 | 4/21/2015 12:05 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS460 | 4/7/2015 10:33 | 4/21/2015 12:19 | 29.707454 | -98.122762 | Comal | USGS Gauge |
| TB05 | 4/7/2015 9:41 | 4/21/2015 15:23 | NA | NA | Comal/ Hays | Trip Blank |
| HCS410 | 6/1/2015 11:03 | 6/15/2015 10:39 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS420 | 6/1/2015 11:30 | 6/15/2015 10:57 | 29.718084 | -98.131644 | Comal | Upper Landa Lake |
| HCS430 | 6/1/2015 10:36 | 6/15/2015 10:26 | 29.709566 | -98.133749 | Comal | Lower Landa Lake |
| HCS440 | 6/1/2015 11:44 | 6/15/2015 11:07 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| FDHCS440 | 6/1/2015 11:44 | 6/15/2015 11:07 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS460 | 6/1/2015 11:58 | 6/15/2015 11:17 | 29.707454 | -98.122762 | Comal | USGS Gauge |
| TB09 | 6/1/2015 10:36 | 6/15/2015 14:01 | NA | NA | Comal/ Hays | Test Blank |
| HCS410 | 8/13/2015 11:01 | 8/27/2015 13:16 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS420 | 8/13/2015 11:08 | 8/27/2015 13:25 | 29.718084 | -98.131644 | Comal | Upper Landa Lake |

| Location / Sample Name | Installed | Retrieved | Latitude (dd) | Longitude (dd) | County | Location Generic Name |
|------------------------|--------------------|---------------------|---------------|----------------|----------------|-----------------------|
| HCS430 | 8/13/2015 10:40 | 8/27/2015 12:55 | 29.709566 | -98.133749 | Comal | Lower Landa Lake |
| HCS440 | 8/13/2015 11:28 | 8/27/2015 13:45 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| FDHCS440 | 8/13/2015 11:28 | 8/27/2015 13:45 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS460 | 8/13/2015 11:36 | 8/27/2015 13:54 | 29.707454 | -98.122762 | Comal | USGS Gauge |
| TB10 | 8/13/2015 10:40 | 8/27/2015 13:54 | NA | NA | Comal/ Hays | Test Blank |
| HCS410 | 10/6/2015 12:19 | 10/20/2015 10:32 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS420 | 10/6/2015 12:37 | 10/20/2015 10:42 | 29.718084 | -98.131644 | Comal | Upper Landa Lake |
| HCS430 | 10/6/2015 23:56 | 10/20/2015 10:16 | 29.709566 | -98.133749 | Comal | Lower Landa Lake |
| HCS440 | 10/6/2015 12:47 | 10/20/2015 10:54 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| FDHCS440 | 10/6/2015 12:47 | 10/20/2015 10:54 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS460 | 10/6/2015 13:03 | 10/20/2015 11:03 | 29.707454 | -98.122762 | Comal | USGS Gauge |
| TB13 | 10/6/2015 12:19 | 10/20/2015 13:03 | NA | NA | Comal/ Hays | Test Blank |
| HCS410 | 12/1/2015 10:28 | 12/15/2015 10:44 | 29.72043 | -98.12525 | Comal | Upper Springs |
| HCS420 | 12/1/2015 10:52 | 12/15/2015 10:52 | 29.718084 | -98.131644 | Comal | Upper Landa Lake |
| HCS430 | 12/1/2015 10:13 | 12/15/2015 10:28 | 29.709566 | -98.133749 | Comal | Lower Landa Lake |
| HCS440 | 12/1/2015 11:06 | 12/15/2015 11:01 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| FDHCS440 | 12/1/2015 11:06 | 12/15/2015 11:01 | 29.710221 | -98.129534 | Comal | Upper Old Channel |
| HCS460 | 12/1/2015 11:23 | 12/15/2015 11:11 | 29.707454 | -98.122762 | Comal | USGS Gauge |
| TB16 | 12/1/2015 10:13 | 12/15/2015 13:28 | NA | NA | Comal/ Hays | Test Blank |

