

# Baseline Soil and Groundwater Assessment Report

Union Reservoir  
Weld County Road 28  
Longmont, Colorado

December 14, 2018  
Terracon Project No. 22187053



**Prepared for:**  
City of Longmont, Colorado

**Prepared by:**  
Terracon Consultants, Inc.  
Longmont, Colorado

[terracon.com](http://terracon.com)

**Terracon**

Environmental



Facilities



Geotechnical



Materials

December 14, 2018

City of Longmont  
385 Kimbark Street  
Longmont, Colorado 80501

Attn: Mr. Jason Elkins  
P: (303) 651-8310  
E: Jason.Elkins@longmontcolorado.gov

Re: Baseline Soil and Groundwater Assessment Report  
Union Reservoir  
Weld County Road 28  
Longmont, Colorado  
Terracon Project No. 22187053

Dear Mr. Elkins:

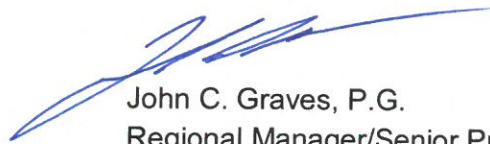
Terracon Consultants, Inc. (Terracon) is pleased to submit our report for Baseline Soil and Groundwater Assessment activities performed at the above referenced site. The report presents data from recent field activities that included the completion of soil borings and the collection groundwater samples for laboratory analysis. Terracon conducted this assessment in general accordance with our proposal (P22187053), dated September 20, 2018.

Terracon appreciates this opportunity to provide environmental services to the City of Longmont. Should you have any questions or require additional information, please do not hesitate to contact our office.

Sincerely,  
**Terracon Consultants, Inc.**



Jaymee L. Binion  
Staff Scientist



John C. Graves, P.G.  
Regional Manager/Senior Principal

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**BASELINE SOIL AND GROUNDWATER ASSESSMENT  
UNION RESERVOIR  
WELD COUNTY ROAD 28  
LONGMONT, COLORADO**

**Terracon Project No. 22187053  
December 14, 2018**

## **1.0 SITE DESCRIPTION**

|                         |                                                                                                                                                                                                                                                                                                                                             |
|-------------------------|---------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Site Name</b>        | Union Reservoir                                                                                                                                                                                                                                                                                                                             |
| <b>Site Address</b>     | South of Weld County Road 28, Longmont, Colorado                                                                                                                                                                                                                                                                                            |
| <b>Site Description</b> | The site is located between Weld County Road 28 to the north and Union Reservoir to the south. The site is undeveloped land sloping to the south. Surrounding areas observed to be rural agricultural land, including a dairy to the northeast. Future oil and gas operations are planned to be located to the north/northeast of the site. |

A Topographic Map showing the site location is included as Exhibit 1 and a Site Diagram is included as Exhibit 2 in Appendix A.

## **2.0 SCOPE OF SERVICES**

In 2012, Terracon was retained by the City of Longmont (City) to assess seventeen plugged and abandoned oil and gas wells located within the City limits. The objective of the 2012 assessment was to provide information concerning the plugging and abandoning of 17 oil and gas (O&G) wellheads located within the City and to assess the potential presence of surficial soil impacts, methane and other gases in the subsurface near the surveyed well locations.

Since 2017, Terracon has been assisting the City with limited soil, groundwater, and soil gas assessments of plugged and abandoned (PA) and active well sites within City limits. Terracon understands that the City of Longmont would like to expand the scope of work from the 2012 and PA projects to include assessing the condition of soil, ground water, and soil gas at future well pad locations, and soil and groundwater baseline conditions of City-owned property and sensitive receptors in areas downgradient of future O&G construction.

The sampling data will be used to establish a baseline data set for the City to assess the potential presence of surficial/subsurface soil impacts and groundwater impacts hydrologically up-gradient of Union Reservoir and potentially down-gradient from the Knight and Olander O&G well pads, being considered for construction to the north/northeast of Union Reservoir. This assessment was not intended to define the extent of potential contamination, but to assess if contaminants of

concern are present in the media in the areas of assessment prior to the installation of the proposed well pads.

## **2.1 Standard of Care**

Terracon's services were performed in a manner consistent with generally accepted practices of the profession undertaken in similar studies in the same geographical area during the same time. Terracon makes no warranties, either express or implied, regarding the findings, conclusions, or recommendations. Please note that Terracon does not warrant the work of laboratories, regulatory agencies, or other third parties supplying information used in the preparation of the report. These services were performed in accordance with the scope of work agreed with you, our client, as reflected in our proposal and were not restricted by ASTM E1903-11.

## **2.2 Additional Scope Limitations**

Findings, conclusions, and recommendations resulting from these services are based upon information derived from the on-site activities and other services performed under this scope of work; such information is subject to change over time. Certain indicators of the presence of hazardous substances, petroleum products, or other constituents may have been latent, inaccessible, unobservable, non-detectable, or not present during these services. We cannot represent that the site contains no hazardous substances, toxic materials, petroleum products, or other latent conditions beyond those identified during this assessment. Subsurface conditions may vary from those encountered at specific borings or wells or during other surveys, tests, assessments, investigations, or exploratory services. The data, interpretations, findings, and our recommendations are based solely upon data obtained at the time and within the scope of these services.

## **2.3 Reliance**

This report has been prepared for the exclusive use of the City of Longmont, and any authorization for use or reliance by any other party (except a governmental entity having jurisdiction over the site) is prohibited without the express written authorization of the City of Longmont and Terracon. Any unauthorized distribution or reuse is at the City of Longmont's sole risk. Notwithstanding the foregoing, reliance by authorized parties will be subject to the terms, conditions, and limitations stated in the proposal, Investigation report, and Terracon's Master Services Agreement (MSA) with the City of Longmont. The limitation of liability defined in the terms and conditions of the MSA is the aggregate limit of Terracon's liability to the City of Longmont and all relying parties unless otherwise agreed in writing.

### 3.0 FIELD INVESTIGATION

#### 3.1 Safety and Subsurface Utilities

Terracon is committed to the safety of all its employees. As such, and in accordance with our Incident and Injury Free® safety goals, Terracon conducted the fieldwork under a site-specific health and safety plan. The plan identified site-specific job hazards and proper pre-task planning procedures. Work was performed using Occupational Safety & Health Administration (OSHA) Level D work attire consisting of hard hats, high-visibility attire, safety glasses, protective gloves, and protective boots. Terracon contacted Colorado 811 and requested location and markings for subsurface utilities that the service was responsible for before commencing intrusive activities at the site.

#### 3.2 Sampling and Analytical Program Summary

Four soil borings were included in the proposed scope of work for this baseline assessment. During a site visit with the client, a mutual decision was reached to reduce the scope of this assessment to three soil borings based on site access. On October 29, 2018, three soil borings (SB-01 through SB-03) were advanced utilizing hand auger methodology to depths of 7 to 8.5 feet below grade surface (bgs). These soil borings were converted to monitoring wells (MW-1 through MW-3). The sample locations were selected to generally represent the areas with the highest potential for detecting constituents of concern based on the proposed location of the well pad and the presumed groundwater flow direction towards Union Reservoir. Refer to the attached Site Diagram (Exhibit 2, Appendix A) for a depiction of the sample locations and pertinent site features. The sampling and analytical program is outlined below.

| <b>SAMPLING AND ANALYTICAL PROGRAM</b> |                                                                                                                                                                                                                                                                                                                                   |
|----------------------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Soil Borings<br/>(Total Depth)</b>  | SB-01 (8.5 feet)<br>SB-02 (7 feet)<br>SB-03 (7 feet)                                                                                                                                                                                                                                                                              |
| <b>Groundwater</b>                     | MW-01, MW-02, and MW-03                                                                                                                                                                                                                                                                                                           |
| <b>Soil Analysis</b>                   | VOCs/TPH-GRO – EPA 8260<br>TPH-DRO/ORO – EPA 8015<br>PAHs – EPA 8270SIM<br>Electrical Conductivity (EC) – EPA 9050A<br>Sodium Adsorption Ratio (SAR) – EPA 3050B<br>pH – EPA 9045D<br>Metals (arsenic, barium, boron, cadmium, chromium III, chromium VI, copper, lead, mercury, nickel, selenium, silver, zinc) – EPA 200.8/6020 |

| <b>SAMPLING AND ANALYTICAL PROGRAM</b> |                                                                                                                                            |
|----------------------------------------|--------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Groundwater Analysis</b>            | VOCs – EPA 8260                                                                                                                            |
|                                        | Dissolved Gases – RSK 175                                                                                                                  |
|                                        | Dissolved Gases CO <sub>2</sub> – EPA 4500CO2 D2011                                                                                        |
|                                        | Metals (arsenic, barium boron, cadmium, chromium III, chromium VI, copper, lead, mercury, nickel, selenium, silver, zinc) – EPA 200.8/6020 |
|                                        | PAHs-8270                                                                                                                                  |
|                                        | Total Dissolved Solids (TDS) – EPA 160.1                                                                                                   |
|                                        | Chloride and Sulfate – EPA 300.0                                                                                                           |

EPA = Environmental Protection Agency; SW-846 analytical methods

VOCs = volatile organic compounds

TPH = total petroleum hydrocarbons

G/D/ORO = gasoline, diesel, and oil range organics

PAH = polycyclic aromatic hydrocarbons

### 3.3 Field Procedures

Soil borings were advanced by a Terracon field professional using a hand auger. The hand auger was cleaned prior to beginning the project and between each soil boring. Soil samples were collected continuously and observed to document soil lithology, color, moisture content and sensory evidence of impairment. The soil samples were field-screened at approximately 1-foot intervals using a MiniRAE photoionization detector (PID) equipped with a 10.6 electron volt ultraviolet lamp source to qualitatively evaluate the potential volatile organic vapors to indicate the presence of VOCs. Terracon calibrated the PID in accordance with the manufacturer's recommendations before the field activities.

Terracon's soil sampling program involved assigning one soil sample from each soil boring for laboratory analysis. The soil sample selected for laboratory analysis was collected from the vadose zone interval directly above the groundwater table observed during boring advancement. The soil samples were collected using Terracon standard operating procedures (SOPs) and field methods.

Following the collection of soil samples, each of the three soil borings were completed as groundwater monitoring wells. Monitoring wells were constructed to approximately 7 to 8.5 feet bgs using 1.0-inch diameter polyvinyl chloride (PVC) with approximately 4 to 6 feet of machine-slotted well screen and 2.5 to 3 feet of PVC riser pipe to the near surface. A silica sand filter pack was placed around the well screen to approximately one foot above the top of well screen, followed by a hydrated bentonite seal above the sand pack filter zone to the near surface. The monitoring wells were fitted with J-plug well caps and bolt-down, flush-mounted well covers set in concrete. Groundwater was measured at depths ranging from 2.6 to 4.5 feet bgs. On October 31, 2018; one groundwater sample was collected from each monitoring well for laboratory analysis using a new, disposable, polypropylene bailer.

Soil and groundwater samples were collected and placed in laboratory-prepared glassware containing the appropriate preservative, labeled, and placed on ice in sample coolers. The sample coolers were either released via chain-of-custody directly from the sampler to a representative of the analytical laboratory or were secured with a custody seal and shipped to the selected analytical laboratory. The sample coolers and completed chain-of-custody forms were relinquished to Pace Analytical in Mount Juliet, Tennessee for analysis for normal turnaround.

## **4.0 RESULTS OF THE FIELD INVESTIGATION**

### **4.1 Geology/Hydrogeology**

Ground-surfacing across the site was observed to be covered with a white powdery substance, indicative of salt accumulations in the near-surface soils. In general, the lithology encountered during boring advancement consisted primarily of clay and silty clay from the ground surface to boring termination at 7 to 8.5 feet bgs. Groundwater was encountered during drilling at approximate depths between 5 to 5.5 feet bgs. Two days following drilling activities, groundwater was observed to have risen to 2.6 to 4.5 bgs. Groundwater is assumed to flow to the south, towards Union Reservoir.

### **4.2 Field Screening**

PID readings above 1 part per million (ppm) were not observed in soil borings completed during this investigation. Staining or odors were not noted during the field investigation.

## **5.0 ANALYTICAL RESULTS**

The laboratory analytical reports and chain-of-custody records are attached in Appendix C. The following sections describe the results of the analytical testing performed as part of this Background Assessment. The constituents of concern concentrations were compared to the May 2016, USEPA, Residential and Industrial RSLs and January 2015 COGCC Table 910-1 (Concentration Levels) for soil. Groundwater analytical results were compared to June 30, 2016 CDPHE Groundwater Quality Standards (GWQSs) and January 2015 COGCC Table 910-1 Groundwater Concentration Levels (910-1 Levels).

### **5.1 Soil Sample Results**

The soil analytical data and corresponding action levels are summarized in Table 1 (Appendix B).



## Baseline Soil and Groundwater Assessment

Union Reservoir ■ Longmont, Colorado  
December 14, 2018 ■ Terracon Project No. 22187053



VOC, PAH, TPH-GRO, TPH-DRO, and TPH-ORO constituents were not reported at concentrations above laboratory detection limits in the soil samples collected during this investigation.

Metal concentrations were reported above laboratory detection limits in the soil samples collected. With the exception of arsenic, reported metal concentrations were observed to be below their applicable action levels and within the expected concentrations for native soils. Arsenic was reported in the soil samples collected from SB-02 (4.13 milligrams per kilogram [mg/kg]) and SB-03 (2.67 mg/kg) above EPA RSLs and COGCC screening level. The Colorado Department of Public Health and Environment (CDPHE) recognizes that arsenic can be naturally occurring and has authored the document titled: *Risk Management Guidance for Evaluating Arsenic Concentrations in Soil*, published 2011, revised in 2014 (CDPHE, 2014). This document states that arsenic has been demonstrated to be naturally occurring in Colorado soils at concentrations significantly higher than the national average. The CDPHE developed an average background concentration of arsenic found in certain native Colorado soils averaging 11 mg/Kg with measured concentrations as low as 3 mg/Kg and as high as 19 mg/Kg. Arsenic concentrations at the site are within published background concentrations.

Physical properties of the soil including the sodium adsorption ratio (SAR), electrical conductivity (EC), and pH were also measured by the laboratory. The physical (inorganic) properties measured for each soil sample are presented below:

### Inorganics in Soils

| Soil Boring  | SAR          | EC<br>(mS/cm) | pH   |
|--------------|--------------|---------------|------|
| SB-01        | 13.4         | 3.10          | 8.83 |
| SB-02        | 10.8         | 2.28          | 8.74 |
| SB-03        | <b>24.6</b>  | <b>9.60</b>   | 8.81 |
| COGCC Levels | Less than 12 | Less than 4   | 6-9  |

The sodium adsorption ratio and electrical conductivity measured in soil boring SB-03 were observed to exceed the COGCC screening levels. The SAR levels indicate that sodium is displacing calcium and magnesium in the soil profile. The measured SAR and EC levels indicate that the ability for the soil in this area to support plant growth is limited. Clay soils, such as those observed during drilling, also tend to have higher measured EC results.

## **5.2 Groundwater Sample Results**

The groundwater analytical data and corresponding action levels are summarized in Table 2 (Appendix B).

Dissolved boron was detected in the groundwater samples from MW-01 (1,110 micrograms per liter [ $\mu\text{g/L}$ ]), MW-02 (1,030  $\mu\text{g/L}$ ), and MW-03 (2,950  $\mu\text{g/L}$ ); exceeding the CDPHE regulatory limit of 750  $\mu\text{g/L}$ . Dissolved selenium was detected in the groundwater samples from MW-02 (465  $\mu\text{g/L}$ ) and MW-03 (1,240  $\mu\text{g/L}$ ); exceeding the CDPHE regulatory limit of 50  $\mu\text{g/L}$ . Other dissolved metals detected above the laboratory detection levels were reported below their respective regulatory limits.

The VOC p-Isopropyltoluene was reported above the laboratory detection level in the groundwater sample collected from MW-01. Neither CDPHE nor COGCC have established a groundwater standard for this constituent.

PAHs fluorene and phenanthrene were detected in the groundwater samples collected during this investigation; however, neither concentrations exceeded CDPHE groundwater standards.

Inorganic constituents analyzed for the groundwater samples included total dissolved solids (TDS), chlorides, and sulfates. The COGCC has defined the groundwater standard exceedance concentrations for chloride and sulfate to be a regional background concentration with a multiplier of 1.25. Terracon utilized 2018 analytical data for chloride and sulfate from the sites sampled during the City of Longmont 2018 Annual Groundwater Quality Monitoring sampling event (Terracon Project No. 22187009) to calculate respective regional background concentrations.

As shown on Table 2, a comparison of the reported concentrations of sulfates and chlorides from this investigation to the background concentrations from previous sampling events in and around the City of Longmont appears to indicate that this area of Longmont has a much higher background concentration of chloride and sulfate than previously measured at other locations. This indicates that the background concentrations may not be applicable to this site.

For TDS, the maximum allowable concentration allowed by CDPHE is also dependent on the background level but can theoretically be any concentration. A TDS concentration of 400,000  $\mu\text{g/L}$  is generally used as a maximum allowable concentration as the background level is established. TDS concentrations reported during this investigation ranged from 19,700,000 to 77,700,000  $\mu\text{g/L}$ .

## **6.0 FINDINGS AND CONCLUSIONS**

### **6.1 Findings**

Based on the scope of services described in this report and subject to the limitations described herein, Terracon's findings include the following.

- n Three soil borings (SB-01 through SB-03) were advanced to approximate depths of 7 to 8.5 feet bgs on October 29, 2018, to characterize the subsurface lithology and to collect soil and groundwater samples for laboratory analysis.
- n In general, the lithology encountered during drilling consisted primarily of clay and silty clay. A white powdery precipitate was observed on the ground surface within the area of the investigation. Groundwater measurements conducted two days following drilling indicated a depth to water ranging from 2.6 to 4.5 feet bgs.
- n Laboratory analysis of soil samples did not indicate elevated concentrations of constituents of concern but did indicate an elevated soil adsorption ratio and electrical conductivity in SB-03 located on the western portion of the investigation area.
- n Laboratory analysis of the groundwater samples indicated the presence of VOCs and PAHs above laboratory detection levels, but below regulatory action levels.
- n Laboratory analysis of groundwater samples indicated that select metals including boron and selenium were detected above regulatory action levels for groundwater for use as a domestic water supply or irrigation water. Additional metals concentrations were reported above laboratory detection levels but below regulatory action levels.
- n Inorganic constituents including total dissolved solids, chlorides, and sulfates were observed to be significantly higher than previous concentrations reported in and around the City of Longmont and were observed to be the highest at MW-03 located on the western portion of the investigation area.

### **6.2 Conclusions**

Inorganic constituent analytical data that included an elevated sodium adsorption ratio and electrical conductivity in one soil sample; apparent elevated total dissolved solids, chlorides, and sulfates in the groundwater samples; and site observations of salt accumulations on the soil surface point to this area being in an alkali area of Longmont. This can be naturally occurring as this would be considered a semi-arid region and it receives relatively little rainfall to leach

## Baseline Soil and Groundwater Assessment

Union Reservoir ■ Longmont, Colorado  
December 14, 2018 ■ Terracon Project No. 22187053



naturally-occurring salts from the soil. The observed clay soils and relatively shallow depth to groundwater could also be contributors. The observed analytical results could also be the result of agricultural activities in the area up-gradient of the site including potential over-application of composts or manure on agricultural fields to the north or over-application of magnesium chloride on County Road 28.

## 7.0 RECOMMENDATIONS

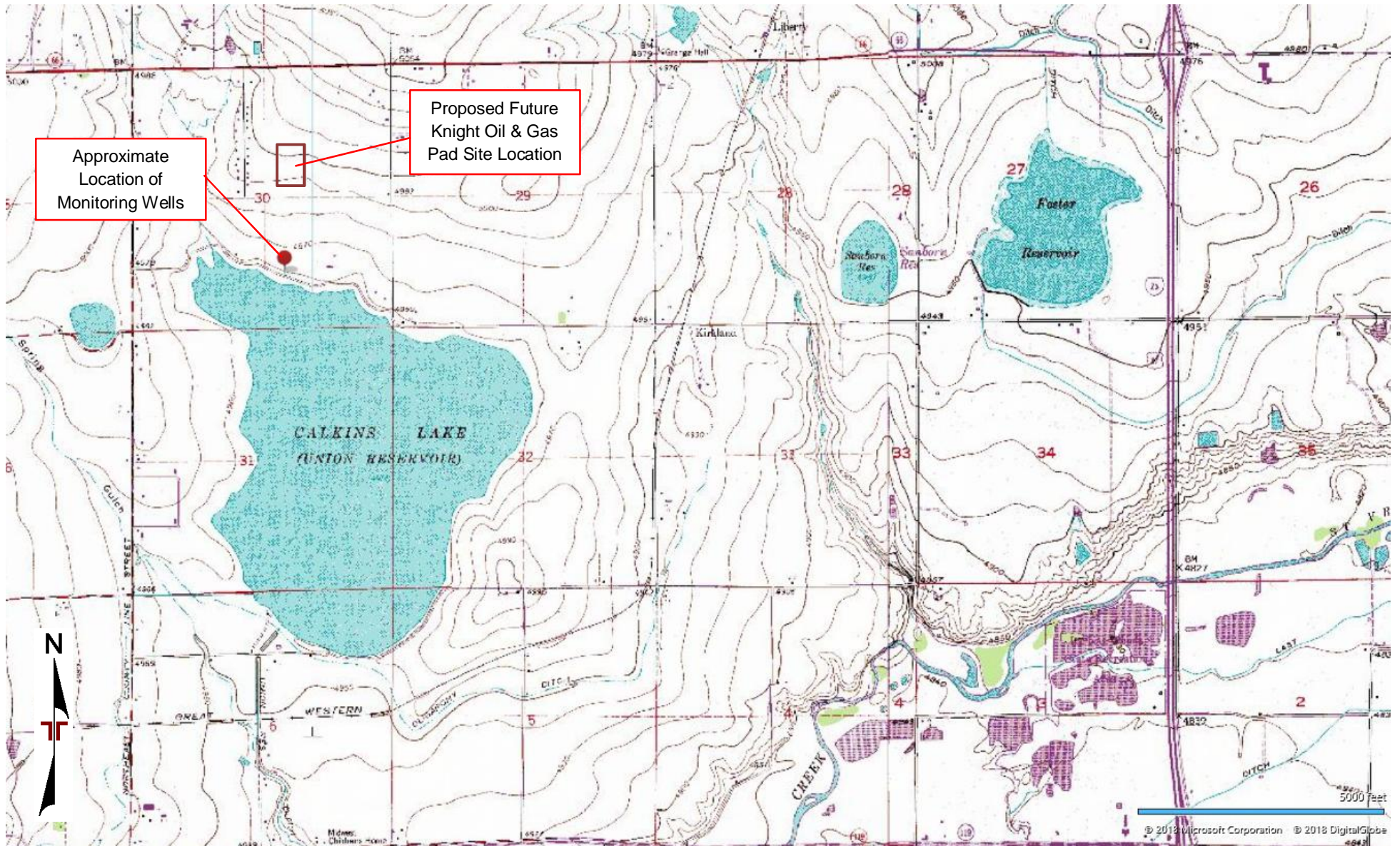
The objective of this investigation was to evaluate the presence of constituents of concern in the on-site soil and groundwater above relevant laboratory detection limits and/or regulatory limits prior to the onset of O&G operations up-gradient of the site.

Based on the scope of services, limitations, and conclusions of this assessment, Terracon recommends continued periodic groundwater monitoring to establish a baseline for the inorganic constituents reported during this investigation to establish a baseline prior to the oil and gas facilities in the area becoming operational.

**APPENDIX A – EXHIBITS**

Exhibit 1 – Topographic Map

Exhibit 2 – Site Diagram



TOPOGRAPHIC MAP IMAGE  
 COURTESY OF THE U.S.  
 GEOLOGICAL SURVEY

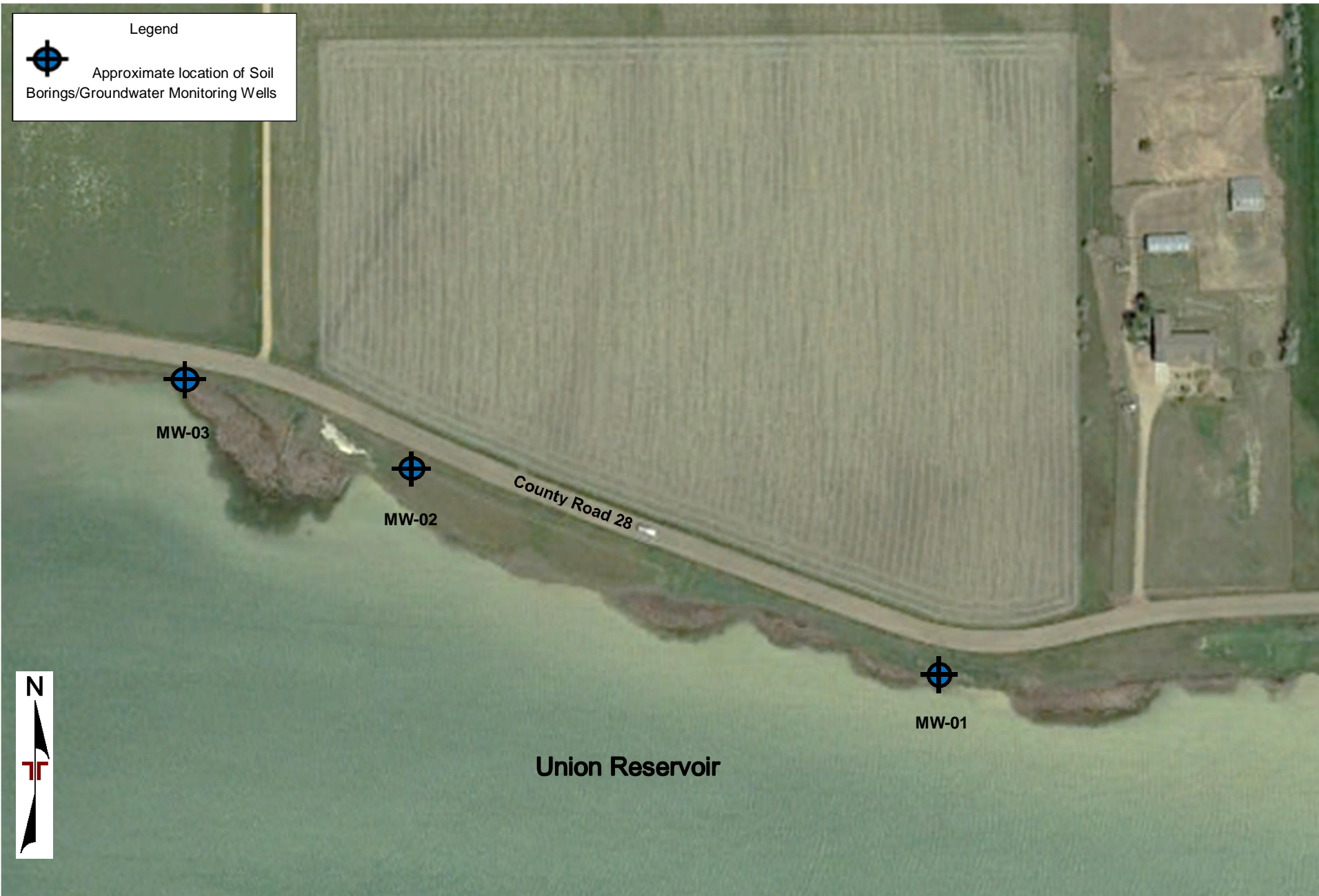
|                         |                         |
|-------------------------|-------------------------|
| Project Manager:<br>MJS | Project No.<br>22187053 |
| Drawn by:<br>JLB        | Scale:<br>SHOWN         |
| Checked by:<br>MJS      | File Name:<br>Site      |
| Approved by:<br>JCG     | Date:<br>12/5/18        |


**Terracon**  
 1901 Sharp Point Dr, Ste C  
 Fort Collins, CO 80525-4429

TOPOGRAPHIC MAP

Union Reservoir Baseline Assessment  
 Weld County Road 28  
 Longmont, Colorado

Exhibit  
 1



Legend  
 Approximate location of Soil Borings/Groundwater Monitoring Wells



AERIAL PHOTOGRAPHY PROVIDED BY  
 MICROSOFT BING MAPS  
 DIAGRAM IS FOR GENERAL LOCATION ONLY,  
 AND IS NOT INTENDED FOR CONSTRUCTION  
 PURPOSES

|                  |     |              |          |
|------------------|-----|--------------|----------|
| Project Manager: | MJS | Project No.: | 20187115 |
| Drawn by:        | JLB | Scale:       | AS SHOWN |
| Checked by:      | MJS | File Name:   | Site     |
| Approved by:     | JCG | Date:        | 12/5/18  |