SAN MARCOS SURFACE WATER

| Location / Sample Name | Date Sampled | Time Sampled | Latitude (dd) | Longitude (dd) | County | Location Generic Name |
|------------------------|--------------|--------------|---------------|----------------|--------|-----------------------|
| HSM 110 | 3/25/2015 | 10:28 | 29.893566 | -97.927631 | Hays | Sink Creek |
| FDHSM 110 | 3/25/2015 | 10:28 | 29.890258 | -97.934568 | Hays | Spring Lake |
| HSM 120 | 3/25/2015 | 10:51 | 29.889831 | -97.935957 | Hays | Sessoms Creek |

| | | | | | | |
|-----------|-----------|-------|-----------|------------|------|---------------|
| HSM 130 | 3/25/2015 | 11:17 | 29.883955 | -97.935295 | Hays | City Park |
| HSM 140 | 3/25/2015 | 11:30 | 29.880016 | -97.932977 | Hays | Rio Vista Dam |
| HSM 150 | 3/25/2015 | 12:25 | 29.87469 | -97.931603 | Hays | 1-35 Reach |
| HSM 160 | 3/25/2015 | 12:50 | 29.87469 | -97.931603 | Hays | 1-35 Reach |
| HSM 170 | 3/25/2015 | 13:18 | 29.868809 | -97.930378 | Hays | Capes Dam |
| TB04 | 3/25/2015 | NA | NA | NA | Hays | Test Blank |
| HSM 110 | 9/17/2015 | 10:53 | 29.893566 | -97.927631 | Hays | Sink Creek |
| FDHSM 110 | 9/17/2015 | 10:53 | 29.890258 | -97.934568 | Hays | Spring Lake |
| HSM 120 | 9/17/2015 | 11:30 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM 130 | 9/17/2015 | 11:57 | 29.883955 | -97.935295 | Hays | City Park |
| HSM 140 | 9/17/2015 | 12:24 | 29.880016 | -97.932977 | Hays | Rio Vista Dam |
| HSM 150 | 9/17/2015 | 12:51 | 29.87469 | -97.931603 | Hays | 1-35 Reach |
| HSM 160 | 9/17/2015 | 13:18 | 29.87469 | -97.931603 | Hays | 1-35 Reach |
| HSM 170 | 9/17/2015 | 13:47 | 29.868809 | -97.930378 | Hays | Capes Dam |
| TB12 | 9/17/2015 | NA | NA | NA | Hays | Test Blank |

SAN MARCOS STORM WATER

| Location / Sample Name | Date Sampled | Time Sampled | Latitude (dd) | Longitude (dd) | County | Location Generic Name |
|-----------------------------------|-------------------------|-------------------------|--------------------------|---------------------------|---------------|----------------------------------|
| HSM210 Lead | 5/5/2015 | 22:10 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM230 Lead | 5/5/2015 | 22:20 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM231 Lead | 5/5/2015 | 22:11 | 29.886254 | -97.935891 | Hays | Dog Beach Outflow |
| HSM240 Lead | 5/5/2015 | 22:49 | 29.884145 | -97.935554 | Hays | City Park |
| HSM250 Lead | 5/5/2015 | 22:27 | 29.882213 | -97.934867 | Hays | Purgatory Creek |
| HSM260 Lead | 5/5/2015 | 22:46 | 29.87484 | -97.931713 | Hays | 1-35 Reach |
| HSM270 Lead | 5/5/2015 | 23:06 | 29.868809 | -97.930378 | Hays | Capes Dam |
| HSM210 Peak | 5/5/2015 | 23:24 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM230 Peak | 5/5/2015 | 23:35 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM231 Peak | 5/5/2015 | 23:50 | 29.886254 | -97.935891 | Hays | Dog Beach Outflow |
| HSM240 Peak | 5/6/2015 | 00:05 | 29.884145 | -97.935554 | Hays | City Park |
| HSM250 Peak | 5/5/2015 | 23:30 | 29.882213 | -97.934867 | Hays | Purgatory Creek |
| HSM260 Peak | 5/5/2015 | 23:49 | 29.87484 | -97.931713 | Hays | 1-35 Reach |
| HSM270 Peak | 5/6/2015 | 00:06 | 29.868809 | -97.930378 | Hays | Capes Dam |
| HSM210 Trail | 5/6/2015 | 1:30 | 29.893566 | -97.927631 | Hays | Sink Creek |
| FDHSM210 Trail | 5/6/2015 | 1:30 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM230 Trail | 5/6/2015 | 2:15 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| FDHSM230 Trail | 5/6/2015 | 2:15 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM231 Trail | 5/6/2015 | 2:40 | 29.886254 | -97.935891 | Hays | Dog Beach Outflow |
| FDHSM231 Trail | 5/6/2015 | 2:40 | 29.886254 | -97.935891 | Hays | Dog Beach Outflow |
| HSM240 Trail | 5/6/2015 | 3:05 | 29.884145 | -97.935554 | Hays | City Park |
| HSM250 Trail | 5/6/2015 | 3:58 | 29.882213 | -97.934867 | Hays | Purgatory Creek |

| Location / Sample Name | Date Sampled | Time Sampled | Latitude (dd) | Longitude (dd) | County | Location Generic Name |
|------------------------|--------------|--------------|---------------|----------------|----------------|-----------------------|
| HSM260 Trail | 5/6/2015 | 4:20 | 29.87484 | -97.931713 | Hays | 1-35 Reach |
| HSM270 Trail | 5/6/2015 | 4:40 | 29.868809 | -97.930378 | Hays | Capes Dam |
| TB06 | 5/6/2015 | NA | NA | NA | Hays | VOC Blank |
| HSM210 Lead | 10/23/2015 | 15:25 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM230 Lead | 10/23/2015 | 15:41 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM231 Lead | 10/23/2015 | 16:02 | 29.886254 | -97.935891 | Hays | Dog Beach Outflow |
| HSM240 Lead | 10/23/2015 | 15:30 | 29.884145 | -97.935554 | Hays | City Park |
| HSM250 Lead | 10/23/2015 | 16:21 | 29.882213 | -97.934867 | Hays | Purgatory Creek |
| HSM260 Lead | 10/23/2015 | 15:58 | 29.87484 | -97.931713 | Hays | 1-35 Reach |
| HSM270 Lead | 10/23/2015 | 16:15 | 29.868809 | -97.930378 | Hays | Capes Dam |
| HSM210 Peak | 10/23/2015 | 17:02 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM230 Peak | 10/23/2015 | 17:16 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM231 Peak | 10/23/2015 | 17:06 | 29.886254 | -97.935891 | Hays | Dog Beach Outflow |
| HSM240 Peak | 10/23/2015 | 17:34 | 29.884145 | -97.935554 | Hays | City Park |
| HSM250 Peak | 10/23/2015 | 17:05 | 29.882213 | -97.934867 | Hays | Purgatory Creek |
| HSM260 Peak | 10/23/2015 | 17:25 | 29.87484 | -97.931713 | Hays | 1-35 Reach |
| HSM270 Peak | 10/23/2015 | 17:45 | 29.868809 | -97.930378 | Hays | Capes Dam |
| HSM210 Trail | 10/23/2015 | 20:35 | 29.893566 | -97.927631 | Hays | Sink Creek |
| FDHSM210 Trail | 10/23/2015 | 20:35 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM230 Trail | 10/23/2015 | 21:01 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| FDHSM230 Trail | 10/23/2015 | 21:01 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM231 Trail | 10/23/2015 | 21:21 | 29.886254 | -97.935891 | Hays | Dog Beach Outflow |
| FDHSM231 Trail | 10/23/2015 | 21:21 | 29.886254 | -97.935891 | Hays | Dog Beach Outflow |
| HSM240 Trail | 10/23/2015 | 20:51 | 29.884145 | -97.935554 | Hays | City Park |
| HSM250 Trail | 10/23/2015 | 21:06 | 29.882213 | -97.934867 | Hays | Purgatory Creek |
| HSM260 Trail | 10/23/2015 | 21:35 | 29.87484 | -97.931713 | Hays | 1-35 Reach |
| HSM270 Trail | 10/23/2015 | 21:55 | 29.868809 | -97.930378 | Hays | Capes Dam |
| TB15 | 10/23/2015 | NA | NA | NA | Comal/ Hays | VOC Blank |