**SITE DIAGRAM**  
 Union Reservoir Baseline Assessment  
 Weld County Road 28  
 Longmont, Colorado

|         |
|---------|
| Exhibit |
| 2       |

## **APPENDIX B – TABLES**

Table 1 – Soil Analytical Summary

Table 2 – Groundwater Analytical Summary



**Table 1**  
**Summary of Soil Analytical Results**  
**Union Reservoir Baseline Sampling**  
**Longmont, Colorado**  
**Terracon Project No. 22187053**

| Sample ID and Depth                 |                 |                |                            | SB-01 (5) | SB-02 (5)   | SB-03 (5)   |
|-------------------------------------|-----------------|----------------|----------------------------|-----------|-------------|-------------|
| Collection Date                     |                 |                |                            | 10/29/18  | 10/29/18    | 10/29/18    |
| Parameter                           | Residential RSL | Industrial RSL | COGCC Concentration Limits | mg/kg     | mg/kg       | mg/kg       |
| <b>Metals</b>                       |                 |                |                            |           |             |             |
| Arsenic                             | <b>0.68</b>     | <b>3.0</b>     | <b>0.39</b>                | <2        | <b>4.13</b> | <b>2.67</b> |
| Barium                              | <b>15,000</b>   | <b>220,000</b> | <b>15,000</b>              | 138       | 132         | 191         |
| Boron                               | NE              | NE             | NE                         | 12.3      | 14.3        | 17.7        |
| Chromium III                        | NE              | NE             | 120,000                    | 19.2      | 21          | 21.6        |
| Copper                              | <b>3,100</b>    | <b>47,000</b>  | <b>3,100</b>               | 9.81      | 10.5        | 10.8        |
| Lead                                | <b>400</b>      | <b>800</b>     | <b>400</b>                 | 11.8      | 13.1        | 12.4        |
| Selenium                            | <b>390</b>      | <b>5,800</b>   | <b>390</b>                 | <2        | 2.45        | 2.83        |
| Nickel                              | <b>1,500</b>    | <b>22,000</b>  | <b>1,600</b>               | 15.2      | 17.9        | 17.5        |
| Zinc                                | <b>23,000</b>   | <b>350,000</b> | <b>23,000</b>              | 47.9      | 53.1        | 54.2        |
| <b>Inorganics</b>                   |                 |                |                            |           |             |             |
| Sodium Adsorption Rate              | NE              | NE             | <12                        | 13.4      | 10.8        | 24.6        |
| Electrical Conductivity (mmhos/cm)* | NE              | NE             | <4                         | 3.1       | 2.28        | 9.6         |
| pH                                  | NE              | NE             | 6-9                        | 8.83      | 8.74        | 8.81        |

Only detected analytes shown (detected concentrations are **bold**)

RSL = EPA Regional Screening Level (November 2018)

NE = Not Established

mmhos/cm = millimhos per centimeter

\* Electrical Conductivity COGCC regulatory values at <4 mmhos/cm or 2 times established background values

COGCC = Colorado Oil and Gas Conservation Commission Table 910-1 (May 2018)

**Table 2**  
**Groundwater Analytical Summary**  
**Union Reservoir Baseline Sampling**  
**Longmont, Colorado**  
**Terracon Project No. 22187053**

| Sample ID                    |                                                                               |                                                                         |                                               | MW-01      | MW-02      | MW-03      |
|------------------------------|-------------------------------------------------------------------------------|-------------------------------------------------------------------------|-----------------------------------------------|------------|------------|------------|
| Collection Date              |                                                                               |                                                                         |                                               | 10/31/18   | 10/31/18   | 10/31/18   |
| Parameter                    | CDPHE Reg. 41<br>Groundwater -<br>Human Health<br>Drinking Water <sup>1</sup> | CDPHE Reg. 41<br>Groundwater -<br>Agricultural<br>Standard <sup>2</sup> | COGCC<br>Concentration<br>Levels <sup>3</sup> | µg/L       | µg/L       | µg/L       |
| <b>Metals</b>                |                                                                               |                                                                         |                                               |            |            |            |
| Barium                       | <b>2,000</b>                                                                  | NE                                                                      | NE                                            | 71.5       | 44.7       | 101        |
| Boron                        | NE                                                                            | 750                                                                     | NE                                            | 1,110      | 1,030      | 2,950      |
| Selenium                     | 50                                                                            | 20                                                                      | NE                                            | 35.8       | 465        | 1,240      |
| Copper                       | 1,000                                                                         | 200                                                                     | NE                                            | 16.7       | <10        | <10        |
| Nickel                       | 100                                                                           | 200                                                                     | NE                                            | 10.5       | <10        | <10        |
| <b>VOCs (8260B)</b>          |                                                                               |                                                                         |                                               |            |            |            |
| p-Isopropyltoluene           | NE                                                                            | NE                                                                      | NE                                            | 1.22       | <1         | <1         |
| <b>PAHs (8270)</b>           |                                                                               |                                                                         |                                               |            |            |            |
| Fluorene                     | 280                                                                           | NE                                                                      | NE                                            | 0.0583     | 0.0647     | <0.05      |
| Phenanthrene                 | NE                                                                            | NE                                                                      | NE                                            | <0.05      | 0.0827     | <0.05      |
| <b>Inorganic Parameters</b>  |                                                                               |                                                                         |                                               |            |            |            |
| Total Dissolved Solids (TDS) | <b>400,000-no limit</b>                                                       | NE                                                                      | NE                                            | 20,000,000 | 19,700,000 | 77,700,000 |
| Chloride                     | <b>250,000</b>                                                                | NE                                                                      | 76,210*                                       | 357,000    | 569,000    | 1,830,000  |
| Sulfate                      | <b>250,000</b>                                                                | NE                                                                      | 757,630*                                      | 15,500,000 | 14,800,000 | 50,300,000 |

1) CDPHE GW Quality Standards – Regulation 41 Table A, Ground Water Human Health Drinking Water Standards (December 2016)

2) CDPHE GW Quality Standards - Regulation 41 Table 3, Agricultural Standards

3) COGCC Concentration Levels = COGCC Table 910-1 (May 2018)

\* The COGCC cleanup standard for chloride and sulfate is 1.25 x background. Background concentrations from unimpacted wells were used to average and calculate background concentrations.

Only detected analytes shown (detected concentrations are **bold**)

NE = Not Established

VOC = Volatile Organic Compounds

COGCC = Colorado Oil and Gas Conservation Commission

**APPENDIX C – ANALYTICAL REPORTS AND CHAINS OF  
CUSTODY**

November 07, 2018

## Terracon Consultants, Inc - Longmont, CO

Sample Delivery Group: L1039710  
Samples Received: 10/31/2018  
Project Number: 22187053  
Description: Union Reservoir

Report To: Michael Skridulis  
1242 Bramwood Place  
Longmont, CO 80501

Entire Report Reviewed By:



Olivia Studebaker  
Project Manager

Results relate only to the items tested or calibrated and are reported as rounded values. This test report shall not be reproduced, except in full, without written approval of the laboratory. Where applicable, sampling conducted by Pace National is performed per guidance provided in laboratory standard operating procedures: 060302, 060303, and 060304.



|                                                             |           |
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| <b>Cp: Cover Page</b>                                       | <b>1</b>  |
| <b>Tc: Table of Contents</b>                                | <b>2</b>  |
| <b>Ss: Sample Summary</b>                                   | <b>3</b>  |
| <b>Cn: Case Narrative</b>                                   | <b>4</b>  |
| <b>Sr: Sample Results</b>                                   | <b>5</b>  |
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| SB-02 (5) L1039710-02                                       | 8         |
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| Wet Chemistry by Method 9045D                               | 16        |
| Wet Chemistry by Method 9050AMod                            | 17        |
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|                 |
|-----------------|
| <sup>1</sup> Cp |
| <sup>2</sup> Tc |
| <sup>3</sup> Ss |
| <sup>4</sup> Cn |
| <sup>5</sup> Sr |
| <sup>6</sup> Qc |
| <sup>7</sup> Gl |
| <sup>8</sup> Al |
| <sup>9</sup> Sc |

# SAMPLE SUMMARY



## SB-01 (5) L1039710-01 Solid

Collected by  
M. Skridulis  
Collected date/time  
10/29/18 12:20  
Received date/time  
10/31/18 08:45

| Method                                                      | Batch     | Dilution | Preparation date/time | Analysis date/time | Analyst |
|-------------------------------------------------------------|-----------|----------|-----------------------|--------------------|---------|
| Calculated Results                                          | WG1190126 | 1        | 11/04/18 21:47        | 11/05/18 17:23     | ST      |
| Calculated Results                                          | WG1189483 | 1        | 11/01/18 13:33        | 11/04/18 15:19     | ST      |
| Wet Chemistry by Method 3060A/7196A                         | WG1190085 | 1        | 11/02/18 10:46        | 11/02/18 16:00     | MLW     |
| Wet Chemistry by Method 9045D                               | WG1190008 | 1        | 11/01/18 16:55        | 11/01/18 17:54     | KBW     |
| Wet Chemistry by Method 9050AMod                            | WG1189493 | 1        | 11/02/18 19:07        | 11/03/18 15:00     | TCC     |
| Mercury by Method 7471A                                     | WG1189737 | 1        | 11/01/18 10:41        | 11/01/18 19:13     | TCT     |
| Metals (ICP) by Method 6010B                                | WG1189483 | 1        | 11/01/18 13:33        | 11/04/18 15:19     | ST      |
| Volatile Organic Compounds (GC) by Method 8015D/GRO         | WG1191554 | 1        | 11/01/18 09:33        | 11/06/18 03:03     | DWR     |
| Volatile Organic Compounds (GC/MS) by Method 8260B          | WG1190073 | 1        | 11/01/18 09:33        | 11/01/18 22:17     | JHH     |
| Semi-Volatile Organic Compounds (GC) by Method 8015         | WG1191819 | 1        | 11/05/18 20:55        | 11/06/18 15:52     | KME     |
| Semi Volatile Organic Compounds (GC/MS) by Method 8270C-SIM | WG1191152 | 1        | 11/04/18 10:56        | 11/04/18 21:29     | DMG     |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

## SB-02 (5) L1039710-02 Solid

Collected by  
M. Skridulis  
Collected date/time  
10/29/18 13:05  
Received date/time  
10/31/18 08:45

| Method                                                      | Batch     | Dilution | Preparation date/time | Analysis date/time | Analyst |
|-------------------------------------------------------------|-----------|----------|-----------------------|--------------------|---------|
| Calculated Results                                          | WG1190126 | 1        | 11/04/18 21:47        | 11/05/18 17:26     | ST      |
| Calculated Results                                          | WG1189483 | 1        | 11/01/18 13:33        | 11/04/18 15:22     | ST      |
| Wet Chemistry by Method 3060A/7196A                         | WG1190085 | 1        | 11/02/18 10:46        | 11/02/18 16:00     | MLW     |
| Wet Chemistry by Method 9045D                               | WG1190008 | 1        | 11/01/18 16:55        | 11/01/18 17:54     | KBW     |
| Wet Chemistry by Method 9050AMod                            | WG1189493 | 1        | 11/02/18 19:07        | 11/03/18 15:00     | TCC     |
| Mercury by Method 7471A                                     | WG1189737 | 1        | 11/01/18 10:41        | 11/01/18 19:24     | TCT     |
| Metals (ICP) by Method 6010B                                | WG1189483 | 1        | 11/01/18 13:33        | 11/04/18 15:22     | ST      |
| Volatile Organic Compounds (GC) by Method 8015D/GRO         | WG1191336 | 1        | 11/01/18 09:33        | 11/05/18 00:58     | JAH     |
| Volatile Organic Compounds (GC/MS) by Method 8260B          | WG1190073 | 1        | 11/01/18 09:33        | 11/01/18 22:37     | JHH     |
| Semi-Volatile Organic Compounds (GC) by Method 8015         | WG1191819 | 1        | 11/05/18 20:55        | 11/06/18 16:08     | KME     |
| Semi Volatile Organic Compounds (GC/MS) by Method 8270C-SIM | WG1191152 | 1        | 11/04/18 10:56        | 11/04/18 21:50     | DMG     |

## SB-03 (5) L1039710-03 Solid

Collected by  
M. Skridulis  
Collected date/time  
10/29/18 13:50  
Received date/time  
10/31/18 08:45

| Method                                                      | Batch     | Dilution | Preparation date/time | Analysis date/time | Analyst |
|-------------------------------------------------------------|-----------|----------|-----------------------|--------------------|---------|
| Calculated Results                                          | WG1190126 | 1        | 11/04/18 21:47        | 11/05/18 22:30     | ST      |
| Calculated Results                                          | WG1189483 | 1        | 11/01/18 13:33        | 11/04/18 15:25     | ST      |
| Wet Chemistry by Method 3060A/7196A                         | WG1190085 | 1        | 11/02/18 10:46        | 11/02/18 16:00     | MLW     |
| Wet Chemistry by Method 9045D                               | WG1190008 | 1        | 11/01/18 16:55        | 11/01/18 17:54     | KBW     |
| Wet Chemistry by Method 9050AMod                            | WG1189493 | 1        | 11/02/18 19:07        | 11/03/18 15:00     | TCC     |
| Mercury by Method 7471A                                     | WG1189737 | 1        | 11/01/18 10:41        | 11/01/18 19:26     | TCT     |
| Metals (ICP) by Method 6010B                                | WG1189483 | 1        | 11/01/18 13:33        | 11/04/18 15:25     | ST      |
| Volatile Organic Compounds (GC) by Method 8015D/GRO         | WG1191336 | 1        | 11/01/18 09:33        | 11/05/18 01:20     | JAH     |
| Volatile Organic Compounds (GC/MS) by Method 8260B          | WG1190073 | 1        | 11/01/18 09:33        | 11/01/18 22:57     | JHH     |
| Volatile Organic Compounds (GC/MS) by Method 8260B          | WG1192300 | 1        | 11/01/18 09:33        | 11/06/18 23:54     | JAH     |
| Semi-Volatile Organic Compounds (GC) by Method 8015         | WG1191819 | 1        | 11/05/18 20:55        | 11/06/18 16:23     | KME     |
| Semi Volatile Organic Compounds (GC/MS) by Method 8270C-SIM | WG1191152 | 1        | 11/04/18 10:56        | 11/04/18 22:11     | DMG     |



All sample aliquots were received at the correct temperature, in the proper containers, with the appropriate preservatives, and within method specified holding times, unless qualified or notated within the report. Where applicable, all MDL (LOD) and RDL (LOQ) values reported for environmental samples have been corrected for the dilution factor used in the analysis. All Method and Batch Quality Control are within established criteria except where addressed in this case narrative, a non-conformance form or properly qualified within the sample results. By my digital signature below, I affirm to the best of my knowledge, all problems/anomalies observed by the laboratory as having the potential to affect the quality of the data have been identified by the laboratory, and no information or data have been knowingly withheld that would affect the quality of the data.

Olivia Studebaker  
Project Manager

- <sup>1</sup> Cp
- <sup>2</sup> Tc
- <sup>3</sup> Ss
- <sup>4</sup> Cn
- <sup>5</sup> Sr
- <sup>6</sup> Qc
- <sup>7</sup> Gl
- <sup>8</sup> Al
- <sup>9</sup> Sc



Calculated Results

| Analyte                 | Result | Qualifier | Dilution | Analysis date / time | Batch     |
|-------------------------|--------|-----------|----------|----------------------|-----------|
| Sodium Adsorption Ratio | 13.4   |           | 1        | 11/05/2018 17:23     | WG1190126 |

1 Cp

2 Tc

Calculated Results

| Analyte             | Result mg/kg | Qualifier | RDL mg/kg | Dilution | Analysis date / time | Batch                     |
|---------------------|--------------|-----------|-----------|----------|----------------------|---------------------------|
| Chromium, Trivalent | 19.2         |           | 1.00      | 1        | 11/04/2018 15:19     | <a href="#">WG1189483</a> |

3 Ss

4 Cn

Wet Chemistry by Method 3060A/7196A

| Analyte              | Result mg/kg | Qualifier | RDL mg/kg | Dilution | Analysis date / time | Batch                     |
|----------------------|--------------|-----------|-----------|----------|----------------------|---------------------------|
| Chromium, Hexavalent | ND           |           | 2.00      | 1        | 11/02/2018 16:00     | <a href="#">WG1190085</a> |

5 Sr

6 Qc

Wet Chemistry by Method 9045D

| Analyte | Result su | Qualifier | Dilution | Analysis date / time | Batch                     |
|---------|-----------|-----------|----------|----------------------|---------------------------|
| pH      | 8.83      | <u>T8</u> | 1        | 11/01/2018 17:54     | <a href="#">WG1190008</a> |

7 Gl

8 Al

Sample Narrative:

L1039710-01 WG1190008: 8.83 at 20.3C

9 Sc

Wet Chemistry by Method 9050AMod

| Analyte              | Result umhos/cm | Qualifier | RDL umhos/cm | Dilution | Analysis date / time | Batch                     |
|----------------------|-----------------|-----------|--------------|----------|----------------------|---------------------------|
| Specific Conductance | 3100            |           | 10.0         | 1        | 11/03/2018 15:00     | <a href="#">WG1189493</a> |

Mercury by Method 7471A

| Analyte | Result mg/kg | Qualifier | RDL mg/kg | Dilution | Analysis date / time | Batch                     |
|---------|--------------|-----------|-----------|----------|----------------------|---------------------------|
| Mercury | ND           |           | 0.0200    | 1        | 11/01/2018 19:13     | <a href="#">WG1189737</a> |

Metals (ICP) by Method 6010B

| Analyte  | Result mg/kg | Qualifier | RDL mg/kg | Dilution | Analysis date / time | Batch                     |
|----------|--------------|-----------|-----------|----------|----------------------|---------------------------|
| Arsenic  | ND           |           | 2.00      | 1        | 11/04/2018 15:19     | <a href="#">WG1189483</a> |
| Barium   | 138          |           | 0.500     | 1        | 11/04/2018 15:19     | <a href="#">WG1189483</a> |
| Boron    | 12.3         |           | 10.0      | 1        | 11/04/2018 15:19     | <a href="#">WG1189483</a> |
| Cadmium  | ND           |           | 0.500     | 1        | 11/04/2018 15:19     | <a href="#">WG1189483</a> |
| Chromium | 19.2         |           | 1.00      | 1        | 11/04/2018 15:19     | <a href="#">WG1189483</a> |
| Copper   | 9.81         |           | 2.00      | 1        | 11/04/2018 15:19     | <a href="#">WG1189483</a> |
| Lead     | 11.8         |           | 0.500     | 1        | 11/04/2018 15:19     | <a href="#">WG1189483</a> |
| Nickel   | 15.2         |           | 2.00      | 1        | 11/04/2018 15:19     | <a href="#">WG1189483</a> |
| Selenium | ND           |           | 2.00      | 1        | 11/04/2018 15:19     | <a href="#">WG1189483</a> |
| Silver   | ND           |           | 1.00      | 1        | 11/04/2018 15:19     | <a href="#">WG1189483</a> |
| Zinc     | 47.9         |           | 5.00      | 1        | 11/04/2018 15:19     | <a href="#">WG1189483</a> |

Volatile Organic Compounds (GC) by Method 8015D/GRO

| Analyte                           | Result mg/kg | Qualifier | RDL mg/kg | Dilution | Analysis date / time | Batch                     |
|-----------------------------------|--------------|-----------|-----------|----------|----------------------|---------------------------|
| TPH (GC/FID) Low Fraction         | ND           |           | 0.100     | 1        | 11/06/2018 03:03     | <a href="#">WG1191554</a> |
| (S) a, a, a-Trifluorotoluene(FID) | 98.7         |           | 77.0-120  |          | 11/06/2018 03:03     | <a href="#">WG1191554</a> |





Collected date/time: 10/29/18 12:20

L1039710

## Volatile Organic Compounds (GC/MS) by Method 8260B

| Analyte                        | Result<br>mg/kg | Qualifier | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch     |
|--------------------------------|-----------------|-----------|--------------|----------|-------------------------|-----------|
| Acetone                        | ND              |           | 0.0250       | 1        | 11/01/2018 22:17        | WG1190073 |
| Acrylonitrile                  | ND              |           | 0.0125       | 1        | 11/01/2018 22:17        | WG1190073 |
| Benzene                        | ND              |           | 0.00100      | 1        | 11/01/2018 22:17        | WG1190073 |
| Bromobenzene                   | ND              |           | 0.0125       | 1        | 11/01/2018 22:17        | WG1190073 |
| Bromodichloromethane           | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| Bromoform                      | ND              |           | 0.0250       | 1        | 11/01/2018 22:17        | WG1190073 |
| Bromomethane                   | ND              |           | 0.0125       | 1        | 11/01/2018 22:17        | WG1190073 |
| n-Butylbenzene                 | ND              |           | 0.0125       | 1        | 11/01/2018 22:17        | WG1190073 |
| sec-Butylbenzene               | ND              |           | 0.0125       | 1        | 11/01/2018 22:17        | WG1190073 |
| tert-Butylbenzene              | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| Carbon tetrachloride           | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| Chlorobenzene                  | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| Chlorodibromomethane           | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| Chloroethane                   | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| Chloroform                     | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| Chloromethane                  | ND              |           | 0.0125       | 1        | 11/01/2018 22:17        | WG1190073 |
| 2-Chlorotoluene                | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| 4-Chlorotoluene                | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,2-Dibromo-3-Chloropropane    | ND              |           | 0.0250       | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,2-Dibromoethane              | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| Dibromomethane                 | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,2-Dichlorobenzene            | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,3-Dichlorobenzene            | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,4-Dichlorobenzene            | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| Dichlorodifluoromethane        | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,1-Dichloroethane             | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,2-Dichloroethane             | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,1-Dichloroethene             | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| cis-1,2-Dichloroethene         | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| trans-1,2-Dichloroethene       | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,2-Dichloropropane            | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,1-Dichloropropene            | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,3-Dichloropropane            | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| cis-1,3-Dichloropropene        | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| trans-1,3-Dichloropropene      | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| 2,2-Dichloropropane            | ND              | J4        | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| Di-isopropyl ether             | ND              |           | 0.00100      | 1        | 11/01/2018 22:17        | WG1190073 |
| Ethylbenzene                   | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| Hexachloro-1,3-butadiene       | ND              |           | 0.0250       | 1        | 11/01/2018 22:17        | WG1190073 |
| Isopropylbenzene               | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| p-Isopropyltoluene             | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| 2-Butanone (MEK)               | ND              |           | 0.0250       | 1        | 11/01/2018 22:17        | WG1190073 |
| Methylene Chloride             | ND              |           | 0.0250       | 1        | 11/01/2018 22:17        | WG1190073 |
| 4-Methyl-2-pentanone (MIBK)    | ND              |           | 0.0250       | 1        | 11/01/2018 22:17        | WG1190073 |
| Methyl tert-butyl ether        | ND              | J4        | 0.00100      | 1        | 11/01/2018 22:17        | WG1190073 |
| Naphthalene                    | ND              |           | 0.0125       | 1        | 11/01/2018 22:17        | WG1190073 |
| n-Propylbenzene                | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| Styrene                        | ND              |           | 0.0125       | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,1,1,2-Tetrachloroethane      | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,1,2,2-Tetrachloroethane      | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,1,2-Trichlorotrifluoroethane | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| Tetrachloroethene              | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| Toluene                        | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,2,3-Trichlorobenzene         | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,2,4-Trichlorobenzene         | ND              |           | 0.0125       | 1        | 11/01/2018 22:17        | WG1190073 |
| 1,1,1-Trichloroethane          | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | WG1190073 |

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



Collected date/time: 10/29/18 12:20

L1039710

Volatile Organic Compounds (GC/MS) by Method 8260B

| Analyte                  | Result<br>mg/kg | Qualifier | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch                     |
|--------------------------|-----------------|-----------|--------------|----------|-------------------------|---------------------------|
| 1,1,2-Trichloroethane    | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | <a href="#">WG1190073</a> |
| Trichloroethene          | ND              |           | 0.00100      | 1        | 11/01/2018 22:17        | <a href="#">WG1190073</a> |
| Trichlorofluoromethane   | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | <a href="#">WG1190073</a> |
| 1,2,3-Trichloropropane   | ND              |           | 0.0125       | 1        | 11/01/2018 22:17        | <a href="#">WG1190073</a> |
| 1,2,4-Trimethylbenzene   | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | <a href="#">WG1190073</a> |
| 1,2,3-Trimethylbenzene   | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | <a href="#">WG1190073</a> |
| 1,3,5-Trimethylbenzene   | ND              |           | 0.00500      | 1        | 11/01/2018 22:17        | <a href="#">WG1190073</a> |
| Vinyl chloride           | ND              |           | 0.00250      | 1        | 11/01/2018 22:17        | <a href="#">WG1190073</a> |
| Xylenes, Total           | ND              |           | 0.00650      | 1        | 11/01/2018 22:17        | <a href="#">WG1190073</a> |
| (S) Toluene-d8           | 121             |           | 75.0-131     |          | 11/01/2018 22:17        | <a href="#">WG1190073</a> |
| (S) Dibromofluoromethane | 99.4            |           | 65.0-129     |          | 11/01/2018 22:17        | <a href="#">WG1190073</a> |
| (S) 4-Bromofluorobenzene | 80.4            |           | 67.0-138     |          | 11/01/2018 22:17        | <a href="#">WG1190073</a> |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Semi-Volatile Organic Compounds (GC) by Method 8015

| Analyte              | Result<br>mg/kg | Qualifier | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch                     |
|----------------------|-----------------|-----------|--------------|----------|-------------------------|---------------------------|
| C10-C28 Diesel Range | ND              |           | 4.00         | 1        | 11/06/2018 15:52        | <a href="#">WG1191819</a> |
| C28-C40 Oil Range    | ND              |           | 4.00         | 1        | 11/06/2018 15:52        | <a href="#">WG1191819</a> |
| (S) o-Terphenyl      | 69.8            |           | 18.0-148     |          | 11/06/2018 15:52        | <a href="#">WG1191819</a> |

Semi Volatile Organic Compounds (GC/MS) by Method 8270C-SIM

| Analyte                | Result<br>mg/kg | Qualifier | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch                     |
|------------------------|-----------------|-----------|--------------|----------|-------------------------|---------------------------|
| Anthracene             | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Acenaphthene           | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Acenaphthylene         | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Benzo(a)anthracene     | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Benzo(a)pyrene         | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Benzo(b)fluoranthene   | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Benzo(g,h,i)perylene   | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Benzo(k)fluoranthene   | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Chrysene               | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Dibenz(a,h)anthracene  | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Fluoranthene           | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Fluorene               | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Indeno(1,2,3-cd)pyrene | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Naphthalene            | ND              |           | 0.0200       | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Phenanthrene           | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| Pyrene                 | ND              |           | 0.00600      | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| 1-Methylnaphthalene    | ND              |           | 0.0200       | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| 2-Methylnaphthalene    | ND              |           | 0.0200       | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| 2-Chloronaphthalene    | ND              |           | 0.0200       | 1        | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| (S) p-Terphenyl-d14    | 71.6            |           | 23.0-120     |          | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| (S) Nitrobenzene-d5    | 88.5            |           | 14.0-149     |          | 11/04/2018 21:29        | <a href="#">WG1191152</a> |
| (S) 2-Fluorobiphenyl   | 79.9            |           | 34.0-125     |          | 11/04/2018 21:29        | <a href="#">WG1191152</a> |



Calculated Results

| Analyte                 | Result | Qualifier | Dilution | Analysis date / time | Batch     |
|-------------------------|--------|-----------|----------|----------------------|-----------|
| Sodium Adsorption Ratio | 10.8   |           | 1        | 11/05/2018 17:26     | WG1190126 |

1 Cp

2 Tc

Calculated Results

| Analyte             | Result | Qualifier | RDL  | Dilution | Analysis date / time | Batch                     |
|---------------------|--------|-----------|------|----------|----------------------|---------------------------|
| Chromium, Trivalent | 21.0   |           | 1.00 | 1        | 11/04/2018 15:22     | <a href="#">WG1189483</a> |

3 Ss

4 Cn

Wet Chemistry by Method 3060A/7196A

| Analyte              | Result | Qualifier | RDL  | Dilution | Analysis date / time | Batch                     |
|----------------------|--------|-----------|------|----------|----------------------|---------------------------|
| Chromium, Hexavalent | ND     |           | 2.00 | 1        | 11/02/2018 16:00     | <a href="#">WG1190085</a> |

5 Sr

6 Qc

Wet Chemistry by Method 9045D

| Analyte | Result | Qualifier | Dilution | Analysis date / time | Batch                     |
|---------|--------|-----------|----------|----------------------|---------------------------|
| pH      | 8.74   | <u>T8</u> | 1        | 11/01/2018 17:54     | <a href="#">WG1190008</a> |

7 Gl

8 Al

Sample Narrative:

L1039710-02 WG1190008: 8.74 at 20.3C

9 Sc

Wet Chemistry by Method 9050AMod

| Analyte              | Result | Qualifier | RDL  | Dilution | Analysis date / time | Batch                     |
|----------------------|--------|-----------|------|----------|----------------------|---------------------------|
| Specific Conductance | 2280   |           | 10.0 | 1        | 11/03/2018 15:00     | <a href="#">WG1189493</a> |

Mercury by Method 7471A

| Analyte | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|---------|--------|-----------|--------|----------|----------------------|---------------------------|
| Mercury | ND     |           | 0.0200 | 1        | 11/01/2018 19:24     | <a href="#">WG1189737</a> |

Metals (ICP) by Method 6010B

| Analyte  | Result | Qualifier | RDL   | Dilution | Analysis date / time | Batch                     |
|----------|--------|-----------|-------|----------|----------------------|---------------------------|
| Arsenic  | 4.13   |           | 2.00  | 1        | 11/04/2018 15:22     | <a href="#">WG1189483</a> |
| Barium   | 132    |           | 0.500 | 1        | 11/04/2018 15:22     | <a href="#">WG1189483</a> |
| Boron    | 14.3   |           | 10.0  | 1        | 11/04/2018 15:22     | <a href="#">WG1189483</a> |
| Cadmium  | ND     |           | 0.500 | 1        | 11/04/2018 15:22     | <a href="#">WG1189483</a> |
| Chromium | 21.0   |           | 1.00  | 1        | 11/04/2018 15:22     | <a href="#">WG1189483</a> |
| Copper   | 10.5   |           | 2.00  | 1        | 11/04/2018 15:22     | <a href="#">WG1189483</a> |
| Lead     | 13.1   |           | 0.500 | 1        | 11/04/2018 15:22     | <a href="#">WG1189483</a> |
| Nickel   | 17.9   |           | 2.00  | 1        | 11/04/2018 15:22     | <a href="#">WG1189483</a> |
| Selenium | 2.45   |           | 2.00  | 1        | 11/04/2018 15:22     | <a href="#">WG1189483</a> |
| Silver   | ND     |           | 1.00  | 1        | 11/04/2018 15:22     | <a href="#">WG1189483</a> |
| Zinc     | 53.1   |           | 5.00  | 1        | 11/04/2018 15:22     | <a href="#">WG1189483</a> |

Volatile Organic Compounds (GC) by Method 8015D/GRO

| Analyte                           | Result | Qualifier | RDL      | Dilution | Analysis date / time | Batch                     |
|-----------------------------------|--------|-----------|----------|----------|----------------------|---------------------------|
| TPH (GC/FID) Low Fraction         | ND     |           | 0.100    | 1        | 11/05/2018 00:58     | <a href="#">WG1191336</a> |
| (S) a, a, a-Trifluorotoluene(FID) | 94.7   |           | 77.0-120 |          | 11/05/2018 00:58     | <a href="#">WG1191336</a> |



Volatile Organic Compounds (GC/MS) by Method 8260B

| Analyte                        | Result<br>mg/kg | Qualifier | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch     |
|--------------------------------|-----------------|-----------|--------------|----------|-------------------------|-----------|
| Acetone                        | ND              |           | 0.0250       | 1        | 11/01/2018 22:37        | WG1190073 |
| Acrylonitrile                  | ND              |           | 0.0125       | 1        | 11/01/2018 22:37        | WG1190073 |
| Benzene                        | ND              |           | 0.00100      | 1        | 11/01/2018 22:37        | WG1190073 |
| Bromobenzene                   | ND              |           | 0.0125       | 1        | 11/01/2018 22:37        | WG1190073 |
| Bromodichloromethane           | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| Bromoform                      | ND              |           | 0.0250       | 1        | 11/01/2018 22:37        | WG1190073 |
| Bromomethane                   | ND              |           | 0.0125       | 1        | 11/01/2018 22:37        | WG1190073 |
| n-Butylbenzene                 | ND              |           | 0.0125       | 1        | 11/01/2018 22:37        | WG1190073 |
| sec-Butylbenzene               | ND              |           | 0.0125       | 1        | 11/01/2018 22:37        | WG1190073 |
| tert-Butylbenzene              | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| Carbon tetrachloride           | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| Chlorobenzene                  | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| Chlorodibromomethane           | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| Chloroethane                   | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| Chloroform                     | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| Chloromethane                  | ND              |           | 0.0125       | 1        | 11/01/2018 22:37        | WG1190073 |
| 2-Chlorotoluene                | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| 4-Chlorotoluene                | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,2-Dibromo-3-Chloropropane    | ND              |           | 0.0250       | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,2-Dibromoethane              | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| Dibromomethane                 | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,2-Dichlorobenzene            | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,3-Dichlorobenzene            | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,4-Dichlorobenzene            | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| Dichlorodifluoromethane        | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,1-Dichloroethane             | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,2-Dichloroethane             | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,1-Dichloroethene             | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| cis-1,2-Dichloroethene         | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| trans-1,2-Dichloroethene       | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,2-Dichloropropane            | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,1-Dichloropropene            | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,3-Dichloropropane            | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| cis-1,3-Dichloropropene        | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| trans-1,3-Dichloropropene      | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| 2,2-Dichloropropane            | ND              | J4        | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| Di-isopropyl ether             | ND              |           | 0.00100      | 1        | 11/01/2018 22:37        | WG1190073 |
| Ethylbenzene                   | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| Hexachloro-1,3-butadiene       | ND              |           | 0.0250       | 1        | 11/01/2018 22:37        | WG1190073 |
| Isopropylbenzene               | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| p-Isopropyltoluene             | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| 2-Butanone (MEK)               | ND              |           | 0.0250       | 1        | 11/01/2018 22:37        | WG1190073 |
| Methylene Chloride             | ND              |           | 0.0250       | 1        | 11/01/2018 22:37        | WG1190073 |
| 4-Methyl-2-pentanone (MIBK)    | ND              |           | 0.0250       | 1        | 11/01/2018 22:37        | WG1190073 |
| Methyl tert-butyl ether        | ND              | J4        | 0.00100      | 1        | 11/01/2018 22:37        | WG1190073 |
| Naphthalene                    | ND              |           | 0.0125       | 1        | 11/01/2018 22:37        | WG1190073 |
| n-Propylbenzene                | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| Styrene                        | ND              |           | 0.0125       | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,1,1,2-Tetrachloroethane      | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,1,2,2-Tetrachloroethane      | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,1,2-Trichlorotrifluoroethane | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| Tetrachloroethene              | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| Toluene                        | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,2,3-Trichlorobenzene         | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,2,4-Trichlorobenzene         | ND              |           | 0.0125       | 1        | 11/01/2018 22:37        | WG1190073 |
| 1,1,1-Trichloroethane          | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | WG1190073 |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc



Volatile Organic Compounds (GC/MS) by Method 8260B

| Analyte                  | Result<br>mg/kg | Qualifier | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch                     |
|--------------------------|-----------------|-----------|--------------|----------|-------------------------|---------------------------|
| 1,1,2-Trichloroethane    | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | <a href="#">WG1190073</a> |
| Trichloroethene          | ND              |           | 0.00100      | 1        | 11/01/2018 22:37        | <a href="#">WG1190073</a> |
| Trichlorofluoromethane   | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | <a href="#">WG1190073</a> |
| 1,2,3-Trichloropropane   | ND              |           | 0.0125       | 1        | 11/01/2018 22:37        | <a href="#">WG1190073</a> |
| 1,2,4-Trimethylbenzene   | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | <a href="#">WG1190073</a> |
| 1,2,3-Trimethylbenzene   | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | <a href="#">WG1190073</a> |
| 1,3,5-Trimethylbenzene   | ND              |           | 0.00500      | 1        | 11/01/2018 22:37        | <a href="#">WG1190073</a> |
| Vinyl chloride           | ND              |           | 0.00250      | 1        | 11/01/2018 22:37        | <a href="#">WG1190073</a> |
| Xylenes, Total           | ND              |           | 0.00650      | 1        | 11/01/2018 22:37        | <a href="#">WG1190073</a> |
| (S) Toluene-d8           | 124             |           | 75.0-131     |          | 11/01/2018 22:37        | <a href="#">WG1190073</a> |
| (S) Dibromofluoromethane | 101             |           | 65.0-129     |          | 11/01/2018 22:37        | <a href="#">WG1190073</a> |
| (S) 4-Bromofluorobenzene | 84.6            |           | 67.0-138     |          | 11/01/2018 22:37        | <a href="#">WG1190073</a> |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Semi-Volatile Organic Compounds (GC) by Method 8015

| Analyte              | Result<br>mg/kg | Qualifier | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch                     |
|----------------------|-----------------|-----------|--------------|----------|-------------------------|---------------------------|
| C10-C28 Diesel Range | ND              |           | 4.00         | 1        | 11/06/2018 16:08        | <a href="#">WG1191819</a> |
| C28-C40 Oil Range    | ND              |           | 4.00         | 1        | 11/06/2018 16:08        | <a href="#">WG1191819</a> |
| (S) o-Terphenyl      | 61.8            |           | 18.0-148     |          | 11/06/2018 16:08        | <a href="#">WG1191819</a> |

Semi Volatile Organic Compounds (GC/MS) by Method 8270C-SIM

| Analyte                | Result<br>mg/kg | Qualifier | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch                     |
|------------------------|-----------------|-----------|--------------|----------|-------------------------|---------------------------|
| Anthracene             | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Acenaphthene           | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Acenaphthylene         | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Benzo(a)anthracene     | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Benzo(a)pyrene         | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Benzo(b)fluoranthene   | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Benzo(g,h,i)perylene   | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Benzo(k)fluoranthene   | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Chrysene               | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Dibenz(a,h)anthracene  | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Fluoranthene           | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Fluorene               | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Indeno(1,2,3-cd)pyrene | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Naphthalene            | ND              |           | 0.0200       | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Phenanthrene           | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| Pyrene                 | ND              |           | 0.00600      | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| 1-Methylnaphthalene    | ND              |           | 0.0200       | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| 2-Methylnaphthalene    | ND              |           | 0.0200       | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| 2-Chloronaphthalene    | ND              |           | 0.0200       | 1        | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| (S) p-Terphenyl-d14    | 72.8            |           | 23.0-120     |          | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| (S) Nitrobenzene-d5    | 86.1            |           | 14.0-149     |          | 11/04/2018 21:50        | <a href="#">WG1191152</a> |
| (S) 2-Fluorobiphenyl   | 79.2            |           | 34.0-125     |          | 11/04/2018 21:50        | <a href="#">WG1191152</a> |



Calculated Results

| Analyte                 | Result | Qualifier | Dilution | Analysis date / time | Batch     |
|-------------------------|--------|-----------|----------|----------------------|-----------|
| Sodium Adsorption Ratio | 24.6   |           | 1        | 11/05/2018 22:30     | WG1190126 |

1 Cp

2 Tc

Calculated Results

| Analyte             | Result | Qualifier | RDL  | Dilution | Analysis date / time | Batch                     |
|---------------------|--------|-----------|------|----------|----------------------|---------------------------|
| Chromium, Trivalent | 21.6   |           | 1.00 | 1        | 11/04/2018 15:25     | <a href="#">WG1189483</a> |

3 Ss

4 Cn

Wet Chemistry by Method 3060A/7196A

| Analyte              | Result | Qualifier | RDL  | Dilution | Analysis date / time | Batch                     |
|----------------------|--------|-----------|------|----------|----------------------|---------------------------|
| Chromium, Hexavalent | ND     |           | 2.00 | 1        | 11/02/2018 16:00     | <a href="#">WG1190085</a> |

5 Sr

6 Qc

Wet Chemistry by Method 9045D

| Analyte | Result | Qualifier | Dilution | Analysis date / time | Batch                     |
|---------|--------|-----------|----------|----------------------|---------------------------|
| pH      | 8.81   | <u>T8</u> | 1        | 11/01/2018 17:54     | <a href="#">WG1190008</a> |

7 Gl

8 Al

Sample Narrative:

L1039710-03 WG1190008: 8.81 at 20.2C

9 Sc

Wet Chemistry by Method 9050AMod

| Analyte              | Result | Qualifier | RDL  | Dilution | Analysis date / time | Batch                     |
|----------------------|--------|-----------|------|----------|----------------------|---------------------------|
| Specific Conductance | 9600   |           | 10.0 | 1        | 11/03/2018 15:00     | <a href="#">WG1189493</a> |

Mercury by Method 7471A

| Analyte | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|---------|--------|-----------|--------|----------|----------------------|---------------------------|
| Mercury | ND     |           | 0.0200 | 1        | 11/01/2018 19:26     | <a href="#">WG1189737</a> |

Metals (ICP) by Method 6010B

| Analyte  | Result | Qualifier | RDL   | Dilution | Analysis date / time | Batch                     |
|----------|--------|-----------|-------|----------|----------------------|---------------------------|
| Arsenic  | 2.67   |           | 2.00  | 1        | 11/04/2018 15:25     | <a href="#">WG1189483</a> |
| Barium   | 191    |           | 0.500 | 1        | 11/04/2018 15:25     | <a href="#">WG1189483</a> |
| Boron    | 17.7   |           | 10.0  | 1        | 11/04/2018 15:25     | <a href="#">WG1189483</a> |
| Cadmium  | ND     |           | 0.500 | 1        | 11/04/2018 15:25     | <a href="#">WG1189483</a> |
| Chromium | 21.6   |           | 1.00  | 1        | 11/04/2018 15:25     | <a href="#">WG1189483</a> |
| Copper   | 10.8   |           | 2.00  | 1        | 11/04/2018 15:25     | <a href="#">WG1189483</a> |
| Lead     | 12.4   |           | 0.500 | 1        | 11/04/2018 15:25     | <a href="#">WG1189483</a> |
| Nickel   | 17.5   |           | 2.00  | 1        | 11/04/2018 15:25     | <a href="#">WG1189483</a> |
| Selenium | 2.83   |           | 2.00  | 1        | 11/04/2018 15:25     | <a href="#">WG1189483</a> |
| Silver   | ND     |           | 1.00  | 1        | 11/04/2018 15:25     | <a href="#">WG1189483</a> |
| Zinc     | 54.2   |           | 5.00  | 1        | 11/04/2018 15:25     | <a href="#">WG1189483</a> |

Volatile Organic Compounds (GC) by Method 8015D/GRO

| Analyte                           | Result | Qualifier | RDL      | Dilution | Analysis date / time | Batch                     |
|-----------------------------------|--------|-----------|----------|----------|----------------------|---------------------------|
| TPH (GC/FID) Low Fraction         | ND     |           | 0.100    | 1        | 11/05/2018 01:20     | <a href="#">WG1191336</a> |
| (S) a, a, a-Trifluorotoluene(FID) | 94.1   |           | 77.0-120 |          | 11/05/2018 01:20     | <a href="#">WG1191336</a> |



Volatile Organic Compounds (GC/MS) by Method 8260B

| Analyte                        | Result<br>mg/kg | Qualifier | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch                     |
|--------------------------------|-----------------|-----------|--------------|----------|-------------------------|---------------------------|
| Acetone                        | ND              |           | 0.0250       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Acrylonitrile                  | ND              |           | 0.0125       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Benzene                        | ND              |           | 0.00100      | 1        | 11/06/2018 23:54        | <a href="#">WG1192300</a> |
| Bromobenzene                   | ND              |           | 0.0125       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Bromodichloromethane           | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Bromoform                      | ND              |           | 0.0250       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Bromomethane                   | ND              |           | 0.0125       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| n-Butylbenzene                 | ND              |           | 0.0125       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| sec-Butylbenzene               | ND              |           | 0.0125       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| tert-Butylbenzene              | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Carbon tetrachloride           | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Chlorobenzene                  | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Chlorodibromomethane           | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Chloroethane                   | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Chloroform                     | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Chloromethane                  | ND              |           | 0.0125       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 2-Chlorotoluene                | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 4-Chlorotoluene                | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,2-Dibromo-3-Chloropropane    | ND              |           | 0.0250       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,2-Dibromoethane              | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Dibromomethane                 | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,2-Dichlorobenzene            | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,3-Dichlorobenzene            | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,4-Dichlorobenzene            | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Dichlorodifluoromethane        | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,1-Dichloroethane             | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,2-Dichloroethane             | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,1-Dichloroethene             | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| cis-1,2-Dichloroethene         | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| trans-1,2-Dichloroethene       | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,2-Dichloropropane            | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,1-Dichloropropene            | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,3-Dichloropropane            | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| cis-1,3-Dichloropropene        | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| trans-1,3-Dichloropropene      | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 2,2-Dichloropropane            | ND              | J4        | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Di-isopropyl ether             | ND              |           | 0.00100      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Ethylbenzene                   | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Hexachloro-1,3-butadiene       | ND              |           | 0.0250       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Isopropylbenzene               | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| p-Isopropyltoluene             | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 2-Butanone (MEK)               | ND              |           | 0.0250       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Methylene Chloride             | ND              |           | 0.0250       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 4-Methyl-2-pentanone (MIBK)    | ND              |           | 0.0250       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Methyl tert-butyl ether        | ND              | J4        | 0.00100      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Naphthalene                    | ND              |           | 0.0125       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| n-Propylbenzene                | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Styrene                        | ND              |           | 0.0125       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,1,1,2-Tetrachloroethane      | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,1,2,2-Tetrachloroethane      | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,1,2-Trichlorotrifluoroethane | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Tetrachloroethene              | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Toluene                        | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,2,3-Trichlorobenzene         | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,2,4-Trichlorobenzene         | ND              |           | 0.0125       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,1,1-Trichloroethane          | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc



Collected date/time: 10/29/18 13:50

L1039710

Volatile Organic Compounds (GC/MS) by Method 8260B

| Analyte                  | Result<br>mg/kg | Qualifier | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch                     |
|--------------------------|-----------------|-----------|--------------|----------|-------------------------|---------------------------|
| 1,1,2-Trichloroethane    | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Trichloroethene          | ND              |           | 0.00100      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Trichlorofluoromethane   | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,2,3-Trichloropropane   | ND              |           | 0.0125       | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,2,4-Trimethylbenzene   | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,2,3-Trimethylbenzene   | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| 1,3,5-Trimethylbenzene   | ND              |           | 0.00500      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Vinyl chloride           | ND              |           | 0.00250      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| Xylenes, Total           | ND              |           | 0.00650      | 1        | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| (S) Toluene-d8           | 126             |           | 75.0-131     |          | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| (S) Toluene-d8           | 97.2            |           | 75.0-131     |          | 11/06/2018 23:54        | <a href="#">WG1192300</a> |
| (S) Dibromofluoromethane | 99.4            |           | 65.0-129     |          | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| (S) Dibromofluoromethane | 122             |           | 65.0-129     |          | 11/06/2018 23:54        | <a href="#">WG1192300</a> |
| (S) 4-Bromofluorobenzene | 84.1            |           | 67.0-138     |          | 11/01/2018 22:57        | <a href="#">WG1190073</a> |
| (S) 4-Bromofluorobenzene | 104             |           | 67.0-138     |          | 11/06/2018 23:54        | <a href="#">WG1192300</a> |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Semi-Volatile Organic Compounds (GC) by Method 8015

| Analyte              | Result<br>mg/kg | Qualifier | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch                     |
|----------------------|-----------------|-----------|--------------|----------|-------------------------|---------------------------|
| C10-C28 Diesel Range | ND              |           | 4.00         | 1        | 11/06/2018 16:23        | <a href="#">WG1191819</a> |
| C28-C40 Oil Range    | ND              |           | 4.00         | 1        | 11/06/2018 16:23        | <a href="#">WG1191819</a> |
| (S) o-Terphenyl      | 79.4            |           | 18.0-148     |          | 11/06/2018 16:23        | <a href="#">WG1191819</a> |

Semi Volatile Organic Compounds (GC/MS) by Method 8270C-SIM

| Analyte                | Result<br>mg/kg | Qualifier | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch                     |
|------------------------|-----------------|-----------|--------------|----------|-------------------------|---------------------------|
| Anthracene             | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Acenaphthene           | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Acenaphthylene         | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Benzo(a)anthracene     | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Benzo(a)pyrene         | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Benzo(b)fluoranthene   | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Benzo(g,h,i)perylene   | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Benzo(k)fluoranthene   | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Chrysene               | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Dibenz(a,h)anthracene  | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Fluoranthene           | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Fluorene               | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Indeno(1,2,3-cd)pyrene | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Naphthalene            | ND              |           | 0.0200       | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Phenanthrene           | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| Pyrene                 | ND              |           | 0.00600      | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| 1-Methylnaphthalene    | ND              |           | 0.0200       | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| 2-Methylnaphthalene    | ND              |           | 0.0200       | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| 2-Chloronaphthalene    | ND              |           | 0.0200       | 1        | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| (S) p-Terphenyl-d14    | 76.0            |           | 23.0-120     |          | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| (S) Nitrobenzene-d5    | 88.2            |           | 14.0-149     |          | 11/04/2018 22:11        | <a href="#">WG1191152</a> |
| (S) 2-Fluorobiphenyl   | 80.4            |           | 34.0-125     |          | 11/04/2018 22:11        | <a href="#">WG1191152</a> |





Method Blank (MB)

(MB) R3356436-1 11/02/18 15:50

| Analyte             | MB Result | MB Qualifier | MB MDL | MB RDL |
|---------------------|-----------|--------------|--------|--------|
| Chromium,Hexavalent | U         |              | 0.640  | 2.00   |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc

L1039204-04 Original Sample (OS) • Duplicate (DUP)

(OS) L1039204-04 11/02/18 15:52 • (DUP) R3356436-3 11/02/18 15:52

| Analyte             | Original Result | DUP Result | Dilution | DUP RPD | DUP Qualifier | DUP RPD Limits |
|---------------------|-----------------|------------|----------|---------|---------------|----------------|
| Chromium,Hexavalent | ND              | 0.000      | 1        | 0.000   |               | 20             |

L1039719-01 Original Sample (OS) • Duplicate (DUP)

(OS) L1039719-01 11/02/18 16:01 • (DUP) R3356436-8 11/02/18 16:02

| Analyte             | Original Result | DUP Result | Dilution | DUP RPD | DUP Qualifier | DUP RPD Limits |
|---------------------|-----------------|------------|----------|---------|---------------|----------------|
| Chromium,Hexavalent | U               | 0.000      | 1        | 0.000   |               | 20             |

Laboratory Control Sample (LCS)

(LCS) R3356436-2 11/02/18 15:51

| Analyte             | Spike Amount | LCS Result | LCS Rec. | Rec. Limits | LCS Qualifier |
|---------------------|--------------|------------|----------|-------------|---------------|
| Chromium,Hexavalent | 24.0         | 23.0       | 95.8     | 80.0-120    |               |

L1039278-04 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1039278-04 11/02/18 15:55 • (MS) R3356436-4 11/02/18 15:56 • (MSD) R3356436-5 11/02/18 15:56

| Analyte             | Spike Amount (dry) | Original Result (dry) | MS Result (dry) | MSD Result (dry) | MS Rec. | MSD Rec. | Dilution | Rec. Limits | MS Qualifier | MSD Qualifier | RPD  | RPD Limits |
|---------------------|--------------------|-----------------------|-----------------|------------------|---------|----------|----------|-------------|--------------|---------------|------|------------|
| Chromium,Hexavalent | 23.3               | U                     | 15.8            | 16.1             | 68.0    | 69.2     | 1        | 75.0-125    | <u>J6</u>    | <u>J6</u>     | 1.75 | 20         |



L1039278-04 Original Sample (OS) • Matrix Spike (MS)

(OS) L1039278-04 11/02/18 15:55 • (MS) R3356436-7 11/02/18 15:57

| Analyte             | Spike Amount (dry)<br>mg/kg | Original Result (dry)<br>mg/kg | MS Result (dry)<br>mg/kg | MS Rec.<br>% | Dilution | Rec. Limits<br>% | <u>MS Qualifier</u> |
|---------------------|-----------------------------|--------------------------------|--------------------------|--------------|----------|------------------|---------------------|
| Chromium,Hexavalent | 794                         | U                              | 781                      | 98.3         | 50       | 75.0-125         |                     |

- <sup>1</sup>Cp
- <sup>2</sup>Tc
- <sup>3</sup>Ss
- <sup>4</sup>Cn
- <sup>5</sup>Sr
- <sup>6</sup>Qc
- <sup>7</sup>Gl
- <sup>8</sup>Al
- <sup>9</sup>Sc



L1039670-05 Original Sample (OS) • Duplicate (DUP)

(OS) L1039670-05 11/01/18 17:54 • (DUP) R3356129-2 11/01/18 17:54

| Analyte | Original Result | DUP Result | Dilution | DUP RPD | DUP Qualifier | DUP RPD Limits |
|---------|-----------------|------------|----------|---------|---------------|----------------|
|         | su              | su         |          | %       |               | %              |
| pH      | 8.41            | 8.37       | 1        | 0.477   |               | 1              |

Sample Narrative:

OS: 8.41 at 20.7C  
DUP: 8.37 at 20.5C

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

L1039742-02 Original Sample (OS) • Duplicate (DUP)

(OS) L1039742-02 11/01/18 17:54 • (DUP) R3356129-3 11/01/18 17:54

| Analyte | Original Result | DUP Result | Dilution | DUP RPD | DUP Qualifier | DUP RPD Limits |
|---------|-----------------|------------|----------|---------|---------------|----------------|
|         | su              | su         |          | %       |               | %              |
| pH      | 6.05            | 6.02       | 1        | 0.497   |               | 1              |

Sample Narrative:

OS: 6.05 at 20.2C  
DUP: 6.02 at 19.9C

6 Qc

7 Gl

8 Al

9 Sc

Laboratory Control Sample (LCS)

(LCS) R3356129-1 11/01/18 17:54

| Analyte | Spike Amount | LCS Result | LCS Rec. | Rec. Limits | LCS Qualifier |
|---------|--------------|------------|----------|-------------|---------------|
|         | su           | su         | %        | %           |               |
| pH      | 10.0         | 9.95       | 99.5     | 99.0-101    |               |

Sample Narrative:

LCS: 9.95 at 18.7C



Method Blank (MB)

(MB) R3356621-1 11/03/18 15:00

| Analyte              | MB Result<br>umhos/cm | MB Qualifier | MB MDL<br>umhos/cm | MB RDL<br>umhos/cm |
|----------------------|-----------------------|--------------|--------------------|--------------------|
| Specific Conductance | U                     |              | 10.0               | 10.0               |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc

L1039204-06 Original Sample (OS) • Duplicate (DUP)

(OS) L1039204-06 11/03/18 15:00 • (DUP) R3356621-3 11/03/18 15:00

| Analyte              | Original Result<br>umhos/cm | DUP Result<br>umhos/cm | Dilution | DUP RPD<br>% | DUP Qualifier | DUP RPD<br>Limits |
|----------------------|-----------------------------|------------------------|----------|--------------|---------------|-------------------|
| Specific Conductance | 10300                       | 10400                  | 1        | 0.966        |               | 20                |

Laboratory Control Sample (LCS)

(LCS) R3356621-2 11/03/18 15:00

| Analyte              | Spike Amount<br>umhos/cm | LCS Result<br>umhos/cm | LCS Rec.<br>% | Rec. Limits<br>% | LCS Qualifier |
|----------------------|--------------------------|------------------------|---------------|------------------|---------------|
| Specific Conductance | 1160                     | 1170                   | 101           | 85.0-115         |               |



Method Blank (MB)

(MB) R3356158-1 11/01/18 18:55

| Analyte | MB Result | MB Qualifier | MB MDL  | MB RDL |
|---------|-----------|--------------|---------|--------|
| Mercury | U         |              | 0.00280 | 0.0200 |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3356158-2 11/01/18 18:58 • (LCSD) R3356158-3 11/01/18 19:00

| Analyte | Spike Amount | LCS Result | LCSD Result | LCS Rec. | LCSD Rec. | Rec. Limits | LCS Qualifier | LCSD Qualifier | RPD   | RPD Limits |
|---------|--------------|------------|-------------|----------|-----------|-------------|---------------|----------------|-------|------------|
| Mercury | 0.300        | 0.242      | 0.242       | 80.5     | 80.8      | 80.0-120    |               |                | 0.301 | 20         |