SAN MARCOS SEDIMENT

| Location / Sample Name | Date Sampled | Time Sampled | Latitude (dd) | Longitude (dd) | County | Location Generic Name |
|------------------------|--------------|--------------|---------------|----------------|--------|-----------------------|
| HSM 310 | 6/5/2015 | 10:03 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM 320 | 6/5/2015 | 10:48 | 29.890258 | -97.934568 | Hays | Spring Lake |
| HSM 330 | 6/5/2015 | 11:03 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM 340 | 6/5/2015 | 11:21 | 29.883955 | -97.935295 | Hays | City Park |
| HSM 350 | 6/5/2015 | 12:24 | 29.880016 | -97.932977 | Hays | Rio Vista Dam |
| HSM 360 | 6/5/2015 | 13:13 | 29.880016 | -97.932977 | Hays | Rio Vista Dam |

| | | | | | | |
|-----------|----------|-------|-----------|------------|------|------------|
| HSM 370 | 6/5/2015 | 13:51 | 29.87469 | -97.931603 | Hays | 1-35 Reach |
| FDHSM 370 | 6/5/2015 | 13:51 | 29.868809 | -97.930378 | Hays | Capes Dam |
| TB08 | 6/5/2015 | NA | NA | NA | Hays | VOC Blank |

SAN MARCOS PDS

| Location / Sample Name | Installed | Retrieved | Latitude (dd) | Longitude (dd) | County | Location Generic Name |
|-----------------------------------|-------------------|--------------------|--------------------------|---------------------------|---------------|--------------------------------------|
| HSM 410 | 2/3/2015 11:32 | 2/17/2015 9:45 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM 420 | 2/3/2015 11:45 | 2/17/2015 9:55 | 29.890258 | -97.934568 | Hays | Spring Lake |
| HSM 430 | 2/3/2015 11:55 | 2/17/2015 10:01 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| FDHSM 430 | 2/3/2015 11:55 | 2/17/2015 10:01 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM 440 | 2/3/2015 12:06 | 2/17/2015 10:08 | 29.883955 | -97.935295 | Hays | City Park |
| HSM 450 | 2/3/2015 12:22 | 2/17/2015 10:22 | 29.880016 | -97.932977 | Hays | Rio Vista Dam |
| HSM 460 | 2/3/2015 12:33 | 2/17/2015 10:35 | 29.87469 | -97.931603 | Hays | 1-35 Reach |
| HSM 470 | 2/3/2015 12:45 | 2/17/2015 10:47 | 29.868809 | -97.930378 | Hays | Capes Dam |
| HSM 410 | 4/7/2015 11:14 | 4/21/2015 13:51 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM 420 | 4/7/2015 11:26 | 4/21/2015 14:02 | 29.890258 | -97.934568 | Hays | Spring Lake |
| HSM 430 | 4/7/2015 11:36 | 4/21/2015 14:21 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| FDHSM 430 | 4/7/2015 11:36 | 4/21/2015 14:21 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM 440 | 4/7/2015 11:49 | 4/21/2015 14:35 | 29.883955 | -97.935295 | Hays | City Park |
| HSM 450 | 4/7/2015 12:31 | 4/21/2015 14:52 | 29.880016 | -97.932977 | Hays | Rio Vista Dam |
| HSM 460 | 4/7/2015 12:50 | 4/21/2015 15:11 | 29.87469 | -97.931603 | Hays | 1-35 Reach |
| HSM 470 | 4/7/2015 13:02 | 4/21/2015 15:23 | 29.868809 | -97.930378 | Hays | Capes Dam |
| HSM 410 | 6/1/2015 13:47 | 6/15/2015 12:14 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM 420 | 6/1/2015 13:58 | 6/15/2015 12:26 | 29.890258 | -97.934568 | Hays | Spring Lake |
| HSM 430 | 6/1/2015 14:08 | 6/15/2015 12:34 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| FDHSM 430 | 6/1/2015 14:08 | 6/15/2015 12:34 | 29.889831 | -97.935957 | Hays | Sessoms Creek |