<sup>5</sup> Sr

<sup>6</sup> Qc

L1039759-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1039759-03 11/01/18 19:03 • (MS) R3356158-4 11/01/18 19:05 • (MSD) R3356158-5 11/01/18 19:08

| Analyte | Spike Amount | Original Result | MS Result | MSD Result | MS Rec. | MSD Rec. | Dilution | Rec. Limits | MS Qualifier | MSD Qualifier | RPD  | RPD Limits |
|---------|--------------|-----------------|-----------|------------|---------|----------|----------|-------------|--------------|---------------|------|------------|
| Mercury | 0.300        | 0.0236          | 0.216     | 0.229      | 64.2    | 68.3     | 1        | 75.0-125    | <u>J6</u>    | <u>J6</u>     | 5.46 | 20         |

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc



Method Blank (MB)

(MB) R3356764-1 11/04/18 14:40

| Analyte  | MB Result<br>mg/kg | MB Qualifier | MB MDL<br>mg/kg | MB RDL<br>mg/kg |
|----------|--------------------|--------------|-----------------|-----------------|
| Arsenic  | U                  |              | 0.460           | 2.00            |
| Barium   | U                  |              | 0.170           | 0.500           |
| Boron    | U                  |              | 1.26            | 10.0            |
| Cadmium  | U                  |              | 0.0700          | 0.500           |
| Chromium | U                  |              | 0.140           | 1.00            |
| Copper   | U                  |              | 0.530           | 2.00            |
| Lead     | U                  |              | 0.190           | 0.500           |
| Nickel   | U                  |              | 0.490           | 2.00            |
| Selenium | U                  |              | 0.620           | 2.00            |
| Silver   | U                  |              | 0.120           | 1.00            |
| Zinc     | U                  |              | 0.590           | 5.00            |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3356764-2 11/04/18 14:42 • (LCSD) R3356764-3 11/04/18 14:45

| Analyte  | Spike Amount<br>mg/kg | LCS Result<br>mg/kg | LCSD Result<br>mg/kg | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | LCS Qualifier | LCSD Qualifier | RPD<br>% | RPD Limits<br>% |
|----------|-----------------------|---------------------|----------------------|---------------|----------------|------------------|---------------|----------------|----------|-----------------|
| Arsenic  | 100                   | 101                 | 101                  | 101           | 101            | 80.0-120         |               |                | 0.601    | 20              |
| Barium   | 100                   | 103                 | 103                  | 103           | 103            | 80.0-120         |               |                | 0.0200   | 20              |
| Boron    | 100                   | 95.6                | 96.1                 | 95.6          | 96.1           | 80.0-120         |               |                | 0.527    | 20              |
| Cadmium  | 100                   | 102                 | 102                  | 102           | 102            | 80.0-120         |               |                | 0.00246  | 20              |
| Chromium | 100                   | 95.8                | 96.6                 | 95.8          | 96.6           | 80.0-120         |               |                | 0.821    | 20              |
| Copper   | 100                   | 97.1                | 97.8                 | 97.1          | 97.8           | 80.0-120         |               |                | 0.676    | 20              |
| Lead     | 100                   | 99.0                | 99.3                 | 99.0          | 99.3           | 80.0-120         |               |                | 0.316    | 20              |
| Nickel   | 100                   | 99.2                | 99.3                 | 99.2          | 99.3           | 80.0-120         |               |                | 0.131    | 20              |
| Selenium | 100                   | 101                 | 101                  | 101           | 101            | 80.0-120         |               |                | 0.127    | 20              |
| Silver   | 20.0                  | 17.8                | 18.0                 | 88.8          | 89.8           | 80.0-120         |               |                | 1.13     | 20              |
| Zinc     | 100                   | 99.0                | 99.0                 | 99.0          | 99.0           | 80.0-120         |               |                | 0.0345   | 20              |

L1039694-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1039694-01 11/04/18 14:47 • (MS) R3356764-6 11/04/18 14:55 • (MSD) R3356764-7 11/04/18 14:58

| Analyte | Spike Amount (dry)<br>mg/kg | Original Result (dry)<br>mg/kg | MS Result (dry)<br>mg/kg | MSD Result (dry)<br>mg/kg | MS Rec.<br>% | MSD Rec.<br>% | Dilution | Rec. Limits<br>% | MS Qualifier | MSD Qualifier | RPD<br>% | RPD Limits<br>% |
|---------|-----------------------------|--------------------------------|--------------------------|---------------------------|--------------|---------------|----------|------------------|--------------|---------------|----------|-----------------|
| Arsenic | 106                         | U                              | 98.9                     | 99.9                      | 93.0         | 93.9          | 1        | 75.0-125         |              |               | 1.02     | 20              |
| Barium  | 106                         | 98.8                           | 205                      | 216                       | 99.8         | 110           | 1        | 75.0-125         |              |               | 5.13     | 20              |
| Boron   | 106                         | U                              | 93.4                     | 93.9                      | 87.9         | 88.3          | 1        | 75.0-125         |              |               | 0.482    | 20              |
| Cadmium | 106                         | 0.120                          | 101                      | 102                       | 95.3         | 95.9          | 1        | 75.0-125         |              |               | 0.630    | 20              |



L1039694-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1039694-01 11/04/18 14:47 • (MS) R3356764-6 11/04/18 14:55 • (MSD) R3356764-7 11/04/18 14:58

| Analyte  | Spike Amount (dry)<br>mg/kg | Original Result (dry)<br>mg/kg | MS Result (dry)<br>mg/kg | MSD Result (dry)<br>mg/kg | MS Rec.<br>% | MSD Rec.<br>% | Dilution | Rec. Limits<br>% | MS Qualifier | MSD Qualifier | RPD<br>% | RPD Limits<br>% |
|----------|-----------------------------|--------------------------------|--------------------------|---------------------------|--------------|---------------|----------|------------------|--------------|---------------|----------|-----------------|
| Chromium | 106                         | 19.0                           | 111                      | 115                       | 86.1         | 89.9          | 1        | 75.0-125         |              |               | 3.57     | 20              |
| Copper   | 106                         | 8.00                           | 110                      | 111                       | 95.9         | 96.4          | 1        | 75.0-125         |              |               | 0.496    | 20              |
| Lead     | 106                         | 60.3                           | 164                      | 168                       | 97.5         | 101           | 1        | 75.0-125         |              |               | 2.42     | 20              |
| Nickel   | 106                         | 4.71                           | 110                      | 111                       | 98.7         | 100           | 1        | 75.0-125         |              |               | 1.34     | 20              |
| Selenium | 106                         | U                              | 98.3                     | 99.0                      | 92.5         | 93.2          | 1        | 75.0-125         |              |               | 0.743    | 20              |
| Silver   | 21.3                        | U                              | 18.1                     | 18.1                      | 85.2         | 85.0          | 1        | 75.0-125         |              |               | 0.319    | 20              |
| Zinc     | 106                         | 77.8                           | 171                      | 173                       | 87.7         | 89.7          | 1        | 75.0-125         |              |               | 1.25     | 20              |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc



Method Blank (MB)

(MB) R3356926-3 11/04/18 20:30

| Analyte                            | MB Result | MB Qualifier | MB MDL | MB RDL   |
|------------------------------------|-----------|--------------|--------|----------|
| TPH (GC/FID) Low Fraction          | 0.0463    | ↓            | 0.0217 | 0.100    |
| (S)<br>a,a,a-Trifluorotoluene(FID) | 98.2      |              |        | 77.0-120 |

1 Cp

2 Tc

3 Ss

4 Cn

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3356926-1 11/04/18 19:24 • (LCSD) R3356926-2 11/04/18 19:46

| Analyte                            | Spike Amount | LCS Result | LCSD Result | LCS Rec. | LCSD Rec. | Rec. Limits | LCS Qualifier | LCSD Qualifier | RPD  | RPD Limits |
|------------------------------------|--------------|------------|-------------|----------|-----------|-------------|---------------|----------------|------|------------|
| TPH (GC/FID) Low Fraction          | 5.50         | 5.75       | 5.58        | 104      | 101       | 72.0-127    |               |                | 2.91 | 20         |
| (S)<br>a,a,a-Trifluorotoluene(FID) |              |            |             | 102      | 102       | 77.0-120    |               |                |      |            |

5 Sr

6 Qc

7 Gl

L1039465-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1039465-01 11/05/18 05:03 • (MS) R3356926-4 11/05/18 05:25 • (MSD) R3356926-5 11/05/18 05:47

| Analyte                            | Spike Amount | Original Result | MS Result | MSD Result | MS Rec. | MSD Rec. | Dilution | Rec. Limits | MS Qualifier | MSD Qualifier | RPD  | RPD Limits |
|------------------------------------|--------------|-----------------|-----------|------------|---------|----------|----------|-------------|--------------|---------------|------|------------|
| TPH (GC/FID) Low Fraction          | 5.50         | ND              | 50.1      | 52.4       | 34.3    | 35.9     | 26       | 10.0-151    |              |               | 4.55 | 28         |
| (S)<br>a,a,a-Trifluorotoluene(FID) |              |                 |           |            | 94.4    | 94.6     |          | 77.0-120    |              |               |      |            |

8 Al

9 Sc

Sample Narrative:

OS: Only received a MEOH vial.





Method Blank (MB)

(MB) R3357277-3 11/06/18 01:02

| Analyte                                       | MB Result<br>mg/kg | MB Qualifier | MB MDL<br>mg/kg | MB RDL<br>mg/kg |
|-----------------------------------------------|--------------------|--------------|-----------------|-----------------|
| TPH (GC/FID) Low Fraction                     | U                  |              | 0.0217          | 0.100           |
| <sup>(S)</sup><br>a,a,a-Trifluorotoluene(FID) | 100                |              |                 | 77.0-120        |

1 Cp

2 Tc

3 Ss

4 Cn

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3357277-1 11/05/18 23:49 • (LCSD) R3357277-2 11/06/18 00:13

| Analyte                                       | Spike Amount<br>mg/kg | LCS Result<br>mg/kg | LCSD Result<br>mg/kg | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | LCS Qualifier | LCSD Qualifier | RPD<br>% | RPD Limits<br>% |
|-----------------------------------------------|-----------------------|---------------------|----------------------|---------------|----------------|------------------|---------------|----------------|----------|-----------------|
| TPH (GC/FID) Low Fraction                     | 5.50                  | 5.81                | 5.74                 | 106           | 104            | 72.0-127         |               |                | 1.20     | 20              |
| <sup>(S)</sup><br>a,a,a-Trifluorotoluene(FID) |                       |                     |                      | 105           | 105            | 77.0-120         |               |                |          |                 |

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



Method Blank (MB)

(MB) R3357449-2 11/01/18 20:46

| Analyte                     | MB Result<br>mg/kg | MB Qualifier | MB MDL<br>mg/kg | MB RDL<br>mg/kg |
|-----------------------------|--------------------|--------------|-----------------|-----------------|
| Acetone                     | U                  |              | 0.0137          | 0.0250          |
| Acrylonitrile               | U                  |              | 0.00190         | 0.0125          |
| Benzene                     | 0.000967           | U            | 0.000400        | 0.00100         |
| Bromobenzene                | U                  |              | 0.00105         | 0.0125          |
| Bromodichloromethane        | U                  |              | 0.000788        | 0.00250         |
| Bromoform                   | U                  |              | 0.00598         | 0.0250          |
| Bromomethane                | U                  |              | 0.00370         | 0.0125          |
| n-Butylbenzene              | U                  |              | 0.00384         | 0.0125          |
| sec-Butylbenzene            | U                  |              | 0.00253         | 0.0125          |
| tert-Butylbenzene           | U                  |              | 0.00155         | 0.00500         |
| Carbon tetrachloride        | U                  |              | 0.00108         | 0.00500         |
| Chlorobenzene               | U                  |              | 0.000573        | 0.00250         |
| Chlorodibromomethane        | U                  |              | 0.000450        | 0.00250         |
| Chloroethane                | U                  |              | 0.00108         | 0.00500         |
| Chloroform                  | U                  |              | 0.000415        | 0.00250         |
| Chloromethane               | U                  |              | 0.00139         | 0.0125          |
| 2-Chlorotoluene             | U                  |              | 0.000920        | 0.00250         |
| 4-Chlorotoluene             | U                  |              | 0.00113         | 0.00500         |
| 1,2-Dibromo-3-Chloropropane | U                  |              | 0.00510         | 0.0250          |
| 1,2-Dibromoethane           | U                  |              | 0.000525        | 0.00250         |
| Dibromomethane              | U                  |              | 0.00100         | 0.00500         |
| 1,2-Dichlorobenzene         | U                  |              | 0.00145         | 0.00500         |
| 1,3-Dichlorobenzene         | U                  |              | 0.00170         | 0.00500         |
| 1,4-Dichlorobenzene         | U                  |              | 0.00197         | 0.00500         |
| Dichlorodifluoromethane     | U                  |              | 0.000818        | 0.00250         |
| 1,1-Dichloroethane          | U                  |              | 0.000575        | 0.00250         |
| 1,2-Dichloroethane          | U                  |              | 0.000475        | 0.00250         |
| 1,1-Dichloroethene          | U                  |              | 0.000500        | 0.00250         |
| cis-1,2-Dichloroethene      | U                  |              | 0.000690        | 0.00250         |
| trans-1,2-Dichloroethene    | U                  |              | 0.00143         | 0.00500         |
| 1,2-Dichloropropane         | U                  |              | 0.00127         | 0.00500         |
| 1,1-Dichloropropene         | U                  |              | 0.000700        | 0.00250         |
| 1,3-Dichloropropane         | U                  |              | 0.00175         | 0.00500         |
| cis-1,3-Dichloropropene     | U                  |              | 0.000678        | 0.00250         |
| trans-1,3-Dichloropropene   | U                  |              | 0.00153         | 0.00500         |
| 2,2-Dichloropropane         | U                  |              | 0.000793        | 0.00250         |
| Di-isopropyl ether          | U                  |              | 0.000350        | 0.00100         |
| Ethylbenzene                | U                  |              | 0.000530        | 0.00250         |
| Hexachloro-1,3-butadiene    | U                  |              | 0.0127          | 0.0250          |
| Isopropylbenzene            | U                  |              | 0.000863        | 0.00250         |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc



Method Blank (MB)

(MB) R3357449-2 11/01/18 20:46

| Analyte                        | MB Result<br>mg/kg | MB Qualifier | MB MDL<br>mg/kg | MB RDL<br>mg/kg |
|--------------------------------|--------------------|--------------|-----------------|-----------------|
| p-Isopropyltoluene             | U                  |              | 0.00233         | 0.00500         |
| 2-Butanone (MEK)               | U                  |              | 0.0125          | 0.0250          |
| Methylene Chloride             | U                  |              | 0.00664         | 0.0250          |
| 4-Methyl-2-pentanone (MIBK)    | U                  |              | 0.0100          | 0.0250          |
| Methyl tert-butyl ether        | U                  |              | 0.000295        | 0.00100         |
| Naphthalene                    | U                  |              | 0.00312         | 0.0125          |
| n-Propylbenzene                | U                  |              | 0.00118         | 0.00500         |
| Styrene                        | U                  |              | 0.00273         | 0.0125          |
| 1,1,1,2-Tetrachloroethane      | U                  |              | 0.000500        | 0.00250         |
| 1,1,2,2-Tetrachloroethane      | U                  |              | 0.000390        | 0.00250         |
| Tetrachloroethene              | U                  |              | 0.000700        | 0.00250         |
| Toluene                        | U                  |              | 0.00125         | 0.00500         |
| 1,1,2-Trichlorotrifluoroethane | U                  |              | 0.000675        | 0.00250         |
| 1,2,3-Trichlorobenzene         | U                  |              | 0.000625        | 0.00250         |
| 1,2,4-Trichlorobenzene         | U                  |              | 0.00482         | 0.0125          |
| 1,1,1-Trichloroethane          | U                  |              | 0.000275        | 0.00250         |
| 1,1,2-Trichloroethane          | U                  |              | 0.000883        | 0.00250         |
| Trichloroethene                | U                  |              | 0.000400        | 0.00100         |
| Trichlorofluoromethane         | U                  |              | 0.000500        | 0.00250         |
| 1,2,3-Trichloropropane         | U                  |              | 0.00510         | 0.0125          |
| 1,2,3-Trimethylbenzene         | U                  |              | 0.00115         | 0.00500         |
| 1,2,4-Trimethylbenzene         | U                  |              | 0.00116         | 0.00500         |
| 1,3,5-Trimethylbenzene         | U                  |              | 0.00108         | 0.00500         |
| Vinyl chloride                 | U                  |              | 0.000683        | 0.00250         |
| Xylenes, Total                 | U                  |              | 0.00478         | 0.00650         |
| (S) Toluene-d8                 | 125                |              |                 | 75.0-131        |
| (S) Dibromofluoromethane       | 98.2               |              |                 | 65.0-129        |
| (S) 4-Bromofluorobenzene       | 84.2               |              |                 | 67.0-138        |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc

Laboratory Control Sample (LCS)

(LCS) R3357449-1 11/01/18 19:36

| Analyte              | Spike Amount<br>mg/kg | LCS Result<br>mg/kg | LCS Rec.<br>% | Rec. Limits<br>% | LCS Qualifier |
|----------------------|-----------------------|---------------------|---------------|------------------|---------------|
| Acetone              | 0.625                 | 0.692               | 111           | 10.0-160         |               |
| Acrylonitrile        | 0.625                 | 0.677               | 108           | 45.0-153         |               |
| Benzene              | 0.125                 | 0.126               | 101           | 70.0-123         |               |
| Bromobenzene         | 0.125                 | 0.115               | 92.3          | 73.0-121         |               |
| Bromodichloromethane | 0.125                 | 0.124               | 99.5          | 73.0-121         |               |



Laboratory Control Sample (LCS)

(LCS) R3357449-1 11/01/18 19:36

| Analyte                     | Spike Amount<br>mg/kg | LCS Result<br>mg/kg | LCS Rec.<br>% | Rec. Limits<br>% | <u>LCS Qualifier</u> |
|-----------------------------|-----------------------|---------------------|---------------|------------------|----------------------|
| Bromoform                   | 0.125                 | 0.0839              | 67.1          | 64.0-132         |                      |
| Bromomethane                | 0.125                 | 0.135               | 108           | 56.0-147         |                      |
| n-Butylbenzene              | 0.125                 | 0.129               | 103           | 68.0-135         |                      |
| sec-Butylbenzene            | 0.125                 | 0.128               | 102           | 74.0-130         |                      |
| tert-Butylbenzene           | 0.125                 | 0.121               | 97.1          | 75.0-127         |                      |
| Carbon tetrachloride        | 0.125                 | 0.121               | 97.2          | 66.0-128         |                      |
| Chlorobenzene               | 0.125                 | 0.125               | 99.7          | 76.0-128         |                      |
| Chlorodibromomethane        | 0.125                 | 0.107               | 86.0          | 74.0-127         |                      |
| Chloroethane                | 0.125                 | 0.131               | 105           | 61.0-134         |                      |
| Chloroform                  | 0.125                 | 0.138               | 111           | 72.0-123         |                      |
| Chloromethane               | 0.125                 | 0.133               | 106           | 51.0-138         |                      |
| 2-Chlorotoluene             | 0.125                 | 0.116               | 92.6          | 75.0-124         |                      |
| 4-Chlorotoluene             | 0.125                 | 0.124               | 99.0          | 75.0-124         |                      |
| 1,2-Dibromo-3-Chloropropane | 0.125                 | 0.112               | 90.0          | 59.0-130         |                      |
| 1,2-Dibromoethane           | 0.125                 | 0.126               | 101           | 74.0-128         |                      |
| Dibromomethane              | 0.125                 | 0.126               | 101           | 75.0-122         |                      |
| 1,2-Dichlorobenzene         | 0.125                 | 0.138               | 110           | 76.0-124         |                      |
| 1,3-Dichlorobenzene         | 0.125                 | 0.124               | 99.6          | 76.0-125         |                      |
| 1,4-Dichlorobenzene         | 0.125                 | 0.117               | 93.9          | 77.0-121         |                      |
| Dichlorodifluoromethane     | 0.125                 | 0.149               | 120           | 43.0-156         |                      |
| 1,1-Dichloroethane          | 0.125                 | 0.145               | 116           | 70.0-127         |                      |
| 1,2-Dichloroethane          | 0.125                 | 0.118               | 94.5          | 65.0-131         |                      |
| 1,1-Dichloroethene          | 0.125                 | 0.136               | 108           | 65.0-131         |                      |
| cis-1,2-Dichloroethene      | 0.125                 | 0.145               | 116           | 73.0-125         |                      |
| trans-1,2-Dichloroethene    | 0.125                 | 0.133               | 106           | 71.0-125         |                      |
| 1,2-Dichloropropane         | 0.125                 | 0.130               | 104           | 74.0-125         |                      |
| 1,1-Dichloropropene         | 0.125                 | 0.130               | 104           | 73.0-125         |                      |
| 1,3-Dichloropropane         | 0.125                 | 0.131               | 105           | 80.0-125         |                      |
| cis-1,3-Dichloropropene     | 0.125                 | 0.117               | 94.0          | 76.0-127         |                      |
| trans-1,3-Dichloropropene   | 0.125                 | 0.107               | 86.0          | 73.0-127         |                      |
| 2,2-Dichloropropane         | 0.125                 | 0.176               | 140           | 59.0-135         | J4                   |
| Di-isopropyl ether          | 0.125                 | 0.141               | 113           | 60.0-136         |                      |
| Ethylbenzene                | 0.125                 | 0.147               | 118           | 74.0-126         |                      |
| Hexachloro-1,3-butadiene    | 0.125                 | 0.133               | 107           | 57.0-150         |                      |
| Isopropylbenzene            | 0.125                 | 0.119               | 94.9          | 72.0-127         |                      |
| p-Isopropyltoluene          | 0.125                 | 0.128               | 103           | 72.0-133         |                      |
| 2-Butanone (MEK)            | 0.625                 | 0.601               | 96.1          | 30.0-160         |                      |
| Methylene Chloride          | 0.125                 | 0.135               | 108           | 68.0-123         |                      |
| 4-Methyl-2-pentanone (MIBK) | 0.625                 | 0.665               | 106           | 56.0-143         |                      |
| Methyl tert-butyl ether     | 0.125                 | 0.166               | 133           | 66.0-132         | J4                   |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc



Laboratory Control Sample (LCS)

(LCS) R3357449-1 11/01/18 19:36

| Analyte                         | Spike Amount<br>mg/kg | LCS Result<br>mg/kg | LCS Rec.<br>% | Rec. Limits<br>% | <u>LCS Qualifier</u> |
|---------------------------------|-----------------------|---------------------|---------------|------------------|----------------------|
| Naphthalene                     | 0.125                 | 0.155               | 124           | 59.0-130         |                      |
| n-Propylbenzene                 | 0.125                 | 0.120               | 96.3          | 74.0-126         |                      |
| Styrene                         | 0.125                 | 0.0991              | 79.3          | 72.0-127         |                      |
| 1,1,1,2-Tetrachloroethane       | 0.125                 | 0.137               | 110           | 74.0-129         |                      |
| 1,1,2,2-Tetrachloroethane       | 0.125                 | 0.119               | 94.9          | 68.0-128         |                      |
| Tetrachloroethene               | 0.125                 | 0.144               | 115           | 70.0-136         |                      |
| Toluene                         | 0.125                 | 0.141               | 113           | 75.0-121         |                      |
| 1,1,2-Trichlorotrifluoroethane  | 0.125                 | 0.153               | 122           | 61.0-139         |                      |
| 1,2,3-Trichlorobenzene          | 0.125                 | 0.171               | 136           | 59.0-139         |                      |
| 1,2,4-Trichlorobenzene          | 0.125                 | 0.142               | 113           | 62.0-137         |                      |
| 1,1,1-Trichloroethane           | 0.125                 | 0.126               | 101           | 69.0-126         |                      |
| 1,1,2-Trichloroethane           | 0.125                 | 0.134               | 107           | 78.0-123         |                      |
| Trichloroethene                 | 0.125                 | 0.106               | 84.9          | 76.0-126         |                      |
| Trichlorofluoromethane          | 0.125                 | 0.153               | 123           | 61.0-142         |                      |
| 1,2,3-Trichloropropane          | 0.125                 | 0.122               | 97.6          | 67.0-129         |                      |
| 1,2,3-Trimethylbenzene          | 0.125                 | 0.130               | 104           | 74.0-124         |                      |
| 1,2,4-Trimethylbenzene          | 0.125                 | 0.123               | 98.2          | 70.0-126         |                      |
| 1,3,5-Trimethylbenzene          | 0.125                 | 0.121               | 96.7          | 73.0-127         |                      |
| Vinyl chloride                  | 0.125                 | 0.132               | 106           | 63.0-134         |                      |
| Xylenes, Total                  | 0.375                 | 0.407               | 109           | 72.0-127         |                      |
| <i>(S) Toluene-d8</i>           |                       |                     | 113           | 75.0-131         |                      |
| <i>(S) Dibromofluoromethane</i> |                       |                     | 110           | 65.0-129         |                      |
| <i>(S) 4-Bromofluorobenzene</i> |                       |                     | 87.3          | 67.0-138         |                      |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc

L1039759-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1039759-03 11/02/18 02:00 • (MS) R3357449-3 11/01/18 21:06 • (MSD) R3357449-4 11/01/18 21:36

| Analyte              | Spike Amount<br>(dry)<br>mg/kg | Original Result<br>(dry)<br>mg/kg | MS Result (dry)<br>mg/kg | MSD Result<br>(dry)<br>mg/kg | MS Rec.<br>% | MSD Rec.<br>% | Dilution | Rec. Limits<br>% | <u>MS Qualifier</u> | <u>MSD Qualifier</u> | RPD<br>% | RPD Limits<br>% |
|----------------------|--------------------------------|-----------------------------------|--------------------------|------------------------------|--------------|---------------|----------|------------------|---------------------|----------------------|----------|-----------------|
| Acetone              | 0.751                          | U                                 | 0.653                    | 0.686                        | 86.9         | 91.4          | 1        | 10.0-160         |                     |                      | 5.02     | 40              |
| Acrylonitrile        | 0.751                          | U                                 | 0.589                    | 0.602                        | 78.5         | 80.2          | 1        | 10.0-160         |                     |                      | 2.18     | 40              |
| Benzene              | 0.150                          | 0.000667                          | 0.120                    | 0.124                        | 79.2         | 82.1          | 1        | 10.0-149         |                     |                      | 3.58     | 37              |
| Bromobenzene         | 0.150                          | U                                 | 0.101                    | 0.111                        | 67.6         | 74.2          | 1        | 10.0-156         |                     |                      | 9.32     | 38              |
| Bromodichloromethane | 0.150                          | U                                 | 0.115                    | 0.119                        | 76.6         | 79.3          | 1        | 10.0-143         |                     |                      | 3.53     | 37              |
| Bromoform            | 0.150                          | U                                 | 0.0699                   | 0.0749                       | 46.5         | 49.9          | 1        | 10.0-146         |                     |                      | 6.96     | 36              |
| Bromomethane         | 0.150                          | U                                 | 0.0752                   | 0.0843                       | 50.1         | 56.1          | 1        | 10.0-149         |                     |                      | 11.4     | 38              |
| n-Butylbenzene       | 0.150                          | U                                 | 0.122                    | 0.134                        | 81.5         | 89.5          | 1        | 10.0-160         |                     |                      | 9.38     | 40              |
| sec-Butylbenzene     | 0.150                          | U                                 | 0.123                    | 0.130                        | 81.8         | 86.5          | 1        | 10.0-159         |                     |                      | 5.63     | 39              |
| tert-Butylbenzene    | 0.150                          | U                                 | 0.111                    | 0.119                        | 74.1         | 79.4          | 1        | 10.0-156         |                     |                      | 7.00     | 39              |



L1039759-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1039759-03 11/02/18 02:00 • (MS) R3357449-3 11/01/18 21:06 • (MSD) R3357449-4 11/01/18 21:36

| Analyte                     | Spike Amount (dry)<br>mg/kg | Original Result (dry)<br>mg/kg | MS Result (dry)<br>mg/kg | MSD Result (dry)<br>mg/kg | MS Rec.<br>% | MSD Rec.<br>% | Dilution | Rec. Limits<br>% | MS Qualifier | MSD Qualifier | RPD<br>% | RPD Limits<br>% |
|-----------------------------|-----------------------------|--------------------------------|--------------------------|---------------------------|--------------|---------------|----------|------------------|--------------|---------------|----------|-----------------|
| Carbon tetrachloride        | 0.150                       | U                              | 0.108                    | 0.109                     | 71.6         | 72.8          | 1        | 10.0-145         |              |               | 1.61     | 37              |
| Chlorobenzene               | 0.150                       | U                              | 0.117                    | 0.127                     | 77.8         | 84.3          | 1        | 10.0-152         |              |               | 8.01     | 39              |
| Chlorodibromomethane        | 0.150                       | U                              | 0.0916                   | 0.100                     | 61.0         | 66.9          | 1        | 10.0-146         |              |               | 9.24     | 37              |
| Chloroethane                | 0.150                       | U                              | 0.0869                   | 0.0874                    | 57.8         | 58.2          | 1        | 10.0-146         |              |               | 0.618    | 40              |
| Chloroform                  | 0.150                       | U                              | 0.124                    | 0.128                     | 82.3         | 85.2          | 1        | 10.0-146         |              |               | 3.47     | 37              |
| Chloromethane               | 0.150                       | U                              | 0.120                    | 0.111                     | 79.8         | 73.6          | 1        | 10.0-159         |              |               | 8.08     | 37              |
| 2-Chlorotoluene             | 0.150                       | U                              | 0.104                    | 0.110                     | 69.5         | 73.5          | 1        | 10.0-159         |              |               | 5.59     | 38              |
| 4-Chlorotoluene             | 0.150                       | U                              | 0.109                    | 0.117                     | 72.3         | 78.2          | 1        | 10.0-155         |              |               | 7.89     | 39              |
| 1,2-Dibromo-3-Chloropropane | 0.150                       | U                              | 0.0852                   | 0.0899                    | 56.7         | 59.8          | 1        | 10.0-151         |              |               | 5.40     | 39              |
| 1,2-Dibromoethane           | 0.150                       | U                              | 0.114                    | 0.120                     | 75.6         | 79.9          | 1        | 10.0-148         |              |               | 5.46     | 34              |
| Dibromomethane              | 0.150                       | U                              | 0.115                    | 0.119                     | 76.5         | 79.1          | 1        | 10.0-147         |              |               | 3.40     | 35              |
| 1,2-Dichlorobenzene         | 0.150                       | U                              | 0.124                    | 0.132                     | 82.3         | 88.1          | 1        | 10.0-155         |              |               | 6.73     | 37              |
| 1,3-Dichlorobenzene         | 0.150                       | U                              | 0.117                    | 0.125                     | 77.6         | 83.3          | 1        | 10.0-153         |              |               | 7.12     | 38              |
| 1,4-Dichlorobenzene         | 0.150                       | U                              | 0.110                    | 0.120                     | 73.4         | 80.2          | 1        | 10.0-151         |              |               | 8.91     | 38              |
| Dichlorodifluoromethane     | 0.150                       | U                              | 0.181                    | 0.155                     | 120          | 103           | 1        | 10.0-160         |              |               | 15.7     | 35              |
| 1,1-Dichloroethane          | 0.150                       | U                              | 0.126                    | 0.133                     | 83.9         | 88.7          | 1        | 10.0-147         |              |               | 5.54     | 37              |
| 1,2-Dichloroethane          | 0.150                       | U                              | 0.109                    | 0.113                     | 72.8         | 75.3          | 1        | 10.0-148         |              |               | 3.39     | 35              |
| 1,1-Dichloroethene          | 0.150                       | U                              | 0.116                    | 0.122                     | 77.5         | 81.5          | 1        | 10.0-155         |              |               | 4.96     | 37              |
| cis-1,2-Dichloroethene      | 0.150                       | U                              | 0.127                    | 0.131                     | 84.9         | 87.5          | 1        | 10.0-149         |              |               | 3.01     | 37              |
| trans-1,2-Dichloroethene    | 0.150                       | U                              | 0.116                    | 0.122                     | 77.1         | 81.3          | 1        | 10.0-150         |              |               | 5.37     | 37              |
| 1,2-Dichloropropane         | 0.150                       | U                              | 0.130                    | 0.125                     | 86.6         | 82.9          | 1        | 10.0-148         |              |               | 4.37     | 37              |
| 1,1-Dichloropropene         | 0.150                       | U                              | 0.121                    | 0.118                     | 80.5         | 78.5          | 1        | 10.0-153         |              |               | 2.41     | 35              |
| 1,3-Dichloropropane         | 0.150                       | U                              | 0.122                    | 0.133                     | 81.4         | 88.3          | 1        | 10.0-154         |              |               | 8.20     | 35              |
| cis-1,3-Dichloropropene     | 0.150                       | U                              | 0.105                    | 0.118                     | 70.2         | 78.5          | 1        | 10.0-151         |              |               | 11.2     | 37              |
| trans-1,3-Dichloropropene   | 0.150                       | U                              | 0.0991                   | 0.105                     | 66.0         | 70.1          | 1        | 10.0-148         |              |               | 5.99     | 37              |
| 2,2-Dichloropropane         | 0.150                       | U                              | 0.145                    | 0.160                     | 96.7         | 106           | 1        | 10.0-138         |              |               | 9.47     | 36              |
| Di-isopropyl ether          | 0.150                       | U                              | 0.122                    | 0.133                     | 81.4         | 88.8          | 1        | 10.0-147         |              |               | 8.72     | 36              |
| Ethylbenzene                | 0.150                       | U                              | 0.136                    | 0.140                     | 90.5         | 93.3          | 1        | 10.0-160         |              |               | 2.98     | 38              |
| Hexachloro-1,3-butadiene    | 0.150                       | U                              | 0.126                    | 0.145                     | 83.6         | 96.8          | 1        | 10.0-160         |              |               | 14.6     | 40              |
| Isopropylbenzene            | 0.150                       | U                              | 0.109                    | 0.117                     | 72.8         | 77.6          | 1        | 10.0-155         |              |               | 6.39     | 38              |
| p-Isopropyltoluene          | 0.150                       | U                              | 0.122                    | 0.131                     | 81.4         | 87.2          | 1        | 10.0-160         |              |               | 6.85     | 40              |
| 2-Butanone (MEK)            | 0.751                       | U                              | 0.605                    | 0.638                     | 80.5         | 84.9          | 1        | 10.0-160         |              |               | 5.33     | 40              |
| Methylene Chloride          | 0.150                       | U                              | 0.121                    | 0.127                     | 80.5         | 84.4          | 1        | 10.0-141         |              |               | 4.73     | 37              |
| 4-Methyl-2-pentanone (MIBK) | 0.751                       | U                              | 0.544                    | 0.550                     | 72.5         | 73.2          | 1        | 10.0-160         |              |               | 1.02     | 35              |
| Methyl tert-butyl ether     | 0.150                       | U                              | 0.144                    | 0.148                     | 95.8         | 98.2          | 1        | 11.0-147         |              |               | 2.53     | 35              |
| Naphthalene                 | 0.150                       | U                              | 0.133                    | 0.137                     | 88.4         | 91.4          | 1        | 10.0-160         |              |               | 3.33     | 36              |
| n-Propylbenzene             | 0.150                       | U                              | 0.109                    | 0.118                     | 72.7         | 78.4          | 1        | 10.0-158         |              |               | 7.62     | 38              |
| Styrene                     | 0.150                       | U                              | 0.0919                   | 0.103                     | 61.2         | 68.8          | 1        | 10.0-160         |              |               | 11.7     | 40              |
| 1,1,1,2-Tetrachloroethane   | 0.150                       | U                              | 0.120                    | 0.123                     | 79.7         | 81.9          | 1        | 10.0-149         |              |               | 2.67     | 39              |

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



L1039759-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1039759-03 11/02/18 02:00 • (MS) R3357449-3 11/01/18 21:06 • (MSD) R3357449-4 11/01/18 21:36

| Analyte                        | Spike Amount (dry)<br>mg/kg | Original Result (dry)<br>mg/kg | MS Result (dry)<br>mg/kg | MSD Result (dry)<br>mg/kg | MS Rec.<br>% | MSD Rec.<br>% | Dilution | Rec. Limits<br>% | MS Qualifier | MSD Qualifier | RPD<br>% | RPD Limits<br>% |
|--------------------------------|-----------------------------|--------------------------------|--------------------------|---------------------------|--------------|---------------|----------|------------------|--------------|---------------|----------|-----------------|
| 1,1,2,2-Tetrachloroethane      | 0.150                       | U                              | 0.101                    | 0.105                     | 67.5         | 70.0          | 1        | 10.0-160         |              |               | 3.62     | 35              |
| Tetrachloroethene              | 0.150                       | U                              | 0.142                    | 0.153                     | 94.8         | 102           | 1        | 10.0-156         |              |               | 7.34     | 39              |
| Toluene                        | 0.150                       | U                              | 0.138                    | 0.144                     | 91.8         | 95.7          | 1        | 10.0-156         |              |               | 4.18     | 38              |
| 1,1,2-Trichlorotrifluoroethane | 0.150                       | U                              | 0.126                    | 0.129                     | 83.9         | 86.1          | 1        | 10.0-160         |              |               | 2.59     | 36              |
| 1,2,3-Trichlorobenzene         | 0.150                       | U                              | 0.154                    | 0.171                     | 103          | 114           | 1        | 10.0-160         |              |               | 10.5     | 40              |
| 1,2,4-Trichlorobenzene         | 0.150                       | U                              | 0.129                    | 0.141                     | 85.9         | 94.1          | 1        | 10.0-160         |              |               | 9.18     | 40              |
| 1,1,1-Trichloroethane          | 0.150                       | U                              | 0.109                    | 0.113                     | 72.3         | 75.2          | 1        | 10.0-144         |              |               | 3.98     | 35              |
| 1,1,2-Trichloroethane          | 0.150                       | U                              | 0.121                    | 0.136                     | 80.5         | 90.3          | 1        | 10.0-160         |              |               | 11.5     | 35              |
| Trichloroethene                | 0.150                       | U                              | 0.0983                   | 0.0995                    | 65.5         | 66.2          | 1        | 10.0-156         |              |               | 1.17     | 38              |
| Trichlorofluoromethane         | 0.150                       | U                              | 0.132                    | 0.138                     | 88.0         | 91.8          | 1        | 10.0-160         |              |               | 4.24     | 40              |
| 1,2,3-Trichloropropane         | 0.150                       | U                              | 0.104                    | 0.103                     | 68.9         | 68.3          | 1        | 10.0-156         |              |               | 0.974    | 35              |
| 1,2,3-Trimethylbenzene         | 0.150                       | U                              | 0.109                    | 0.117                     | 72.9         | 77.6          | 1        | 10.0-160         |              |               | 6.32     | 36              |
| 1,2,4-Trimethylbenzene         | 0.150                       | U                              | 0.109                    | 0.119                     | 72.6         | 78.9          | 1        | 10.0-160         |              |               | 8.33     | 36              |
| 1,3,5-Trimethylbenzene         | 0.150                       | U                              | 0.108                    | 0.117                     | 71.9         | 77.7          | 1        | 10.0-160         |              |               | 7.74     | 38              |
| Vinyl chloride                 | 0.150                       | U                              | 0.0604                   | 0.0611                    | 40.2         | 40.7          | 1        | 10.0-160         |              |               | 1.05     | 37              |
| Xylenes, Total                 | 0.451                       | U                              | 0.374                    | 0.385                     | 82.9         | 85.3          | 1        | 10.0-160         |              |               | 2.85     | 38              |
| (S) Toluene-d8                 |                             |                                |                          |                           | 112          | 119           |          | 75.0-131         |              |               |          |                 |
| (S) Dibromofluoromethane       |                             |                                |                          |                           | 94.8         | 105           |          | 65.0-129         |              |               |          |                 |
| (S) 4-Bromofluorobenzene       |                             |                                |                          |                           | 83.0         | 88.9          |          | 67.0-138         |              |               |          |                 |

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



Method Blank (MB)

(MB) R3357599-2 11/06/18 23:24

| Analyte                  | MB Result<br>mg/kg | MB Qualifier | MB MDL<br>mg/kg | MB RDL<br>mg/kg |
|--------------------------|--------------------|--------------|-----------------|-----------------|
| Benzene                  | U                  |              | 0.000400        | 0.00100         |
| (S) Toluene-d8           | 98.3               |              |                 | 75.0-131        |
| (S) Dibromofluoromethane | 122                |              |                 | 65.0-129        |
| (S) 4-Bromofluorobenzene | 101                |              |                 | 67.0-138        |

Laboratory Control Sample (LCS)

(LCS) R3357599-1 11/06/18 21:59

| Analyte                  | Spike Amount<br>mg/kg | LCS Result<br>mg/kg | LCS Rec.<br>% | Rec. Limits<br>% | LCS Qualifier |
|--------------------------|-----------------------|---------------------|---------------|------------------|---------------|
| Benzene                  | 0.125                 | 0.132               | 106           | 70.0-123         |               |
| (S) Toluene-d8           |                       |                     | 97.8          | 75.0-131         |               |
| (S) Dibromofluoromethane |                       |                     | 121           | 65.0-129         |               |
| (S) 4-Bromofluorobenzene |                       |                     | 101           | 67.0-138         |               |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc





Method Blank (MB)

(MB) R3357494-1 11/06/18 15:09

| Analyte                | MB Result<br>mg/kg | MB Qualifier | MB MDL<br>mg/kg | MB RDL<br>mg/kg |
|------------------------|--------------------|--------------|-----------------|-----------------|
| C10-C28 Diesel Range   | U                  |              | 1.61            | 4.00            |
| C28-C40 Oil Range      | U                  |              | 0.274           | 4.00            |
| <i>(S) o-Terphenyl</i> | 83.2               |              |                 | 18.0-148        |

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3357494-2 11/06/18 15:22 • (LCSD) R3357494-3 11/06/18 15:38

| Analyte                | Spike Amount<br>mg/kg | LCS Result<br>mg/kg | LCSD Result<br>mg/kg | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | LCS Qualifier | LCSD Qualifier | RPD<br>% | RPD Limits<br>% |
|------------------------|-----------------------|---------------------|----------------------|---------------|----------------|------------------|---------------|----------------|----------|-----------------|
| C10-C28 Diesel Range   | 50.0                  | 32.1                | 34.6                 | 64.2          | 69.2           | 50.0-150         |               |                | 7.50     | 20              |
| <i>(S) o-Terphenyl</i> |                       |                     |                      | 80.6          | 91.0           | 18.0-148         |               |                |          |                 |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc



Method Blank (MB)

(MB) R3356952-3 11/04/18 19:43

| Analyte                | MB Result<br>mg/kg | MB Qualifier | MB MDL<br>mg/kg | MB RDL<br>mg/kg |
|------------------------|--------------------|--------------|-----------------|-----------------|
| Anthracene             | U                  |              | 0.00600         | 0.00600         |
| Acenaphthene           | U                  |              | 0.00600         | 0.00600         |
| Acenaphthylene         | U                  |              | 0.00600         | 0.00600         |
| Benzo(a)anthracene     | U                  |              | 0.00600         | 0.00600         |
| Benzo(a)pyrene         | U                  |              | 0.00600         | 0.00600         |
| Benzo(b)fluoranthene   | U                  |              | 0.00600         | 0.00600         |
| Benzo(g,h,i)perylene   | U                  |              | 0.00600         | 0.00600         |
| Benzo(k)fluoranthene   | U                  |              | 0.00600         | 0.00600         |
| Chrysene               | U                  |              | 0.00600         | 0.00600         |
| Dibenz(a,h)anthracene  | U                  |              | 0.00600         | 0.00600         |
| Fluoranthene           | U                  |              | 0.00600         | 0.00600         |
| Fluorene               | U                  |              | 0.00600         | 0.00600         |
| Indeno(1,2,3-cd)pyrene | U                  |              | 0.00600         | 0.00600         |
| Naphthalene            | U                  |              | 0.00200         | 0.0200          |
| Phenanthrene           | U                  |              | 0.00600         | 0.00600         |
| Pyrene                 | U                  |              | 0.00600         | 0.00600         |
| 1-Methylnaphthalene    | U                  |              | 0.00200         | 0.0200          |
| 2-Methylnaphthalene    | U                  |              | 0.00200         | 0.0200          |
| 2-Chloronaphthalene    | U                  |              | 0.00200         | 0.0200          |
| (S) Nitrobenzene-d5    | 89.8               |              |                 | 14.0-149        |
| (S) 2-Fluorobiphenyl   | 83.6               |              |                 | 34.0-125        |
| (S) p-Terphenyl-d14    | 76.3               |              |                 | 23.0-120        |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3356952-1 11/04/18 19:00 • (LCSD) R3356952-2 11/04/18 19:22

| Analyte               | Spike Amount<br>mg/kg | LCS Result<br>mg/kg | LCSD Result<br>mg/kg | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | LCS Qualifier | LCSD Qualifier | RPD<br>% | RPD Limits<br>% |
|-----------------------|-----------------------|---------------------|----------------------|---------------|----------------|------------------|---------------|----------------|----------|-----------------|
| Anthracene            | 0.0800                | 0.0662              | 0.0640               | 82.8          | 80.0           | 50.0-126         |               |                | 3.38     | 20              |
| Acenaphthene          | 0.0800                | 0.0643              | 0.0629               | 80.4          | 78.6           | 50.0-120         |               |                | 2.20     | 20              |
| Acenaphthylene        | 0.0800                | 0.0653              | 0.0638               | 81.6          | 79.8           | 50.0-120         |               |                | 2.32     | 20              |
| Benzo(a)anthracene    | 0.0800                | 0.0629              | 0.0606               | 78.6          | 75.8           | 45.0-120         |               |                | 3.72     | 20              |
| Benzo(a)pyrene        | 0.0800                | 0.0540              | 0.0506               | 67.5          | 63.3           | 42.0-120         |               |                | 6.50     | 20              |
| Benzo(b)fluoranthene  | 0.0800                | 0.0597              | 0.0616               | 74.6          | 77.0           | 42.0-121         |               |                | 3.13     | 20              |
| Benzo(g,h,i)perylene  | 0.0800                | 0.0621              | 0.0598               | 77.6          | 74.8           | 45.0-125         |               |                | 3.77     | 20              |
| Benzo(k)fluoranthene  | 0.0800                | 0.0685              | 0.0638               | 85.6          | 79.8           | 49.0-125         |               |                | 7.11     | 20              |
| Chrysene              | 0.0800                | 0.0674              | 0.0654               | 84.3          | 81.8           | 49.0-122         |               |                | 3.01     | 20              |
| Dibenz(a,h)anthracene | 0.0800                | 0.0626              | 0.0607               | 78.3          | 75.9           | 47.0-125         |               |                | 3.08     | 20              |
| Fluoranthene          | 0.0800                | 0.0723              | 0.0700               | 90.4          | 87.5           | 49.0-129         |               |                | 3.23     | 20              |



Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3356952-1 11/04/18 19:00 • (LCSD) R3356952-2 11/04/18 19:22

| Analyte                | Spike Amount<br>mg/kg | LCS Result<br>mg/kg | LCSD Result<br>mg/kg | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | <u>LCS Qualifier</u> | <u>LCSD Qualifier</u> | RPD<br>% | RPD Limits<br>% |
|------------------------|-----------------------|---------------------|----------------------|---------------|----------------|------------------|----------------------|-----------------------|----------|-----------------|
| Fluorene               | 0.0800                | 0.0636              | 0.0626               | 79.5          | 78.3           | 49.0-120         |                      |                       | 1.58     | 20              |
| Indeno(1,2,3-cd)pyrene | 0.0800                | 0.0633              | 0.0614               | 79.1          | 76.8           | 46.0-125         |                      |                       | 3.05     | 20              |
| Naphthalene            | 0.0800                | 0.0631              | 0.0614               | 78.9          | 76.8           | 50.0-120         |                      |                       | 2.73     | 20              |
| Phenanthrene           | 0.0800                | 0.0653              | 0.0629               | 81.6          | 78.6           | 47.0-120         |                      |                       | 3.74     | 20              |
| Pyrene                 | 0.0800                | 0.0608              | 0.0585               | 76.0          | 73.1           | 43.0-123         |                      |                       | 3.86     | 20              |
| 1-Methylnaphthalene    | 0.0800                | 0.0672              | 0.0658               | 84.0          | 82.3           | 51.0-121         |                      |                       | 2.11     | 20              |
| 2-Methylnaphthalene    | 0.0800                | 0.0623              | 0.0605               | 77.9          | 75.6           | 50.0-120         |                      |                       | 2.93     | 20              |
| 2-Chloronaphthalene    | 0.0800                | 0.0656              | 0.0645               | 82.0          | 80.6           | 50.0-120         |                      |                       | 1.69     | 20              |
| (S) Nitrobenzene-d5    |                       |                     |                      | 92.6          | 95.7           | 14.0-149         |                      |                       |          |                 |
| (S) 2-Fluorobiphenyl   |                       |                     |                      | 83.9          | 87.4           | 34.0-125         |                      |                       |          |                 |
| (S) p-Terphenyl-d14    |                       |                     |                      | 77.6          | 78.9           | 23.0-120         |                      |                       |          |                 |

L1039723-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1039723-01 11/05/18 01:42 • (MS) R3356952-4 11/05/18 02:04 • (MSD) R3356952-5 11/05/18 02:25

| Analyte                | Spike Amount<br>(dry)<br>mg/kg | Original Result<br>(dry)<br>mg/kg | MS Result (dry)<br>mg/kg | MSD Result<br>(dry)<br>mg/kg | MS Rec.<br>% | MSD Rec.<br>% | Dilution | Rec. Limits<br>% | <u>MS Qualifier</u> | <u>MSD Qualifier</u> | RPD<br>% | RPD Limits<br>% |
|------------------------|--------------------------------|-----------------------------------|--------------------------|------------------------------|--------------|---------------|----------|------------------|---------------------|----------------------|----------|-----------------|
| Anthracene             | 0.0932                         | ND                                | 0.0795                   | 0.0626                       | 85.4         | 67.3          | 1        | 10.0-145         |                     |                      | 23.8     | 30              |
| Acenaphthene           | 0.0932                         | ND                                | 0.0735                   | 0.0589                       | 78.9         | 63.3          | 1        | 14.0-127         |                     |                      | 22.0     | 27              |
| Acenaphthylene         | 0.0932                         | ND                                | 0.0752                   | 0.0600                       | 80.7         | 64.4          | 1        | 21.0-124         |                     |                      | 22.6     | 25              |
| Benzo(a)anthracene     | 0.0932                         | ND                                | 0.0718                   | 0.0554                       | 77.1         | 59.5          | 1        | 10.0-139         |                     |                      | 25.8     | 30              |
| Benzo(a)pyrene         | 0.0932                         | ND                                | 0.0738                   | 0.0583                       | 79.3         | 62.6          | 1        | 10.0-141         |                     |                      | 23.4     | 31              |
| Benzo(b)fluoranthene   | 0.0932                         | ND                                | 0.0687                   | 0.0571                       | 73.8         | 61.3          | 1        | 10.0-140         |                     |                      | 18.5     | 36              |
| Benzo(g,h,i)perylene   | 0.0932                         | ND                                | 0.0684                   | 0.0532                       | 73.4         | 57.1          | 1        | 10.0-140         |                     |                      | 24.9     | 33              |
| Benzo(k)fluoranthene   | 0.0932                         | ND                                | 0.0749                   | 0.0571                       | 80.4         | 61.3          | 1        | 10.0-137         |                     |                      | 27.0     | 31              |
| Chrysene               | 0.0932                         | ND                                | 0.0750                   | 0.0590                       | 80.5         | 63.4          | 1        | 10.0-145         |                     |                      | 23.8     | 30              |
| Dibenz(a,h)anthracene  | 0.0932                         | ND                                | 0.0701                   | 0.0541                       | 75.3         | 58.1          | 1        | 10.0-132         |                     |                      | 25.7     | 31              |
| Fluoranthene           | 0.0932                         | ND                                | 0.0837                   | 0.0653                       | 89.9         | 70.1          | 1        | 10.0-153         |                     |                      | 24.7     | 33              |
| Fluorene               | 0.0932                         | ND                                | 0.0722                   | 0.0579                       | 77.5         | 62.1          | 1        | 11.0-130         |                     |                      | 22.0     | 29              |
| Indeno(1,2,3-cd)pyrene | 0.0932                         | ND                                | 0.0696                   | 0.0540                       | 74.8         | 58.0          | 1        | 10.0-137         |                     |                      | 25.2     | 32              |
| Naphthalene            | 0.0932                         | ND                                | 0.0734                   | 0.0589                       | 78.8         | 63.3          | 1        | 10.0-135         |                     |                      | 21.8     | 27              |
| Phenanthrene           | 0.0932                         | ND                                | 0.0744                   | 0.0586                       | 79.9         | 62.9          | 1        | 10.0-144         |                     |                      | 23.8     | 31              |
| Pyrene                 | 0.0932                         | ND                                | 0.0695                   | 0.0548                       | 74.6         | 58.9          | 1        | 10.0-148         |                     |                      | 23.6     | 35              |
| 1-Methylnaphthalene    | 0.0932                         | ND                                | 0.0774                   | 0.0617                       | 83.1         | 66.3          | 1        | 10.0-142         |                     |                      | 22.6     | 28              |
| 2-Methylnaphthalene    | 0.0932                         | ND                                | 0.0711                   | 0.0566                       | 76.4         | 60.8          | 1        | 10.0-137         |                     |                      | 22.8     | 28              |
| 2-Chloronaphthalene    | 0.0932                         | ND                                | 0.0755                   | 0.0612                       | 81.0         | 65.8          | 1        | 29.0-120         |                     |                      | 20.8     | 24              |
| (S) Nitrobenzene-d5    |                                |                                   |                          |                              | 95.2         | 79.0          |          | 14.0-149         |                     |                      |          |                 |
| (S) 2-Fluorobiphenyl   |                                |                                   |                          |                              | 86.7         | 73.4          |          | 34.0-125         |                     |                      |          |                 |
| (S) p-Terphenyl-d14    |                                |                                   |                          |                              | 79.4         | 64.5          |          | 23.0-120         |                     |                      |          |                 |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc



Guide to Reading and Understanding Your Laboratory Report

The information below is designed to better explain the various terms used in your report of analytical results from the Laboratory. This is not intended as a comprehensive explanation, and if you have additional questions please contact your project representative.

Abbreviations and Definitions

|                              |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
|------------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| (dry)                        | Results are reported based on the dry weight of the sample. [this will only be present on a dry report basis for soils].                                                                                                                                                                                                                                                                                                                                                                                                   |
| MDL                          | Method Detection Limit.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    |
| ND                           | Not detected at the Reporting Limit (or MDL where applicable).                                                                                                                                                                                                                                                                                                                                                                                                                                                             |
| RDL                          | Reported Detection Limit.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                  |
| Rec.                         | Recovery.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                  |
| RPD                          | Relative Percent Difference.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                               |
| SDG                          | Sample Delivery Group.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                     |
| (S)                          | Surrogate (Surrogate Standard) - Analytes added to every blank, sample, Laboratory Control Sample/Duplicate and Matrix Spike/Duplicate; used to evaluate analytical efficiency by measuring recovery. Surrogates are not expected to be detected in all environmental media.                                                                                                                                                                                                                                               |
| U                            | Not detected at the Reporting Limit (or MDL where applicable).                                                                                                                                                                                                                                                                                                                                                                                                                                                             |
| Analyte                      | The name of the particular compound or analysis performed. Some Analyses and Methods will have multiple analytes reported.                                                                                                                                                                                                                                                                                                                                                                                                 |
| Dilution                     | If the sample matrix contains an interfering material, the sample preparation volume or weight values differ from the standard, or if concentrations of analytes in the sample are higher than the highest limit of concentration that the laboratory can accurately report, the sample may be diluted for analysis. If a value different than 1 is used in this field, the result reported has already been corrected for this factor.                                                                                    |
| Limits                       | These are the target % recovery ranges or % difference value that the laboratory has historically determined as normal for the method and analyte being reported. Successful QC Sample analysis will target all analytes recovered or duplicated within these ranges.                                                                                                                                                                                                                                                      |
| Original Sample              | The non-spiked sample in the prep batch used to determine the Relative Percent Difference (RPD) from a quality control sample. The Original Sample may not be included within the reported SDG.                                                                                                                                                                                                                                                                                                                            |
| Qualifier                    | This column provides a letter and/or number designation that corresponds to additional information concerning the result reported. If a Qualifier is present, a definition per Qualifier is provided within the Glossary and Definitions page and potentially a discussion of possible implications of the Qualifier in the Case Narrative if applicable.                                                                                                                                                                  |
| Result                       | The actual analytical final result (corrected for any sample specific characteristics) reported for your sample. If there was no measurable result returned for a specific analyte, the result in this column may state "ND" (Not Detected) or "BDL" (Below Detectable Levels). The information in the results column should always be accompanied by either an MDL (Method Detection Limit) or RDL (Reporting Detection Limit) that defines the lowest value that the laboratory could detect or report for this analyte. |
| Case Narrative (Cn)          | A brief discussion about the included sample results, including a discussion of any non-conformances to protocol observed either at sample receipt by the laboratory from the field or during the analytical process. If present, there will be a section in the Case Narrative to discuss the meaning of any data qualifiers used in the report.                                                                                                                                                                          |
| Quality Control Summary (Qc) | This section of the report includes the results of the laboratory quality control analyses required by procedure or analytical methods to assist in evaluating the validity of the results reported for your samples. These analyses are not being performed on your samples typically, but on laboratory generated material.                                                                                                                                                                                              |
| Sample Chain of Custody (Sc) | This is the document created in the field when your samples were initially collected. This is used to verify the time and date of collection, the person collecting the samples, and the analyses that the laboratory is requested to perform. This chain of custody also documents all persons (excluding commercial shippers) that have had control or possession of the samples from the time of collection until delivery to the laboratory for analysis.                                                              |
| Sample Results (Sr)          | This section of your report will provide the results of all testing performed on your samples. These results are provided by sample ID and are separated by the analyses performed on each sample. The header line of each analysis section for each sample will provide the name and method number for the analysis reported.                                                                                                                                                                                             |
| Sample Summary (Ss)          | This section of the Analytical Report defines the specific analyses performed for each sample ID, including the dates and times of preparation and/or analysis.                                                                                                                                                                                                                                                                                                                                                            |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

| Qualifier | Description                                                                                           |
|-----------|-------------------------------------------------------------------------------------------------------|
| J         | The identification of the analyte is acceptable; the reported value is an estimate.                   |
| J4        | The associated batch QC was outside the established quality control range for accuracy.               |
| J6        | The sample matrix interfered with the ability to make any accurate determination; spike value is low. |
| T8        | Sample(s) received past/too close to holding time expiration.                                         |



Pace National is the only environmental laboratory accredited/certified to support your work nationwide from one location. One phone call, one point of contact, one laboratory. No other lab is as accessible or prepared to handle your needs throughout the country. Our capacity and capability from our single location laboratory is comparable to the collective totals of the network laboratories in our industry. The most significant benefit to our one location design is the design of our laboratory campus. The model is conducive to accelerated productivity, decreasing turn-around time, and preventing cross contamination, thus protecting sample integrity. Our focus on premium quality and prompt service allows us to be YOUR LAB OF CHOICE.