| Location / Sample Name | Installed | Retrieved | Latitude (dd) | Longitude (dd) | County | Location Generic Name |
|-----------------------------------|--------------------|---------------------|--------------------------|---------------------------|---------------|--------------------------------------|
| HSM 440 | 6/1/2015 14:23 | 6/15/2015 12:47 | 29.883955 | -97.935295 | Hays | City Park |
| HSM 450 | 6/1/2015 14:46 | 6/15/2015 15:30 | 29.880016 | -97.932977 | Hays | Rio Vista Dam |
| HSM 460 | 6/1/2015 15:34 | 6/15/2015 15:43 | 29.87469 | -97.931603 | Hays | 1-35 Reach |
| HSM 470 | 6/1/2015 15:59 | 6/15/2015 16:01 | 29.868809 | -97.930378 | Hays | Capes Dam |
| HSM 410 | 8/13/2015 13:08 | 8/27/2015 9:48 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM 420 | 8/13/2015 13:21 | 8/27/2015 10:00 | 29.890258 | -97.934568 | Hays | Spring Lake |
| HSM 430 | 8/13/2015 13:27 | 8/27/2015 10:08 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| FDHSM 430 | 8/13/2015 13:27 | 8/27/2015 10:08 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM 440 | 8/13/2015 13:39 | 8/27/2015 10:19 | 29.883955 | -97.935295 | Hays | City Park |
| HSM 450 | 8/13/2015 13:56 | 8/27/2015 10:51 | 29.880016 | -97.932977 | Hays | Rio Vista Dam |
| HSM 460 | 8/13/2015 14:25 | 8/27/2015 11:06 | 29.87469 | -97.931603 | Hays | 1-35 Reach |
| HSM 470 | 8/13/2015 14:38 | 8/27/2015 11:20 | 29.868809 | -97.930378 | Hays | Capes Dam |
| HSM 410 | 10/6/2015 13:45 | 10/20/2015 11:54 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM 420 | 10/6/2015 13:59 | 10/20/2015 12:04 | 29.890258 | -97.934568 | Hays | Spring Lake |
| HSM 430 | 10/6/2015 14:13 | 10/20/2015 12:10 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| FDHSM 430 | 10/6/2015 14:13 | 10/20/2015 12:10 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM 440 | 10/6/2015 14:48 | 10/20/2015 12:22 | 29.883955 | -97.935295 | Hays | City Park |
| HSM 450 | 10/6/2015 14:44 | 10/20/2015 12:38 | 29.880016 | -97.932977 | Hays | Rio Vista Dam |
| HSM 460 | 10/6/2015 14:55 | 10/20/2015 12:53 | 29.87469 | -97.931603 | Hays | 1-35 Reach |
| HSM 470 | 10/6/2015 15:12 | 10/20/2015 13:03 | 29.868809 | -97.930378 | Hays | Capes Dam |
| HSM 410 | 12/1/2015 12:48 | 12/15/2015 12:06 | 29.893566 | -97.927631 | Hays | Sink Creek |
| HSM 420 | 12/1/2015 13:05 | 12/15/2015 12:14 | 29.890258 | -97.934568 | Hays | Spring Lake |
| HSM 430 | 12/1/2015 13:15 | 12/15/2015 12:23 | 29.889831 | -97.935957 | Hays | Sessoms Creek |

| Location / Sample Name | Installed | Retrieved | Latitude (dd) | Longitude (dd) | County | Location Generic Name |
|-----------------------------------|--------------------|---------------------|--------------------------|---------------------------|---------------|--------------------------------------|
| FDHSM 430 | 12/1/2015 13:15 | 12/15/2015 12:23 | 29.889831 | -97.935957 | Hays | Sessoms Creek |
| HSM 440 | 12/1/2015 13:30 | 12/15/2015 12:35 | 29.883955 | -97.935295 | Hays | City Park |
| HSM 450 | 12/1/2015 13:45 | 12/15/2015 13:09 | 29.880016 | -97.932977 | Hays | Rio Vista Dam |
| HSM 460 | 12/1/2015 13:58 | 12/15/2015 13:19 | 29.87469 | -97.931603 | Hays | 1-35 Reach |
| HSM 470 | 12/1/2015 14:12 | 12/15/2015 13:28 | 29.868809 | -97.930378 | Hays | Capes Dam |