\* Not all certifications held by the laboratory are applicable to the results reported in the attached report.  
 \* Accreditation is only applicable to the test methods specified on each scope of accreditation held by Pace National.

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

## State Accreditations

|                         |             |                             |                   |
|-------------------------|-------------|-----------------------------|-------------------|
| Alabama                 | 40660       | Nebraska                    | NE-OS-15-05       |
| Alaska                  | 17-026      | Nevada                      | TN-03-2002-34     |
| Arizona                 | AZ0612      | New Hampshire               | 2975              |
| Arkansas                | 88-0469     | New Jersey-NELAP            | TN002             |
| California              | 2932        | New Mexico <sup>1</sup>     | n/a               |
| Colorado                | TN00003     | New York                    | 11742             |
| Connecticut             | PH-0197     | North Carolina              | Env375            |
| Florida                 | E87487      | North Carolina <sup>1</sup> | DW21704           |
| Georgia                 | NELAP       | North Carolina <sup>3</sup> | 41                |
| Georgia <sup>1</sup>    | 923         | North Dakota                | R-140             |
| Idaho                   | TN00003     | Ohio-VAP                    | CL0069            |
| Illinois                | 200008      | Oklahoma                    | 9915              |
| Indiana                 | C-TN-01     | Oregon                      | TN200002          |
| Iowa                    | 364         | Pennsylvania                | 68-02979          |
| Kansas                  | E-10277     | Rhode Island                | LA000356          |
| Kentucky <sup>1,6</sup> | 90010       | South Carolina              | 84004             |
| Kentucky <sup>2</sup>   | 16          | South Dakota                | n/a               |
| Louisiana               | AI30792     | Tennessee <sup>1,4</sup>    | 2006              |
| Louisiana <sup>1</sup>  | LA180010    | Texas                       | T 104704245-17-14 |
| Maine                   | TN0002      | Texas <sup>5</sup>          | LAB0152           |
| Maryland                | 324         | Utah                        | TN00003           |
| Massachusetts           | M-TN003     | Vermont                     | VT2006            |
| Michigan                | 9958        | Virginia                    | 460132            |
| Minnesota               | 047-999-395 | Washington                  | C847              |
| Mississippi             | TN00003     | West Virginia               | 233               |
| Missouri                | 340         | Wisconsin                   | 9980939910        |
| Montana                 | CERT0086    | Wyoming                     | A2LA              |

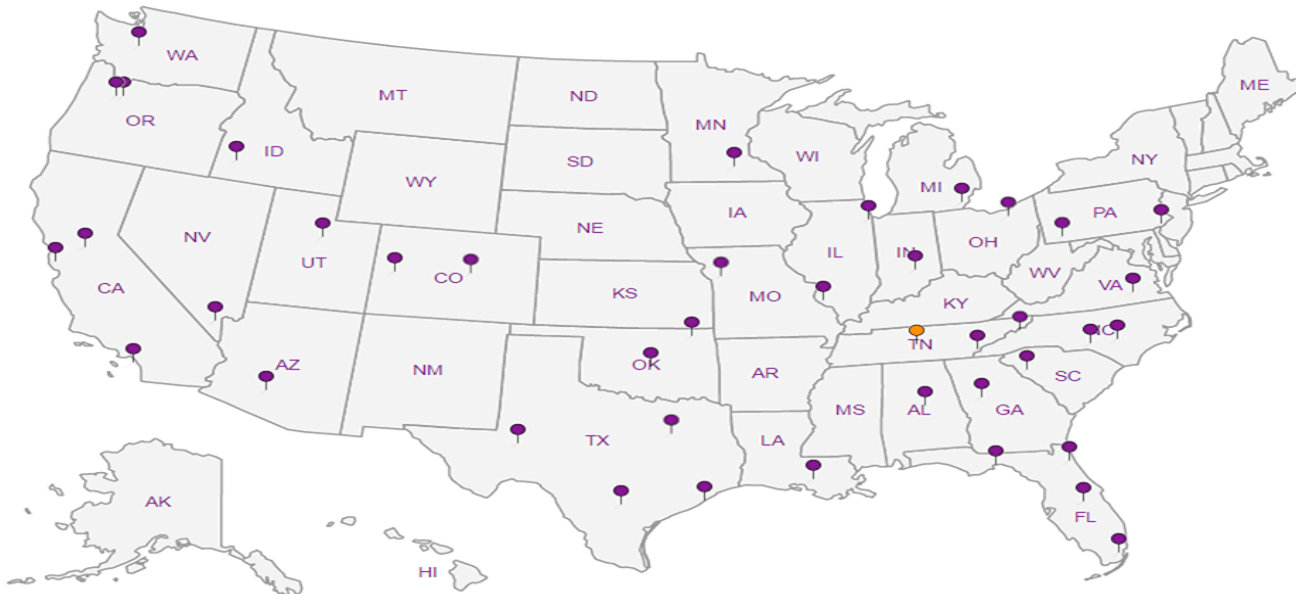
## Third Party Federal Accreditations

|                               |         |                    |               |
|-------------------------------|---------|--------------------|---------------|
| A2LA – ISO 17025              | 1461.01 | AIHA-LAP,LLC EMLAP | 100789        |
| A2LA – ISO 17025 <sup>5</sup> | 1461.02 | DOD                | 1461.01       |
| Canada                        | 1461.01 | USDA               | P330-15-00234 |
| EPA-Crypto                    | TN00003 |                    |               |

<sup>1</sup> Drinking Water <sup>2</sup> Underground Storage Tanks <sup>3</sup> Aquatic Toxicity <sup>4</sup> Chemical/Microbiological <sup>5</sup> Mold <sup>6</sup> Wastewater n/a Accreditation not applicable

## Our Locations

Pace National has sixty-four client support centers that provide sample pickup and/or the delivery of sampling supplies. If you would like assistance from one of our support offices, please contact our main office. Pace National performs all testing at our central laboratory.





November 09, 2018

## Terracon Consultants, Inc - Longmont, CO

Sample Delivery Group: L1040168  
Samples Received: 11/01/2018  
Project Number: 22187053  
Description: Union Reservior

Report To: Michael Skridulis  
1242 Bramwood Place  
Longmont, CO 80501

Entire Report Reviewed By:



Olivia Studebaker  
Project Manager

Results relate only to the items tested or calibrated and are reported as rounded values. This test report shall not be reproduced, except in full, without written approval of the laboratory. Where applicable, sampling conducted by Pace National is performed per guidance provided in laboratory standard operating procedures: 060302, 060303, and 060304.



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# SAMPLE SUMMARY



## MW-01 L1040168-01 GW

Collected by Charles Covington  
Collected date/time 10/31/18 11:55  
Received date/time 11/01/18 08:45

| Method                                                      | Batch     | Dilution | Preparation date/time | Analysis date/time | Analyst |
|-------------------------------------------------------------|-----------|----------|-----------------------|--------------------|---------|
| Calculated Results                                          | WG1190852 | 1        | 11/04/18 08:00        | 11/04/18 11:31     | ST      |
| Gravimetric Analysis by Method 2540 C-2011                  | WG1192072 | 1        | 11/07/18 17:26        | 11/07/18 18:20     | MMF     |
| Wet Chemistry by Method 4500CO2 D-2011                      | WG1191793 | 1        | 11/07/18 13:25        | 11/07/18 13:25     | GB      |
| Wet Chemistry by Method 7196A                               | WG1190723 | 1        | 11/02/18 18:30        | 11/02/18 18:30     | MLW     |
| Wet Chemistry by Method 9056A                               | WG1190237 | 10       | 11/02/18 23:35        | 11/02/18 23:35     | ELN     |
| Wet Chemistry by Method 9056A                               | WG1190237 | 500      | 11/03/18 10:40        | 11/03/18 10:40     | ELN     |
| Mercury by Method 7470A                                     | WG1191176 | 1        | 11/05/18 11:43        | 11/05/18 17:50     | TCT     |
| Metals (ICP) by Method 6010B                                | WG1190852 | 1        | 11/04/18 08:00        | 11/04/18 11:31     | ST      |
| Metals (ICP) by Method 6010B                                | WG1190863 | 1        | 11/03/18 15:06        | 11/05/18 13:27     | ST      |
| Volatile Organic Compounds (GC) by Method RSK175            | WG1190587 | 1        | 11/03/18 07:43        | 11/03/18 07:43     | MEL     |
| Volatile Organic Compounds (GC/MS) by Method 8260B          | WG1190163 | 1        | 11/02/18 03:19        | 11/02/18 03:19     | PP      |
| Semi Volatile Organic Compounds (GC/MS) by Method 8270C-SIM | WG1192224 | 1        | 11/07/18 07:25        | 11/07/18 16:39     | CJR     |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

## MW-02 L1040168-02 GW

Collected by Charles Covington  
Collected date/time 10/31/18 12:35  
Received date/time 11/01/18 08:45

| Method                                                      | Batch     | Dilution | Preparation date/time | Analysis date/time | Analyst |
|-------------------------------------------------------------|-----------|----------|-----------------------|--------------------|---------|
| Calculated Results                                          | WG1190852 | 1        | 11/04/18 08:00        | 11/04/18 11:45     | ST      |
| Gravimetric Analysis by Method 2540 C-2011                  | WG1192072 | 1        | 11/07/18 17:26        | 11/07/18 18:20     | MMF     |
| Wet Chemistry by Method 4500CO2 D-2011                      | WG1191793 | 1        | 11/07/18 13:34        | 11/07/18 13:34     | GB      |
| Wet Chemistry by Method 7196A                               | WG1190723 | 1        | 11/02/18 18:30        | 11/02/18 18:30     | MLW     |
| Wet Chemistry by Method 9056A                               | WG1190237 | 10       | 11/03/18 00:11        | 11/03/18 00:11     | ELN     |
| Wet Chemistry by Method 9056A                               | WG1190237 | 500      | 11/03/18 10:57        | 11/03/18 10:57     | ELN     |
| Mercury by Method 7470A                                     | WG1191176 | 1        | 11/05/18 11:43        | 11/05/18 17:52     | TCT     |
| Metals (ICP) by Method 6010B                                | WG1190852 | 1        | 11/04/18 08:00        | 11/04/18 11:45     | ST      |
| Metals (ICP) by Method 6010B                                | WG1190863 | 1        | 11/03/18 15:06        | 11/05/18 13:30     | ST      |
| Volatile Organic Compounds (GC) by Method RSK175            | WG1190587 | 1        | 11/03/18 07:47        | 11/03/18 07:47     | MEL     |
| Volatile Organic Compounds (GC/MS) by Method 8260B          | WG1190163 | 1        | 11/02/18 03:41        | 11/02/18 03:41     | PP      |
| Semi Volatile Organic Compounds (GC/MS) by Method 8270C-SIM | WG1192224 | 1        | 11/07/18 07:25        | 11/07/18 17:01     | CJR     |

## MW-03 L1040168-03 GW

Collected by Charles Covington  
Collected date/time 10/31/18 13:25  
Received date/time 11/01/18 08:45

| Method                                                      | Batch     | Dilution | Preparation date/time | Analysis date/time | Analyst |
|-------------------------------------------------------------|-----------|----------|-----------------------|--------------------|---------|
| Calculated Results                                          | WG1190852 | 1        | 11/04/18 08:00        | 11/04/18 11:48     | ST      |
| Gravimetric Analysis by Method 2540 C-2011                  | WG1192072 | 1        | 11/07/18 17:26        | 11/07/18 18:20     | MMF     |
| Wet Chemistry by Method 4500CO2 D-2011                      | WG1191793 | 1        | 11/07/18 13:42        | 11/07/18 13:42     | GB      |
| Wet Chemistry by Method 7196A                               | WG1190723 | 1        | 11/02/18 18:31        | 11/02/18 18:31     | MLW     |
| Wet Chemistry by Method 9056A                               | WG1190237 | 1000     | 11/03/18 11:42        | 11/03/18 11:42     | ELN     |
| Wet Chemistry by Method 9056A                               | WG1190237 | 20       | 11/03/18 00:48        | 11/03/18 00:48     | ELN     |
| Mercury by Method 7470A                                     | WG1191176 | 1        | 11/05/18 11:43        | 11/05/18 17:55     | TCT     |
| Metals (ICP) by Method 6010B                                | WG1190852 | 1        | 11/04/18 08:00        | 11/04/18 11:48     | ST      |
| Metals (ICP) by Method 6010B                                | WG1190863 | 10       | 11/03/18 15:06        | 11/05/18 22:38     | ST      |
| Volatile Organic Compounds (GC) by Method RSK175            | WG1190587 | 1        | 11/03/18 07:53        | 11/03/18 07:53     | MEL     |
| Volatile Organic Compounds (GC/MS) by Method 8260B          | WG1190163 | 1        | 11/02/18 04:03        | 11/02/18 04:03     | PP      |
| Semi Volatile Organic Compounds (GC/MS) by Method 8270C-SIM | WG1192224 | 1        | 11/07/18 07:25        | 11/07/18 17:23     | CJR     |



All sample aliquots were received at the correct temperature, in the proper containers, with the appropriate preservatives, and within method specified holding times, unless qualified or notated within the report. Where applicable, all MDL (LOD) and RDL (LOQ) values reported for environmental samples have been corrected for the dilution factor used in the analysis. All Method and Batch Quality Control are within established criteria except where addressed in this case narrative, a non-conformance form or properly qualified within the sample results. By my digital signature below, I affirm to the best of my knowledge, all problems/anomalies observed by the laboratory as having the potential to affect the quality of the data have been identified by the laboratory, and no information or data have been knowingly withheld that would affect the quality of the data.

Olivia Studebaker  
Project Manager

- <sup>1</sup> Cp
- <sup>2</sup> Tc
- <sup>3</sup> Ss
- <sup>4</sup> Cn
- <sup>5</sup> Sr
- <sup>6</sup> Qc
- <sup>7</sup> Gl
- <sup>8</sup> Al
- <sup>9</sup> Sc



Calculated Results

| Analyte             | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|---------------------|--------|-----------|--------|----------|----------------------|---------------------------|
| Chromium, Trivalent | ND     |           | 0.0100 | 1        | 11/04/2018 11:31     | <a href="#">WG1190852</a> |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Gravimetric Analysis by Method 2540 C-2011

| Analyte          | Result | Qualifier | RDL | Dilution | Analysis date / time | Batch                     |
|------------------|--------|-----------|-----|----------|----------------------|---------------------------|
| Dissolved Solids | 20000  |           | 200 | 1        | 11/07/2018 18:20     | <a href="#">WG1192072</a> |

Wet Chemistry by Method 4500CO2 D-2011

| Analyte             | Result | Qualifier | RDL  | Dilution | Analysis date / time | Batch                     |
|---------------------|--------|-----------|------|----------|----------------------|---------------------------|
| Free Carbon Dioxide | ND     | T8        | 20.0 | 1        | 11/07/2018 13:25     | <a href="#">WG1191793</a> |

Sample Narrative:

L1040168-01 WG1191793: Endpoint pH 4.5

Wet Chemistry by Method 7196A

| Analyte              | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|----------------------|--------|-----------|--------|----------|----------------------|---------------------------|
| Chromium, Hexavalent | ND     | T8        | 0.0100 | 1        | 11/02/2018 18:30     | <a href="#">WG1190723</a> |

Wet Chemistry by Method 9056A

| Analyte  | Result | Qualifier | RDL  | Dilution | Analysis date / time | Batch                     |
|----------|--------|-----------|------|----------|----------------------|---------------------------|
| Chloride | 357    |           | 10.0 | 10       | 11/02/2018 23:35     | <a href="#">WG1190237</a> |
| Sulfate  | 15500  |           | 2500 | 500      | 11/03/2018 10:40     | <a href="#">WG1190237</a> |

Mercury by Method 7470A

| Analyte            | Result | Qualifier | RDL      | Dilution | Analysis date / time | Batch                     |
|--------------------|--------|-----------|----------|----------|----------------------|---------------------------|
| Mercury, Dissolved | ND     |           | 0.000200 | 1        | 11/05/2018 17:50     | <a href="#">WG1191176</a> |

Metals (ICP) by Method 6010B

| Analyte             | Result | Qualifier | RDL     | Dilution | Analysis date / time | Batch                     |
|---------------------|--------|-----------|---------|----------|----------------------|---------------------------|
| Arsenic, Dissolved  | ND     |           | 0.0100  | 1        | 11/05/2018 13:27     | <a href="#">WG1190863</a> |
| Barium, Dissolved   | 0.0715 |           | 0.00500 | 1        | 11/05/2018 13:27     | <a href="#">WG1190863</a> |
| Boron, Dissolved    | 1.11   |           | 0.200   | 1        | 11/05/2018 13:27     | <a href="#">WG1190863</a> |
| Cadmium, Dissolved  | ND     |           | 0.00200 | 1        | 11/05/2018 13:27     | <a href="#">WG1190863</a> |
| Chromium            | ND     |           | 0.0100  | 1        | 11/04/2018 11:31     | <a href="#">WG1190852</a> |
| Chromium, Dissolved | ND     |           | 0.0100  | 1        | 11/05/2018 13:27     | <a href="#">WG1190863</a> |
| Copper, Dissolved   | 0.0167 |           | 0.0100  | 1        | 11/05/2018 13:27     | <a href="#">WG1190863</a> |
| Lead, Dissolved     | ND     |           | 0.00500 | 1        | 11/05/2018 13:27     | <a href="#">WG1190863</a> |
| Nickel, Dissolved   | 0.0105 |           | 0.0100  | 1        | 11/05/2018 13:27     | <a href="#">WG1190863</a> |
| Selenium, Dissolved | 0.0358 |           | 0.0100  | 1        | 11/05/2018 13:27     | <a href="#">WG1190863</a> |
| Silver, Dissolved   | ND     |           | 0.00500 | 1        | 11/05/2018 13:27     | <a href="#">WG1190863</a> |
| Zinc, Dissolved     | ND     |           | 0.0500  | 1        | 11/05/2018 13:27     | <a href="#">WG1190863</a> |

Volatile Organic Compounds (GC) by Method RSK175

| Analyte | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|---------|--------|-----------|--------|----------|----------------------|---------------------------|
| Methane | ND     |           | 0.0100 | 1        | 11/03/2018 07:43     | <a href="#">WG1190587</a> |
| Ethane  | ND     |           | 0.0130 | 1        | 11/03/2018 07:43     | <a href="#">WG1190587</a> |



Collected date/time: 10/31/18 11:55

L1040168

Volatile Organic Compounds (GC) by Method RSK175

| Analyte | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|---------|--------|-----------|--------|----------|----------------------|---------------------------|
| Ethene  | ND     |           | 0.0130 | 1        | 11/03/2018 07:43     | <a href="#">WG1190587</a> |

1 Cp

2 Tc

Volatile Organic Compounds (GC/MS) by Method 8260B

| Analyte                     | Result  | Qualifier | RDL     | Dilution | Analysis date / time | Batch                     |
|-----------------------------|---------|-----------|---------|----------|----------------------|---------------------------|
| Acetone                     | ND      |           | 0.0500  | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Acrolein                    | ND      |           | 0.0500  | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Acrylonitrile               | ND      |           | 0.0100  | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Benzene                     | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Bromobenzene                | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Bromodichloromethane        | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Bromoform                   | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Bromomethane                | ND      |           | 0.00500 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| n-Butylbenzene              | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| sec-Butylbenzene            | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| tert-Butylbenzene           | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Carbon tetrachloride        | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Chlorobenzene               | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Chlorodibromomethane        | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Chloroethane                | ND      |           | 0.00500 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Chloroform                  | ND      |           | 0.00500 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Chloromethane               | ND      |           | 0.00250 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 2-Chlorotoluene             | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 4-Chlorotoluene             | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,2-Dibromo-3-Chloropropane | ND      |           | 0.00500 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,2-Dibromoethane           | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Dibromomethane              | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,2-Dichlorobenzene         | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,3-Dichlorobenzene         | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,4-Dichlorobenzene         | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Dichlorodifluoromethane     | ND      |           | 0.00500 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,1-Dichloroethane          | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,2-Dichloroethane          | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,1-Dichloroethene          | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| cis-1,2-Dichloroethene      | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| trans-1,2-Dichloroethene    | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,2-Dichloropropane         | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,1-Dichloropropene         | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,3-Dichloropropane         | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| cis-1,3-Dichloropropene     | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| trans-1,3-Dichloropropene   | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 2,2-Dichloropropane         | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Di-isopropyl ether          | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Ethylbenzene                | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Hexachloro-1,3-butadiene    | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Isopropylbenzene            | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| p-Isopropyltoluene          | 0.00122 |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 2-Butanone (MEK)            | ND      |           | 0.0100  | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Methylene Chloride          | ND      |           | 0.00500 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 4-Methyl-2-pentanone (MIBK) | ND      |           | 0.0100  | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Methyl tert-butyl ether     | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Naphthalene                 | ND      |           | 0.00500 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| n-Propylbenzene             | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| Styrene                     | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,1,1,2-Tetrachloroethane   | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |
| 1,1,2,2-Tetrachloroethane   | ND      |           | 0.00100 | 1        | 11/02/2018 03:19     | <a href="#">WG1190163</a> |

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



Volatile Organic Compounds (GC/MS) by Method 8260B

| Analyte                        | Result<br>mg/l | Qualifier | RDL<br>mg/l | Dilution | Analysis<br>date / time | Batch                     |
|--------------------------------|----------------|-----------|-------------|----------|-------------------------|---------------------------|
| 1,1,2-Trichlorotrifluoroethane | ND             |           | 0.00100     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| Tetrachloroethene              | ND             |           | 0.00100     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| Toluene                        | ND             |           | 0.00100     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| 1,2,3-Trichlorobenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| 1,2,4-Trichlorobenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| 1,1,1-Trichloroethane          | ND             |           | 0.00100     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| 1,1,2-Trichloroethane          | ND             |           | 0.00100     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| Trichloroethene                | ND             |           | 0.00100     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| Trichlorofluoromethane         | ND             |           | 0.00500     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| 1,2,3-Trichloropropane         | ND             |           | 0.00250     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| 1,2,4-Trimethylbenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| 1,2,3-Trimethylbenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| 1,3,5-Trimethylbenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| Vinyl chloride                 | ND             |           | 0.00100     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| Xylenes, Total                 | ND             |           | 0.00300     | 1        | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| (S) Toluene-d8                 | 99.3           |           | 80.0-120    |          | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| (S) Dibromofluoromethane       | 108            |           | 75.0-120    |          | 11/02/2018 03:19        | <a href="#">WG1190163</a> |
| (S) 4-Bromofluorobenzene       | 106            |           | 77.0-126    |          | 11/02/2018 03:19        | <a href="#">WG1190163</a> |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Semi Volatile Organic Compounds (GC/MS) by Method 8270C-SIM

| Analyte                | Result<br>mg/l | Qualifier | RDL<br>mg/l | Dilution | Analysis<br>date / time | Batch                     |
|------------------------|----------------|-----------|-------------|----------|-------------------------|---------------------------|
| Anthracene             | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Acenaphthene           | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Acenaphthylene         | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Benzo(a)anthracene     | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Benzo(a)pyrene         | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Benzo(b)fluoranthene   | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Benzo(g,h,i)perylene   | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Benzo(k)fluoranthene   | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Chrysene               | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Dibenz(a,h)anthracene  | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Fluoranthene           | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Fluorene               | 0.0000583      |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Indeno(1,2,3-cd)pyrene | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Naphthalene            | ND             |           | 0.000250    | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Phenanthrene           | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| Pyrene                 | ND             |           | 0.0000500   | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| 1-Methylnaphthalene    | ND             |           | 0.000250    | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| 2-Methylnaphthalene    | ND             |           | 0.000250    | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| 2-Chloronaphthalene    | ND             |           | 0.000250    | 1        | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| (S) Nitrobenzene-d5    | 99.5           |           | 31.0-160    |          | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| (S) 2-Fluorobiphenyl   | 106            |           | 48.0-148    |          | 11/07/2018 16:39        | <a href="#">WG1192224</a> |
| (S) p-Terphenyl-d14    | 115            |           | 37.0-146    |          | 11/07/2018 16:39        | <a href="#">WG1192224</a> |



Calculated Results

| Analyte             | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|---------------------|--------|-----------|--------|----------|----------------------|---------------------------|
| Chromium, Trivalent | ND     |           | 0.0100 | 1        | 11/04/2018 11:45     | <a href="#">WG1190852</a> |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Gravimetric Analysis by Method 2540 C-2011

| Analyte          | Result | Qualifier | RDL | Dilution | Analysis date / time | Batch                     |
|------------------|--------|-----------|-----|----------|----------------------|---------------------------|
| Dissolved Solids | 19700  |           | 200 | 1        | 11/07/2018 18:20     | <a href="#">WG1192072</a> |

Wet Chemistry by Method 4500CO2 D-2011

| Analyte             | Result | Qualifier | RDL  | Dilution | Analysis date / time | Batch                     |
|---------------------|--------|-----------|------|----------|----------------------|---------------------------|
| Free Carbon Dioxide | ND     | T8        | 20.0 | 1        | 11/07/2018 13:34     | <a href="#">WG1191793</a> |

Sample Narrative:

L1040168-02 WG1191793: Endpoint pH 4.5

Wet Chemistry by Method 7196A

| Analyte              | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|----------------------|--------|-----------|--------|----------|----------------------|---------------------------|
| Chromium, Hexavalent | ND     | T8        | 0.0100 | 1        | 11/02/2018 18:30     | <a href="#">WG1190723</a> |

Wet Chemistry by Method 9056A

| Analyte  | Result | Qualifier | RDL  | Dilution | Analysis date / time | Batch                     |
|----------|--------|-----------|------|----------|----------------------|---------------------------|
| Chloride | 569    |           | 10.0 | 10       | 11/03/2018 00:11     | <a href="#">WG1190237</a> |
| Sulfate  | 14800  |           | 2500 | 500      | 11/03/2018 10:57     | <a href="#">WG1190237</a> |

Mercury by Method 7470A

| Analyte            | Result | Qualifier | RDL      | Dilution | Analysis date / time | Batch                     |
|--------------------|--------|-----------|----------|----------|----------------------|---------------------------|
| Mercury, Dissolved | ND     |           | 0.000200 | 1        | 11/05/2018 17:52     | <a href="#">WG1191176</a> |

Metals (ICP) by Method 6010B

| Analyte             | Result | Qualifier | RDL     | Dilution | Analysis date / time | Batch                     |
|---------------------|--------|-----------|---------|----------|----------------------|---------------------------|
| Arsenic, Dissolved  | ND     |           | 0.0100  | 1        | 11/05/2018 13:30     | <a href="#">WG1190863</a> |
| Barium, Dissolved   | 0.0447 |           | 0.00500 | 1        | 11/05/2018 13:30     | <a href="#">WG1190863</a> |
| Boron, Dissolved    | 1.03   |           | 0.200   | 1        | 11/05/2018 13:30     | <a href="#">WG1190863</a> |
| Cadmium, Dissolved  | ND     |           | 0.00200 | 1        | 11/05/2018 13:30     | <a href="#">WG1190863</a> |
| Chromium            | ND     |           | 0.0100  | 1        | 11/04/2018 11:45     | <a href="#">WG1190852</a> |
| Chromium, Dissolved | ND     |           | 0.0100  | 1        | 11/05/2018 13:30     | <a href="#">WG1190863</a> |
| Copper, Dissolved   | ND     |           | 0.0100  | 1        | 11/05/2018 13:30     | <a href="#">WG1190863</a> |
| Lead, Dissolved     | ND     |           | 0.00500 | 1        | 11/05/2018 13:30     | <a href="#">WG1190863</a> |
| Nickel, Dissolved   | ND     |           | 0.0100  | 1        | 11/05/2018 13:30     | <a href="#">WG1190863</a> |
| Selenium, Dissolved | 0.465  |           | 0.0100  | 1        | 11/05/2018 13:30     | <a href="#">WG1190863</a> |
| Silver, Dissolved   | ND     |           | 0.00500 | 1        | 11/05/2018 13:30     | <a href="#">WG1190863</a> |
| Zinc, Dissolved     | ND     |           | 0.0500  | 1        | 11/05/2018 13:30     | <a href="#">WG1190863</a> |

Volatile Organic Compounds (GC) by Method RSK175

| Analyte | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|---------|--------|-----------|--------|----------|----------------------|---------------------------|
| Methane | ND     |           | 0.0100 | 1        | 11/03/2018 07:47     | <a href="#">WG1190587</a> |
| Ethane  | ND     |           | 0.0130 | 1        | 11/03/2018 07:47     | <a href="#">WG1190587</a> |



Collected date/time: 10/31/18 12:35

L1040168

Volatile Organic Compounds (GC) by Method RSK175

| Analyte | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|---------|--------|-----------|--------|----------|----------------------|---------------------------|
| Ethene  | ND     |           | 0.0130 | 1        | 11/03/2018 07:47     | <a href="#">WG1190587</a> |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Volatile Organic Compounds (GC/MS) by Method 8260B

| Analyte                     | Result | Qualifier | RDL     | Dilution | Analysis date / time | Batch                     |
|-----------------------------|--------|-----------|---------|----------|----------------------|---------------------------|
| Acetone                     | ND     |           | 0.0500  | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Acrolein                    | ND     |           | 0.0500  | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Acrylonitrile               | ND     |           | 0.0100  | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Benzene                     | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Bromobenzene                | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Bromodichloromethane        | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Bromoform                   | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Bromomethane                | ND     |           | 0.00500 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| n-Butylbenzene              | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| sec-Butylbenzene            | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| tert-Butylbenzene           | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Carbon tetrachloride        | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Chlorobenzene               | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Chlorodibromomethane        | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Chloroethane                | ND     |           | 0.00500 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Chloroform                  | ND     |           | 0.00500 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Chloromethane               | ND     |           | 0.00250 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 2-Chlorotoluene             | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 4-Chlorotoluene             | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,2-Dibromo-3-Chloropropane | ND     |           | 0.00500 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,2-Dibromoethane           | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Dibromomethane              | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,2-Dichlorobenzene         | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,3-Dichlorobenzene         | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,4-Dichlorobenzene         | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Dichlorodifluoromethane     | ND     |           | 0.00500 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,1-Dichloroethane          | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,2-Dichloroethane          | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,1-Dichloroethene          | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| cis-1,2-Dichloroethene      | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| trans-1,2-Dichloroethene    | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,2-Dichloropropane         | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,1-Dichloropropene         | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,3-Dichloropropane         | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| cis-1,3-Dichloropropene     | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| trans-1,3-Dichloropropene   | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 2,2-Dichloropropane         | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Di-isopropyl ether          | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Ethylbenzene                | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Hexachloro-1,3-butadiene    | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Isopropylbenzene            | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| p-Isopropyltoluene          | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 2-Butanone (MEK)            | ND     |           | 0.0100  | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Methylene Chloride          | ND     |           | 0.00500 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 4-Methyl-2-pentanone (MIBK) | ND     |           | 0.0100  | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Methyl tert-butyl ether     | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Naphthalene                 | ND     |           | 0.00500 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| n-Propylbenzene             | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| Styrene                     | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,1,1,2-Tetrachloroethane   | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |
| 1,1,2,2-Tetrachloroethane   | ND     |           | 0.00100 | 1        | 11/02/2018 03:41     | <a href="#">WG1190163</a> |



Volatile Organic Compounds (GC/MS) by Method 8260B

| Analyte                        | Result<br>mg/l | Qualifier | RDL<br>mg/l | Dilution | Analysis<br>date / time | Batch                     |
|--------------------------------|----------------|-----------|-------------|----------|-------------------------|---------------------------|
| 1,1,2-Trichlorotrifluoroethane | ND             |           | 0.00100     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| Tetrachloroethene              | ND             |           | 0.00100     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| Toluene                        | ND             |           | 0.00100     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| 1,2,3-Trichlorobenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| 1,2,4-Trichlorobenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| 1,1,1-Trichloroethane          | ND             |           | 0.00100     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| 1,1,2-Trichloroethane          | ND             |           | 0.00100     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| Trichloroethene                | ND             |           | 0.00100     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| Trichlorofluoromethane         | ND             |           | 0.00500     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| 1,2,3-Trichloropropane         | ND             |           | 0.00250     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| 1,2,4-Trimethylbenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| 1,2,3-Trimethylbenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| 1,3,5-Trimethylbenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| Vinyl chloride                 | ND             |           | 0.00100     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| Xylenes, Total                 | ND             |           | 0.00300     | 1        | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| (S) Toluene-d8                 | 97.9           |           | 80.0-120    |          | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| (S) Dibromofluoromethane       | 107            |           | 75.0-120    |          | 11/02/2018 03:41        | <a href="#">WG1190163</a> |
| (S) 4-Bromofluorobenzene       | 110            |           | 77.0-126    |          | 11/02/2018 03:41        | <a href="#">WG1190163</a> |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Semi Volatile Organic Compounds (GC/MS) by Method 8270C-SIM

| Analyte                | Result<br>mg/l | Qualifier | RDL<br>mg/l | Dilution | Analysis<br>date / time | Batch                     |
|------------------------|----------------|-----------|-------------|----------|-------------------------|---------------------------|
| Anthracene             | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Acenaphthene           | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Acenaphthylene         | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Benzo(a)anthracene     | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Benzo(a)pyrene         | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Benzo(b)fluoranthene   | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Benzo(g,h,i)perylene   | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Benzo(k)fluoranthene   | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Chrysene               | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Dibenz(a,h)anthracene  | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Fluoranthene           | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Fluorene               | 0.0000647      |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Indeno(1,2,3-cd)pyrene | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Naphthalene            | ND             |           | 0.000250    | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Phenanthrene           | 0.0000827      |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| Pyrene                 | ND             |           | 0.0000500   | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| 1-Methylnaphthalene    | ND             |           | 0.000250    | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| 2-Methylnaphthalene    | ND             |           | 0.000250    | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| 2-Chloronaphthalene    | ND             |           | 0.000250    | 1        | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| (S) Nitrobenzene-d5    | 104            |           | 31.0-160    |          | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| (S) 2-Fluorobiphenyl   | 106            |           | 48.0-148    |          | 11/07/2018 17:01        | <a href="#">WG1192224</a> |
| (S) p-Terphenyl-d14    | 106            |           | 37.0-146    |          | 11/07/2018 17:01        | <a href="#">WG1192224</a> |





Calculated Results

| Analyte             | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|---------------------|--------|-----------|--------|----------|----------------------|---------------------------|
| Chromium, Trivalent | ND     |           | 0.0100 | 1        | 11/04/2018 11:48     | <a href="#">WG1190852</a> |

1 Cp

2 Tc

Gravimetric Analysis by Method 2540 C-2011

| Analyte          | Result | Qualifier | RDL | Dilution | Analysis date / time | Batch                     |
|------------------|--------|-----------|-----|----------|----------------------|---------------------------|
| Dissolved Solids | 77700  |           | 500 | 1        | 11/07/2018 18:20     | <a href="#">WG1192072</a> |

3 Ss

4 Cn

Wet Chemistry by Method 4500CO2 D-2011

| Analyte             | Result | Qualifier | RDL  | Dilution | Analysis date / time | Batch                     |
|---------------------|--------|-----------|------|----------|----------------------|---------------------------|
| Free Carbon Dioxide | 21.5   | T8        | 20.0 | 1        | 11/07/2018 13:42     | <a href="#">WG1191793</a> |

5 Sr

6 Qc

Sample Narrative:

L1040168-03 WG1191793: Endpoint pH 4.5

7 Gl

Wet Chemistry by Method 7196A

| Analyte              | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|----------------------|--------|-----------|--------|----------|----------------------|---------------------------|
| Chromium, Hexavalent | ND     | T8        | 0.0100 | 1        | 11/02/2018 18:31     | <a href="#">WG1190723</a> |

8 Al

9 Sc

Wet Chemistry by Method 9056A

| Analyte  | Result | Qualifier | RDL  | Dilution | Analysis date / time | Batch                     |
|----------|--------|-----------|------|----------|----------------------|---------------------------|
| Chloride | 1830   |           | 20.0 | 20       | 11/03/2018 00:48     | <a href="#">WG1190237</a> |
| Sulfate  | 50300  |           | 5000 | 1000     | 11/03/2018 11:42     | <a href="#">WG1190237</a> |

Mercury by Method 7470A

| Analyte            | Result | Qualifier | RDL      | Dilution | Analysis date / time | Batch                     |
|--------------------|--------|-----------|----------|----------|----------------------|---------------------------|
| Mercury, Dissolved | ND     |           | 0.000200 | 1        | 11/05/2018 17:55     | <a href="#">WG1191176</a> |

Metals (ICP) by Method 6010B

| Analyte             | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|---------------------|--------|-----------|--------|----------|----------------------|---------------------------|
| Arsenic, Dissolved  | ND     |           | 0.100  | 10       | 11/05/2018 22:38     | <a href="#">WG1190863</a> |
| Barium, Dissolved   | 0.101  |           | 0.0500 | 10       | 11/05/2018 22:38     | <a href="#">WG1190863</a> |
| Boron, Dissolved    | 2.95   |           | 2.00   | 10       | 11/05/2018 22:38     | <a href="#">WG1190863</a> |
| Cadmium, Dissolved  | ND     |           | 0.0200 | 10       | 11/05/2018 22:38     | <a href="#">WG1190863</a> |
| Chromium            | ND     |           | 0.0100 | 1        | 11/04/2018 11:48     | <a href="#">WG1190852</a> |
| Chromium, Dissolved | ND     |           | 0.100  | 10       | 11/05/2018 22:38     | <a href="#">WG1190863</a> |
| Copper, Dissolved   | ND     |           | 0.100  | 10       | 11/05/2018 22:38     | <a href="#">WG1190863</a> |
| Lead, Dissolved     | ND     |           | 0.0500 | 10       | 11/05/2018 22:38     | <a href="#">WG1190863</a> |
| Nickel, Dissolved   | ND     |           | 0.100  | 10       | 11/05/2018 22:38     | <a href="#">WG1190863</a> |
| Selenium, Dissolved | 1.24   |           | 0.100  | 10       | 11/05/2018 22:38     | <a href="#">WG1190863</a> |
| Silver, Dissolved   | ND     |           | 0.0500 | 10       | 11/05/2018 22:38     | <a href="#">WG1190863</a> |
| Zinc, Dissolved     | ND     |           | 0.500  | 10       | 11/05/2018 22:38     | <a href="#">WG1190863</a> |

Volatile Organic Compounds (GC) by Method RSK175

| Analyte | Result | Qualifier | RDL    | Dilution | Analysis date / time | Batch                     |
|---------|--------|-----------|--------|----------|----------------------|---------------------------|
| Methane | ND     |           | 0.0100 | 1        | 11/03/2018 07:53     | <a href="#">WG1190587</a> |
| Ethane  | ND     |           | 0.0130 | 1        | 11/03/2018 07:53     | <a href="#">WG1190587</a> |



Collected date/time: 10/31/18 13:25

L1040168

Volatile Organic Compounds (GC) by Method RSK175

| Analyte | Result | Qualifier | RDL    | Dilution | Analysis         | Batch                     |
|---------|--------|-----------|--------|----------|------------------|---------------------------|
|         | mg/l   |           | mg/l   |          | date / time      |                           |
| Ethene  | ND     |           | 0.0130 | 1        | 11/03/2018 07:53 | <a href="#">WG1190587</a> |

1 Cp

2 Tc

Volatile Organic Compounds (GC/MS) by Method 8260B

| Analyte                     | Result | Qualifier | RDL     | Dilution | Analysis         | Batch                     |
|-----------------------------|--------|-----------|---------|----------|------------------|---------------------------|
|                             | mg/l   |           | mg/l    |          | date / time      |                           |
| Acetone                     | ND     |           | 0.0500  | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Acrolein                    | ND     |           | 0.0500  | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Acrylonitrile               | ND     |           | 0.0100  | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Benzene                     | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Bromobenzene                | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Bromodichloromethane        | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Bromoform                   | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Bromomethane                | ND     |           | 0.00500 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| n-Butylbenzene              | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| sec-Butylbenzene            | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| tert-Butylbenzene           | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Carbon tetrachloride        | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Chlorobenzene               | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Chlorodibromomethane        | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Chloroethane                | ND     |           | 0.00500 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Chloroform                  | ND     |           | 0.00500 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Chloromethane               | ND     |           | 0.00250 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 2-Chlorotoluene             | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 4-Chlorotoluene             | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,2-Dibromo-3-Chloropropane | ND     |           | 0.00500 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,2-Dibromoethane           | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Dibromomethane              | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,2-Dichlorobenzene         | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,3-Dichlorobenzene         | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,4-Dichlorobenzene         | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Dichlorodifluoromethane     | ND     |           | 0.00500 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,1-Dichloroethane          | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,2-Dichloroethane          | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,1-Dichloroethene          | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| cis-1,2-Dichloroethene      | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| trans-1,2-Dichloroethene    | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,2-Dichloropropane         | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,1-Dichloropropene         | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,3-Dichloropropane         | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| cis-1,3-Dichloropropene     | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| trans-1,3-Dichloropropene   | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 2,2-Dichloropropane         | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Di-isopropyl ether          | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Ethylbenzene                | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Hexachloro-1,3-butadiene    | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Isopropylbenzene            | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| p-Isopropyltoluene          | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 2-Butanone (MEK)            | ND     |           | 0.0100  | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Methylene Chloride          | ND     |           | 0.00500 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 4-Methyl-2-pentanone (MIBK) | ND     |           | 0.0100  | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Methyl tert-butyl ether     | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Naphthalene                 | ND     |           | 0.00500 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| n-Propylbenzene             | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| Styrene                     | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,1,1,2-Tetrachloroethane   | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |
| 1,1,2,2-Tetrachloroethane   | ND     |           | 0.00100 | 1        | 11/02/2018 04:03 | <a href="#">WG1190163</a> |

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



Collected date/time: 10/31/18 13:25

L1040168

Volatile Organic Compounds (GC/MS) by Method 8260B

| Analyte                        | Result<br>mg/l | Qualifier | RDL<br>mg/l | Dilution | Analysis<br>date / time | Batch                     |
|--------------------------------|----------------|-----------|-------------|----------|-------------------------|---------------------------|
| 1,1,2-Trichlorotrifluoroethane | ND             |           | 0.00100     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| Tetrachloroethene              | ND             |           | 0.00100     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| Toluene                        | ND             |           | 0.00100     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| 1,2,3-Trichlorobenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| 1,2,4-Trichlorobenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| 1,1,1-Trichloroethane          | ND             |           | 0.00100     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| 1,1,2-Trichloroethane          | ND             |           | 0.00100     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| Trichloroethene                | ND             |           | 0.00100     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| Trichlorofluoromethane         | ND             |           | 0.00500     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| 1,2,3-Trichloropropane         | ND             |           | 0.00250     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| 1,2,4-Trimethylbenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| 1,2,3-Trimethylbenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| 1,3,5-Trimethylbenzene         | ND             |           | 0.00100     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| Vinyl chloride                 | ND             |           | 0.00100     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| Xylenes, Total                 | ND             |           | 0.00300     | 1        | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| (S) Toluene-d8                 | 95.2           |           | 80.0-120    |          | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| (S) Dibromofluoromethane       | 107            |           | 75.0-120    |          | 11/02/2018 04:03        | <a href="#">WG1190163</a> |
| (S) 4-Bromofluorobenzene       | 105            |           | 77.0-126    |          | 11/02/2018 04:03        | <a href="#">WG1190163</a> |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Semi Volatile Organic Compounds (GC/MS) by Method 8270C-SIM

| Analyte                | Result<br>mg/l | Qualifier | RDL<br>mg/l | Dilution | Analysis<br>date / time | Batch                     |
|------------------------|----------------|-----------|-------------|----------|-------------------------|---------------------------|
| Anthracene             | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Acenaphthene           | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Acenaphthylene         | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Benzo(a)anthracene     | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Benzo(a)pyrene         | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Benzo(b)fluoranthene   | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Benzo(g,h,i)perylene   | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Benzo(k)fluoranthene   | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Chrysene               | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Dibenz(a,h)anthracene  | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Fluoranthene           | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Fluorene               | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Indeno(1,2,3-cd)pyrene | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Naphthalene            | ND             |           | 0.000250    | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Phenanthrene           | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| Pyrene                 | ND             |           | 0.0000500   | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| 1-Methylnaphthalene    | ND             |           | 0.000250    | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| 2-Methylnaphthalene    | ND             |           | 0.000250    | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| 2-Chloronaphthalene    | ND             |           | 0.000250    | 1        | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| (S) Nitrobenzene-d5    | 109            |           | 31.0-160    |          | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| (S) 2-Fluorobiphenyl   | 99.5           |           | 48.0-148    |          | 11/07/2018 17:23        | <a href="#">WG1192224</a> |
| (S) p-Terphenyl-d14    | 108            |           | 37.0-146    |          | 11/07/2018 17:23        | <a href="#">WG1192224</a> |



Method Blank (MB)

(MB) R3358295-1 11/07/18 18:20

| Analyte          | MB Result<br>mg/l | MB Qualifier | MB MDL<br>mg/l | MB RDL<br>mg/l |
|------------------|-------------------|--------------|----------------|----------------|
| Dissolved Solids | U                 |              | 2.82           | 10.0           |

1 Cp

2 Tc

3 Ss

L1040168-01 Original Sample (OS) • Duplicate (DUP)

(OS) L1040168-01 11/07/18 18:20 • (DUP) R3358295-3 11/07/18 18:20

| Analyte          | Original Result<br>mg/l | DUP Result<br>mg/l | Dilution | DUP RPD<br>% | DUP Qualifier | DUP RPD<br>Limits<br>% |
|------------------|-------------------------|--------------------|----------|--------------|---------------|------------------------|
| Dissolved Solids | 20000                   | 19700              | 1        | 1.51         |               | 5                      |

4 Cn

5 Sr

Laboratory Control Sample (LCS)

(LCS) R3358295-2 11/07/18 18:20

| Analyte          | Spike Amount<br>mg/l | LCS Result<br>mg/l | LCS Rec.<br>% | Rec. Limits<br>% | LCS Qualifier |
|------------------|----------------------|--------------------|---------------|------------------|---------------|
| Dissolved Solids | 8800                 | 8530               | 96.9          | 85.0-115         |               |

6 Qc

7 Gl

8 Al

9 Sc



Method Blank (MB)

(MB) R3357760-2 11/06/18 17:06

| Analyte             | MB Result | MB Qualifier | MB MDL | MB RDL |
|---------------------|-----------|--------------|--------|--------|
| Free Carbon Dioxide | U         |              | 6.67   | 20.0   |

Sample Narrative:

BLANK: Endpoint pH 4.5

L1040178-01 Original Sample (OS) • Duplicate (DUP)

(OS) L1040178-01 11/06/18 17:15 • (DUP) R3357760-4 11/06/18 17:22

| Analyte             | Original Result | DUP Result | Dilution | DUP RPD | DUP Qualifier | DUP RPD Limits |
|---------------------|-----------------|------------|----------|---------|---------------|----------------|
| Free Carbon Dioxide |                 | ND         | 1        | 0.000   |               | 20             |

Sample Narrative:

OS: Endpoint pH 4.5 headspace

DUP: Endpoint pH 4.5

L1040178-04 Original Sample (OS) • Duplicate (DUP)

(OS) L1040178-04 11/06/18 20:03 • (DUP) R3357760-7 11/06/18 20:11

| Analyte             | Original Result | DUP Result | Dilution | DUP RPD | DUP Qualifier | DUP RPD Limits |
|---------------------|-----------------|------------|----------|---------|---------------|----------------|
| Free Carbon Dioxide |                 | ND         | 1        | 0.000   |               | 20             |

Sample Narrative:

OS: Endpoint pH 4.5 headspace

DUP: Endpoint pH 4.5

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



Method Blank (MB)

(MB) R3356474-1 11/02/18 18:16

| Analyte             | MB Result<br>mg/l | MB Qualifier | MB MDL<br>mg/l | MB RDL<br>mg/l |
|---------------------|-------------------|--------------|----------------|----------------|
| Chromium,Hexavalent | U                 |              | 0.00300        | 0.0100         |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc

L1040409-01 Original Sample (OS) • Duplicate (DUP)

(OS) L1040409-01 11/02/18 18:17 • (DUP) R3356474-3 11/02/18 18:20

| Analyte             | Original Result<br>mg/l | DUP Result<br>mg/l | Dilution | DUP RPD<br>% | DUP Qualifier | DUP RPD<br>Limits<br>% |
|---------------------|-------------------------|--------------------|----------|--------------|---------------|------------------------|
| Chromium,Hexavalent | ND                      | 0.0160             | 5        | 0.000        |               | 20                     |

Sample Narrative:

OS: Diluted due to matrix

Laboratory Control Sample (LCS)

(LCS) R3356474-2 11/02/18 18:16

| Analyte             | Spike Amount<br>mg/l | LCS Result<br>mg/l | LCS Rec.<br>% | Rec. Limits<br>% | LCS Qualifier |
|---------------------|----------------------|--------------------|---------------|------------------|---------------|
| Chromium,Hexavalent | 0.600                | 0.594              | 99.0          | 80.0-120         |               |

L1040409-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1040409-01 11/02/18 18:17 • (MS) R3356474-4 11/02/18 18:20 • (MSD) R3356474-5 11/02/18 18:21

| Analyte             | Spike Amount<br>mg/l | Original Result<br>mg/l | MS Result<br>mg/l | MSD Result<br>mg/l | MS Rec.<br>% | MSD Rec.<br>% | Dilution | Rec. Limits<br>% | MS Qualifier | MSD Qualifier | RPD<br>% | RPD Limits<br>% |
|---------------------|----------------------|-------------------------|-------------------|--------------------|--------------|---------------|----------|------------------|--------------|---------------|----------|-----------------|
| Chromium,Hexavalent | 0.500                | ND                      | 1.60              | 1.58               | 63.3         | 62.6          | 5        | 85.0-115         | <u>J6</u>    | <u>J6</u>     | 1.07     | 20              |

Sample Narrative:

OS: Diluted due to matrix



Method Blank (MB)

(MB) R3356614-1 11/02/18 18:32

| Analyte  | MB Result | MB Qualifier | MB MDL | MB RDL |
|----------|-----------|--------------|--------|--------|
|          | mg/l      |              | mg/l   | mg/l   |
| Chloride | 0.0668    | ↓            | 0.0519 | 1.00   |
| Sulfate  | 0.117     | ↓            | 0.0774 | 5.00   |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc

L1039944-01 Original Sample (OS) • Duplicate (DUP)

(OS) L1039944-01 11/02/18 20:33 • (DUP) R3356614-3 11/02/18 20:52

| Analyte  | Original Result | DUP Result | Dilution | DUP RPD | DUP Qualifier | DUP RPD Limits |
|----------|-----------------|------------|----------|---------|---------------|----------------|
|          | mg/l            | mg/l       |          | %       |               | %              |
| Chloride | 8.25            | 8.37       | 1        | 1.47    |               | 15             |
| Sulfate  | 19.9            | 20.1       | 1        | 0.819   |               | 15             |

L1040213-08 Original Sample (OS) • Duplicate (DUP)

(OS) L1040213-08 11/03/18 04:26 • (DUP) R3356614-6 11/03/18 04:44

| Analyte  | Original Result | DUP Result | Dilution | DUP RPD | DUP Qualifier | DUP RPD Limits |
|----------|-----------------|------------|----------|---------|---------------|----------------|
|          | mg/l            | mg/l       |          | %       |               | %              |
| Chloride | 4.33            | 4.30       | 1        | 0.679   |               | 15             |
| Sulfate  | 21.6            | 21.5       | 1        | 0.256   |               | 15             |

Laboratory Control Sample (LCS)

(LCS) R3356614-2 11/02/18 18:50

| Analyte  | Spike Amount | LCS Result | LCS Rec. | Rec. Limits | LCS Qualifier |
|----------|--------------|------------|----------|-------------|---------------|
|          | mg/l         | mg/l       | %        | %           |               |
| Chloride | 40.0         | 39.4       | 98.5     | 80.0-120    |               |
| Sulfate  | 40.0         | 39.6       | 99.1     | 80.0-120    |               |

L1039944-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1039944-01 11/02/18 20:33 • (MS) R3356614-4 11/02/18 21:10 • (MSD) R3356614-5 11/02/18 21:28

| Analyte  | Spike Amount | Original Result | MS Result | MSD Result | MS Rec. | MSD Rec. | Dilution | Rec. Limits | MS Qualifier | MSD Qualifier | RPD  | RPD Limits |
|----------|--------------|-----------------|-----------|------------|---------|----------|----------|-------------|--------------|---------------|------|------------|
|          | mg/l         | mg/l            | mg/l      | mg/l       | %       | %        |          | %           |              |               | %    | %          |
| Chloride | 50.0         | 8.25            | 60.3      | 62.5       | 104     | 108      | 1        | 80.0-120    |              |               | 3.55 | 15         |
| Sulfate  | 50.0         | 19.9            | 71.2      | 73.1       | 102     | 106      | 1        | 80.0-120    |              |               | 2.60 | 15         |



L1040213-08 Original Sample (OS) • Matrix Spike (MS)

(OS) L1040213-08 11/03/18 04:26 • (MS) R3356614-7 11/03/18 05:02

| Analyte  | Spike Amount<br>mg/l | Original Result<br>mg/l | MS Result<br>mg/l | MS Rec.<br>% | Dilution | Rec. Limits<br>% | <u>MS Qualifier</u> |
|----------|----------------------|-------------------------|-------------------|--------------|----------|------------------|---------------------|
| Chloride | 50.0                 | 4.33                    | 53.2              | 97.7         | 1        | 80.0-120         |                     |
| Sulfate  | 50.0                 | 21.6                    | 68.7              | 94.3         | 1        | 80.0-120         |                     |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc





Method Blank (MB)

(MB) R3357117-1 11/05/18 17:30

| Analyte           | MB Result<br>mg/l | MB Qualifier | MB MDL<br>mg/l | MB RDL<br>mg/l |
|-------------------|-------------------|--------------|----------------|----------------|
| Mercury,Dissolved | U                 |              | 0.0000490      | 0.000200       |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3357117-2 11/05/18 17:33 • (LCSD) R3357117-3 11/05/18 17:35

| Analyte           | Spike Amount<br>mg/l | LCS Result<br>mg/l | LCSD Result<br>mg/l | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | LCS Qualifier | LCSD Qualifier | RPD<br>% | RPD Limits<br>% |
|-------------------|----------------------|--------------------|---------------------|---------------|----------------|------------------|---------------|----------------|----------|-----------------|
| Mercury,Dissolved | 0.00300              | 0.00302            | 0.00289             | 101           | 96.3           | 80.0-120         |               |                | 4.58     | 20              |

L1040453-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1040453-01 11/05/18 17:38 • (MS) R3357117-4 11/05/18 17:45 • (MSD) R3357117-5 11/05/18 17:47

| Analyte           | Spike Amount<br>mg/l | Original Result<br>mg/l | MS Result<br>mg/l | MSD Result<br>mg/l | MS Rec.<br>% | MSD Rec.<br>% | Dilution | Rec. Limits<br>% | MS Qualifier | MSD Qualifier | RPD<br>% | RPD Limits<br>% |
|-------------------|----------------------|-------------------------|-------------------|--------------------|--------------|---------------|----------|------------------|--------------|---------------|----------|-----------------|
| Mercury,Dissolved | 0.00300              | U                       | 0.00303           | 0.00311            | 101          | 104           | 1        | 75.0-125         |              |               | 2.66     | 20              |

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc



Method Blank (MB)

(MB) R3356760-1 11/04/18 11:06

| Analyte  | MB Result<br>mg/l | MB Qualifier | MB MDL<br>mg/l | MB RDL<br>mg/l |
|----------|-------------------|--------------|----------------|----------------|
| Chromium | U                 |              | 0.00140        | 0.0100         |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3356760-2 11/04/18 11:09 • (LCSD) R3356760-3 11/04/18 11:11

| Analyte  | Spike Amount<br>mg/l | LCS Result<br>mg/l | LCSD Result<br>mg/l | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | LCS Qualifier | LCSD Qualifier | RPD<br>% | RPD Limits<br>% |
|----------|----------------------|--------------------|---------------------|---------------|----------------|------------------|---------------|----------------|----------|-----------------|
| Chromium | 1.00                 | 0.971              | 0.967               | 97.1          | 96.7           | 80.0-120         |               |                | 0.325    | 20              |

L1040345-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1040345-01 11/04/18 11:14 • (MS) R3356760-5 11/04/18 11:19 • (MSD) R3356760-6 11/04/18 11:22

| Analyte  | Spike Amount<br>mg/l | Original Result<br>mg/l | MS Result<br>mg/l | MSD Result<br>mg/l | MS Rec.<br>% | MSD Rec.<br>% | Dilution | Rec. Limits<br>% | MS Qualifier | MSD Qualifier | RPD<br>% | RPD Limits<br>% |
|----------|----------------------|-------------------------|-------------------|--------------------|--------------|---------------|----------|------------------|--------------|---------------|----------|-----------------|
| Chromium | 1.00                 | ND                      | 0.983             | 0.976              | 98.3         | 97.6          | 1        | 75.0-125         |              |               | 0.683    | 20              |

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc



Method Blank (MB)

(MB) R3357093-1 11/05/18 12:54

| Analyte            | MB Result | MB Qualifier | MB MDL   | MB RDL  |
|--------------------|-----------|--------------|----------|---------|
|                    | mg/l      |              | mg/l     | mg/l    |
| Arsenic,Dissolved  | U         |              | 0.00650  | 0.0100  |
| Barium,Dissolved   | U         |              | 0.00170  | 0.00500 |
| Boron,Dissolved    | U         |              | 0.0126   | 0.200   |
| Cadmium,Dissolved  | U         |              | 0.000700 | 0.00200 |
| Chromium,Dissolved | U         |              | 0.00140  | 0.0100  |
| Copper,Dissolved   | U         |              | 0.00530  | 0.0100  |
| Lead,Dissolved     | U         |              | 0.00190  | 0.00500 |
| Nickel,Dissolved   | U         |              | 0.00490  | 0.0100  |
| Selenium,Dissolved | U         |              | 0.00740  | 0.0100  |
| Silver,Dissolved   | U         |              | 0.00280  | 0.00500 |
| Zinc,Dissolved     | 0.00943   | ↓            | 0.00590  | 0.0500  |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3357093-2 11/05/18 12:56 • (LCSD) R3357093-3 11/05/18 12:59

| Analyte            | Spike Amount | LCS Result | LCSD Result | LCS Rec. | LCSD Rec. | Rec. Limits | LCS Qualifier | LCSD Qualifier | RPD   | RPD Limits |
|--------------------|--------------|------------|-------------|----------|-----------|-------------|---------------|----------------|-------|------------|
|                    | mg/l         | mg/l       | mg/l        | %        | %         | %           |               |                | %     | %          |
| Arsenic,Dissolved  | 1.00         | 1.02       | 1.01        | 102      | 101       | 80.0-120    |               |                | 1.08  | 20         |
| Barium,Dissolved   | 1.00         | 1.09       | 1.08        | 109      | 108       | 80.0-120    |               |                | 1.03  | 20         |
| Boron,Dissolved    | 1.00         | 1.04       | 1.04        | 104      | 104       | 80.0-120    |               |                | 0.151 | 20         |
| Cadmium,Dissolved  | 1.00         | 0.998      | 0.989       | 99.8     | 98.9      | 80.0-120    |               |                | 0.949 | 20         |
| Chromium,Dissolved | 1.00         | 1.00       | 0.986       | 100      | 98.6      | 80.0-120    |               |                | 1.44  | 20         |
| Copper,Dissolved   | 1.00         | 1.02       | 1.01        | 102      | 101       | 80.0-120    |               |                | 1.34  | 20         |
| Lead,Dissolved     | 1.00         | 1.06       | 1.05        | 106      | 105       | 80.0-120    |               |                | 1.17  | 20         |
| Nickel,Dissolved   | 1.00         | 1.05       | 1.04        | 105      | 104       | 80.0-120    |               |                | 1.15  | 20         |
| Selenium,Dissolved | 1.00         | 0.996      | 0.985       | 99.6     | 98.5      | 80.0-120    |               |                | 1.15  | 20         |
| Silver,Dissolved   | 0.200        | 0.187      | 0.185       | 93.4     | 92.4      | 80.0-120    |               |                | 1.14  | 20         |
| Zinc,Dissolved     | 1.00         | 1.02       | 1.01        | 102      | 101       | 80.0-120    |               |                | 1.07  | 20         |

L1040218-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1040218-01 11/05/18 13:02 • (MS) R3357093-5 11/05/18 13:07 • (MSD) R3357093-6 11/05/18 13:09

| Analyte            | Spike Amount | Original Result | MS Result | MSD Result | MS Rec. | MSD Rec. | Dilution | Rec. Limits | MS Qualifier | MSD Qualifier | RPD   | RPD Limits |
|--------------------|--------------|-----------------|-----------|------------|---------|----------|----------|-------------|--------------|---------------|-------|------------|
|                    | mg/l         | mg/l            | mg/l      | mg/l       | %       | %        |          | %           |              |               | %     | %          |
| Arsenic,Dissolved  | 1.00         | 0.0103          | 1.06      | 1.04       | 105     | 103      | 1        | 75.0-125    |              |               | 2.11  | 20         |
| Barium,Dissolved   | 1.00         | 0.709           | 1.77      | 1.75       | 106     | 104      | 1        | 75.0-125    |              |               | 1.32  | 20         |
| Boron,Dissolved    | 1.00         | ND              | 1.09      | 1.08       | 104     | 104      | 1        | 75.0-125    |              |               | 0.767 | 20         |
| Cadmium,Dissolved  | 1.00         | ND              | 1.02      | 0.999      | 102     | 99.9     | 1        | 75.0-125    |              |               | 2.05  | 20         |
| Chromium,Dissolved | 1.00         | ND              | 1.00      | 0.976      | 100     | 97.6     | 1        | 75.0-125    |              |               | 2.61  | 20         |



[L1040168-01,02,03](#)

L1040218-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1040218-01 11/05/18 13:02 • (MS) R3357093-5 11/05/18 13:07 • (MSD) R3357093-6 11/05/18 13:09

| Analyte            | Spike Amount<br>mg/l | Original Result<br>mg/l | MS Result<br>mg/l | MSD Result<br>mg/l | MS Rec.<br>% | MSD Rec.<br>% | Dilution | Rec. Limits<br>% | MS Qualifier | MSD Qualifier | RPD<br>% | RPD Limits<br>% |
|--------------------|----------------------|-------------------------|-------------------|--------------------|--------------|---------------|----------|------------------|--------------|---------------|----------|-----------------|
| Copper,Dissolved   | 1.00                 | ND                      | 1.04              | 1.02               | 104          | 102           | 1        | 75.0-125         |              |               | 2.21     | 20              |
| Lead,Dissolved     | 1.00                 | 0.00740                 | 1.07              | 1.05               | 106          | 104           | 1        | 75.0-125         |              |               | 2.18     | 20              |
| Nickel,Dissolved   | 1.00                 | ND                      | 1.05              | 1.03               | 105          | 103           | 1        | 75.0-125         |              |               | 1.94     | 20              |
| Selenium,Dissolved | 1.00                 | ND                      | 1.03              | 1.02               | 103          | 102           | 1        | 75.0-125         |              |               | 1.63     | 20              |
| Silver,Dissolved   | 0.200                | ND                      | 0.191             | 0.186              | 95.3         | 93.2          | 1        | 75.0-125         |              |               | 2.23     | 20              |
| Zinc,Dissolved     | 1.00                 | ND                      | 1.02              | 0.997              | 101          | 98.4          | 1        | 75.0-125         |              |               | 2.10     | 20              |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc



Method Blank (MB)

(MB) R3356511-1 11/03/18 07:17

| Analyte | MB Result | MB Qualifier | MB MDL  | MB RDL |
|---------|-----------|--------------|---------|--------|
|         | mg/l      |              | mg/l    | mg/l   |
| Methane | U         |              | 0.00291 | 0.0100 |
| Ethane  | U         |              | 0.00407 | 0.0130 |
| Ethene  | U         |              | 0.00426 | 0.0130 |

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

L1039927-01 Original Sample (OS) • Duplicate (DUP)

(OS) L1039927-01 11/03/18 07:27 • (DUP) R3356511-2 11/03/18 08:19

| Analyte | Original Result | DUP Result | Dilution | DUP RPD | DUP Qualifier | DUP RPD Limits |
|---------|-----------------|------------|----------|---------|---------------|----------------|
|         | mg/l            | mg/l       |          | %       |               | %              |
| Methane | 0.0714          | 0.0711     | 1        | 0.484   |               | 20             |
| Ethane  | ND              | 0.000      | 1        | 0.000   |               | 20             |
| Ethene  | ND              | 0.000      | 1        | 0.000   |               | 20             |

L1039951-01 Original Sample (OS) • Duplicate (DUP)

(OS) L1039951-01 11/03/18 07:41 • (DUP) R3356511-3 11/03/18 08:32

| Analyte | Original Result | DUP Result | Dilution | DUP RPD | DUP Qualifier | DUP RPD Limits |
|---------|-----------------|------------|----------|---------|---------------|----------------|
|         | mg/l            | mg/l       |          | %       |               | %              |
| Methane | 0.0466          | 0.0511     | 1        | 9.11    |               | 20             |
| Ethane  | ND              | 0.000      | 1        | 0.000   |               | 20             |
| Ethene  | ND              | 0.000      | 1        | 0.000   |               | 20             |

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3356511-4 11/03/18 09:12 • (LCSD) R3356511-5 11/03/18 09:16

| Analyte | Spike Amount | LCS Result | LCSD Result | LCS Rec. | LCSD Rec. | Rec. Limits | LCS Qualifier | LCSD Qualifier | RPD   | RPD Limits |
|---------|--------------|------------|-------------|----------|-----------|-------------|---------------|----------------|-------|------------|
|         | mg/l         | mg/l       | mg/l        | %        | %         | %           |               |                | %     | %          |
| Methane | 0.0678       | 0.0740     | 0.0733      | 109      | 108       | 85.0-115    |               |                | 0.867 | 20         |
| Ethane  | 0.129        | 0.114      | 0.114       | 88.1     | 88.6      | 85.0-115    |               |                | 0.581 | 20         |
| Ethene  | 0.127        | 0.114      | 0.113       | 89.7     | 89.0      | 85.0-115    |               |                | 0.738 | 20         |



Method Blank (MB)

(MB) R3357484-3 11/01/18 23:00

| Analyte                     | MB Result<br>mg/l | MB Qualifier | MB MDL<br>mg/l | MB RDL<br>mg/l |
|-----------------------------|-------------------|--------------|----------------|----------------|
| Acetone                     | U                 |              | 0.0100         | 0.0500         |
| Acrolein                    | U                 |              | 0.00887        | 0.0500         |
| Acrylonitrile               | U                 |              | 0.00187        | 0.0100         |
| Benzene                     | U                 |              | 0.000331       | 0.00100        |
| Bromobenzene                | U                 |              | 0.000352       | 0.00100        |
| Bromodichloromethane        | U                 |              | 0.000380       | 0.00100        |
| Bromoform                   | U                 |              | 0.000469       | 0.00100        |
| Bromomethane                | U                 |              | 0.000866       | 0.00500        |
| n-Butylbenzene              | U                 |              | 0.000361       | 0.00100        |
| sec-Butylbenzene            | U                 |              | 0.000365       | 0.00100        |
| tert-Butylbenzene           | U                 |              | 0.000399       | 0.00100        |
| Carbon tetrachloride        | U                 |              | 0.000379       | 0.00100        |
| Chlorobenzene               | U                 |              | 0.000348       | 0.00100        |
| Chlorodibromomethane        | U                 |              | 0.000327       | 0.00100        |
| Chloroethane                | U                 |              | 0.000453       | 0.00500        |
| Chloroform                  | U                 |              | 0.000324       | 0.00500        |
| Chloromethane               | U                 |              | 0.000276       | 0.00250        |
| 2-Chlorotoluene             | U                 |              | 0.000375       | 0.00100        |
| 4-Chlorotoluene             | U                 |              | 0.000351       | 0.00100        |
| 1,2-Dibromo-3-Chloropropane | U                 |              | 0.00133        | 0.00500        |
| 1,2-Dibromoethane           | U                 |              | 0.000381       | 0.00100        |
| Dibromomethane              | U                 |              | 0.000346       | 0.00100        |
| 1,2-Dichlorobenzene         | U                 |              | 0.000349       | 0.00100        |
| 1,3-Dichlorobenzene         | U                 |              | 0.000220       | 0.00100        |
| 1,4-Dichlorobenzene         | U                 |              | 0.000274       | 0.00100        |
| Dichlorodifluoromethane     | U                 |              | 0.000551       | 0.00500        |
| 1,1-Dichloroethane          | U                 |              | 0.000259       | 0.00100        |
| 1,2-Dichloroethane          | U                 |              | 0.000361       | 0.00100        |
| 1,1-Dichloroethene          | U                 |              | 0.000398       | 0.00100        |
| cis-1,2-Dichloroethene      | U                 |              | 0.000260       | 0.00100        |
| trans-1,2-Dichloroethene    | U                 |              | 0.000396       | 0.00100        |
| 1,2-Dichloropropane         | U                 |              | 0.000306       | 0.00100        |
| 1,1-Dichloropropene         | U                 |              | 0.000352       | 0.00100        |
| 1,3-Dichloropropane         | U                 |              | 0.000366       | 0.00100        |
| cis-1,3-Dichloropropene     | U                 |              | 0.000418       | 0.00100        |
| trans-1,3-Dichloropropene   | U                 |              | 0.000419       | 0.00100        |
| 2,2-Dichloropropane         | U                 |              | 0.000321       | 0.00100        |
| Di-isopropyl ether          | U                 |              | 0.000320       | 0.00100        |
| Ethylbenzene                | U                 |              | 0.000384       | 0.00100        |
| Hexachloro-1,3-butadiene    | 0.000498          | U            | 0.000256       | 0.00100        |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc



Method Blank (MB)

(MB) R3357484-3 11/01/18 23:00

| Analyte                        | MB Result<br>mg/l | MB Qualifier | MB MDL<br>mg/l | MB RDL<br>mg/l |
|--------------------------------|-------------------|--------------|----------------|----------------|
| Isopropylbenzene               | U                 |              | 0.000326       | 0.00100        |
| p-Isopropyltoluene             | U                 |              | 0.000350       | 0.00100        |
| 2-Butanone (MEK)               | U                 |              | 0.00393        | 0.0100         |
| Methylene Chloride             | U                 |              | 0.00100        | 0.00500        |
| 4-Methyl-2-pentanone (MIBK)    | U                 |              | 0.00214        | 0.0100         |
| Methyl tert-butyl ether        | U                 |              | 0.000367       | 0.00100        |
| Naphthalene                    | U                 |              | 0.00100        | 0.00500        |
| n-Propylbenzene                | U                 |              | 0.000349       | 0.00100        |
| Styrene                        | U                 |              | 0.000307       | 0.00100        |
| 1,1,1,2-Tetrachloroethane      | U                 |              | 0.000385       | 0.00100        |
| 1,1,2,2-Tetrachloroethane      | U                 |              | 0.000130       | 0.00100        |
| Tetrachloroethene              | U                 |              | 0.000372       | 0.00100        |
| Toluene                        | U                 |              | 0.000412       | 0.00100        |
| 1,1,2-Trichlorotrifluoroethane | U                 |              | 0.000303       | 0.00100        |
| 1,2,3-Trichlorobenzene         | 0.000516          | U            | 0.000230       | 0.00100        |
| 1,2,4-Trichlorobenzene         | 0.000364          | U            | 0.000355       | 0.00100        |
| 1,1,1-Trichloroethane          | U                 |              | 0.000319       | 0.00100        |
| 1,1,2-Trichloroethane          | U                 |              | 0.000383       | 0.00100        |
| Trichloroethene                | U                 |              | 0.000398       | 0.00100        |
| Trichlorofluoromethane         | U                 |              | 0.00120        | 0.00500        |
| 1,2,3-Trichloropropane         | U                 |              | 0.000807       | 0.00250        |
| 1,2,3-Trimethylbenzene         | U                 |              | 0.000321       | 0.00100        |
| 1,2,4-Trimethylbenzene         | U                 |              | 0.000373       | 0.00100        |
| 1,3,5-Trimethylbenzene         | U                 |              | 0.000387       | 0.00100        |
| Vinyl chloride                 | U                 |              | 0.000259       | 0.00100        |
| Xylenes, Total                 | U                 |              | 0.00106        | 0.00300        |
| (S) Toluene-d8                 | 95.5              |              |                | 80.0-120       |
| (S) Dibromofluoromethane       | 112               |              |                | 75.0-120       |
| (S) 4-Bromofluorobenzene       | 105               |              |                | 77.0-126       |

1  
Cp

2  
Tc

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Ss

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Al

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Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3357484-1 11/01/18 21:54 • (LCSD) R3357484-2 11/01/18 22:16

| Analyte       | Spike Amount<br>mg/l | LCS Result<br>mg/l | LCSD Result<br>mg/l | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | LCS Qualifier | LCSD Qualifier | RPD<br>% | RPD Limits<br>% |
|---------------|----------------------|--------------------|---------------------|---------------|----------------|------------------|---------------|----------------|----------|-----------------|
| Acetone       | 0.125                | 0.137              | 0.140               | 109           | 112            | 19.0-160         |               |                | 2.61     | 27              |
| Acrolein      | 0.125                | 0.0890             | 0.0895              | 71.2          | 71.6           | 10.0-160         |               |                | 0.512    | 26              |
| Acrylonitrile | 0.125                | 0.136              | 0.137               | 109           | 109            | 55.0-149         |               |                | 0.487    | 20              |
| Benzene       | 0.0250               | 0.0236             | 0.0236              | 94.2          | 94.2           | 70.0-123         |               |                | 0.0338   | 20              |



Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3357484-1 11/01/18 21:54 • (LCSD) R3357484-2 11/01/18 22:16

| Analyte                     | Spike Amount<br>mg/l | LCS Result<br>mg/l | LCSD Result<br>mg/l | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | <u>LCS Qualifier</u> | <u>LCSD Qualifier</u> | RPD<br>% | RPD Limits<br>% |
|-----------------------------|----------------------|--------------------|---------------------|---------------|----------------|------------------|----------------------|-----------------------|----------|-----------------|
| Bromobenzene                | 0.0250               | 0.0235             | 0.0255              | 94.1          | 102            | 73.0-121         |                      |                       | 8.03     | 20              |
| Bromodichloromethane        | 0.0250               | 0.0255             | 0.0269              | 102           | 108            | 75.0-120         |                      |                       | 5.40     | 20              |
| Bromoform                   | 0.0250               | 0.0251             | 0.0276              | 101           | 110            | 68.0-132         |                      |                       | 9.15     | 20              |
| Bromomethane                | 0.0250               | 0.0234             | 0.0240              | 93.5          | 96.2           | 10.0-160         |                      |                       | 2.85     | 25              |
| n-Butylbenzene              | 0.0250               | 0.0224             | 0.0233              | 89.6          | 93.1           | 73.0-125         |                      |                       | 3.93     | 20              |
| sec-Butylbenzene            | 0.0250               | 0.0229             | 0.0242              | 91.8          | 96.7           | 75.0-125         |                      |                       | 5.29     | 20              |
| tert-Butylbenzene           | 0.0250               | 0.0233             | 0.0249              | 93.1          | 99.8           | 76.0-124         |                      |                       | 6.88     | 20              |
| Carbon tetrachloride        | 0.0250               | 0.0259             | 0.0270              | 103           | 108            | 68.0-126         |                      |                       | 4.39     | 20              |
| Chlorobenzene               | 0.0250               | 0.0221             | 0.0231              | 88.4          | 92.5           | 80.0-121         |                      |                       | 4.52     | 20              |
| Chlorodibromomethane        | 0.0250               | 0.0243             | 0.0252              | 97.2          | 101            | 77.0-125         |                      |                       | 3.53     | 20              |
| Chloroethane                | 0.0250               | 0.0246             | 0.0251              | 98.3          | 100            | 47.0-150         |                      |                       | 2.10     | 20              |
| Chloroform                  | 0.0250               | 0.0255             | 0.0258              | 102           | 103            | 73.0-120         |                      |                       | 1.32     | 20              |
| Chloromethane               | 0.0250               | 0.0266             | 0.0267              | 107           | 107            | 41.0-142         |                      |                       | 0.0504   | 20              |
| 2-Chlorotoluene             | 0.0250               | 0.0239             | 0.0254              | 95.5          | 102            | 76.0-123         |                      |                       | 6.19     | 20              |
| 4-Chlorotoluene             | 0.0250               | 0.0238             | 0.0258              | 95.3          | 103            | 75.0-122         |                      |                       | 7.90     | 20              |
| 1,2-Dibromo-3-Chloropropane | 0.0250               | 0.0212             | 0.0233              | 84.8          | 93.1           | 58.0-134         |                      |                       | 9.39     | 20              |
| 1,2-Dibromoethane           | 0.0250               | 0.0229             | 0.0237              | 91.8          | 94.7           | 80.0-122         |                      |                       | 3.17     | 20              |
| Dibromomethane              | 0.0250               | 0.0247             | 0.0254              | 98.9          | 102            | 80.0-120         |                      |                       | 2.70     | 20              |
| 1,2-Dichlorobenzene         | 0.0250               | 0.0217             | 0.0225              | 86.9          | 90.0           | 79.0-121         |                      |                       | 3.51     | 20              |
| 1,3-Dichlorobenzene         | 0.0250               | 0.0218             | 0.0232              | 87.1          | 92.9           | 79.0-120         |                      |                       | 6.44     | 20              |
| 1,4-Dichlorobenzene         | 0.0250               | 0.0216             | 0.0227              | 86.4          | 90.6           | 79.0-120         |                      |                       | 4.79     | 20              |
| Dichlorodifluoromethane     | 0.0250               | 0.0261             | 0.0268              | 104           | 107            | 51.0-149         |                      |                       | 2.64     | 20              |
| 1,1-Dichloroethane          | 0.0250               | 0.0258             | 0.0264              | 103           | 106            | 70.0-126         |                      |                       | 2.26     | 20              |
| 1,2-Dichloroethane          | 0.0250               | 0.0280             | 0.0285              | 112           | 114            | 70.0-128         |                      |                       | 1.64     | 20              |
| 1,1-Dichloroethene          | 0.0250               | 0.0232             | 0.0237              | 92.9          | 94.7           | 71.0-124         |                      |                       | 1.92     | 20              |
| cis-1,2-Dichloroethene      | 0.0250               | 0.0230             | 0.0238              | 92.2          | 95.1           | 73.0-120         |                      |                       | 3.15     | 20              |
| trans-1,2-Dichloroethene    | 0.0250               | 0.0246             | 0.0251              | 98.4          | 100            | 73.0-120         |                      |                       | 2.05     | 20              |
| 1,2-Dichloropropane         | 0.0250               | 0.0238             | 0.0239              | 95.3          | 95.7           | 77.0-125         |                      |                       | 0.442    | 20              |
| 1,1-Dichloropropene         | 0.0250               | 0.0248             | 0.0252              | 99.3          | 101            | 74.0-126         |                      |                       | 1.46     | 20              |
| 1,3-Dichloropropane         | 0.0250               | 0.0228             | 0.0232              | 91.0          | 92.8           | 80.0-120         |                      |                       | 1.90     | 20              |
| cis-1,3-Dichloropropene     | 0.0250               | 0.0228             | 0.0237              | 91.1          | 94.7           | 80.0-123         |                      |                       | 3.85     | 20              |
| trans-1,3-Dichloropropene   | 0.0250               | 0.0235             | 0.0246              | 94.1          | 98.5           | 78.0-124         |                      |                       | 4.56     | 20              |
| 2,2-Dichloropropane         | 0.0250               | 0.0263             | 0.0261              | 105           | 105            | 58.0-130         |                      |                       | 0.686    | 20              |
| Di-isopropyl ether          | 0.0250               | 0.0271             | 0.0274              | 108           | 110            | 58.0-138         |                      |                       | 1.11     | 20              |
| Ethylbenzene                | 0.0250               | 0.0222             | 0.0227              | 88.8          | 90.9           | 79.0-123         |                      |                       | 2.43     | 20              |
| Hexachloro-1,3-butadiene    | 0.0250               | 0.0192             | 0.0221              | 76.9          | 88.3           | 54.0-138         |                      |                       | 13.8     | 20              |
| Isopropylbenzene            | 0.0250               | 0.0236             | 0.0258              | 94.2          | 103            | 76.0-127         |                      |                       | 9.08     | 20              |
| p-Isopropyltoluene          | 0.0250               | 0.0230             | 0.0240              | 91.9          | 96.1           | 76.0-125         |                      |                       | 4.56     | 20              |
| 2-Butanone (MEK)            | 0.125                | 0.141              | 0.140               | 113           | 112            | 44.0-160         |                      |                       | 0.613    | 20              |
| Methylene Chloride          | 0.0250               | 0.0229             | 0.0238              | 91.7          | 95.3           | 67.0-120         |                      |                       | 3.82     | 20              |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc





Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3357484-1 11/01/18 21:54 • (LCSD) R3357484-2 11/01/18 22:16

| Analyte                        | Spike Amount<br>mg/l | LCS Result<br>mg/l | LCSD Result<br>mg/l | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | LCS Qualifier | LCSD Qualifier | RPD<br>% | RPD Limits<br>% |
|--------------------------------|----------------------|--------------------|---------------------|---------------|----------------|------------------|---------------|----------------|----------|-----------------|
| 4-Methyl-2-pentanone (MIBK)    | 0.125                | 0.133              | 0.137               | 107           | 110            | 68.0-142         |               |                | 2.62     | 20              |
| Methyl tert-butyl ether        | 0.0250               | 0.0245             | 0.0252              | 97.9          | 101            | 68.0-125         |               |                | 2.82     | 20              |
| Naphthalene                    | 0.0250               | 0.0197             | 0.0216              | 78.7          | 86.5           | 54.0-135         |               |                | 9.51     | 20              |
| n-Propylbenzene                | 0.0250               | 0.0235             | 0.0252              | 94.2          | 101            | 77.0-124         |               |                | 6.98     | 20              |
| Styrene                        | 0.0250               | 0.0240             | 0.0263              | 95.8          | 105            | 73.0-130         |               |                | 9.33     | 20              |
| 1,1,1,2-Tetrachloroethane      | 0.0250               | 0.0233             | 0.0246              | 93.2          | 98.4           | 75.0-125         |               |                | 5.40     | 20              |
| 1,1,2,2-Tetrachloroethane      | 0.0250               | 0.0231             | 0.0246              | 92.3          | 98.2           | 65.0-130         |               |                | 6.29     | 20              |
| Tetrachloroethene              | 0.0250               | 0.0209             | 0.0213              | 83.4          | 85.1           | 72.0-132         |               |                | 1.99     | 20              |
| Toluene                        | 0.0250               | 0.0219             | 0.0226              | 87.7          | 90.3           | 79.0-120         |               |                | 2.91     | 20              |
| 1,1,2-Trichlorotrifluoroethane | 0.0250               | 0.0246             | 0.0256              | 98.2          | 102            | 69.0-132         |               |                | 4.10     | 20              |
| 1,2,3-Trichlorobenzene         | 0.0250               | 0.0184             | 0.0213              | 73.7          | 85.4           | 50.0-138         |               |                | 14.7     | 20              |
| 1,2,4-Trichlorobenzene         | 0.0250               | 0.0187             | 0.0209              | 74.7          | 83.8           | 57.0-137         |               |                | 11.5     | 20              |
| 1,1,1-Trichloroethane          | 0.0250               | 0.0275             | 0.0280              | 110           | 112            | 73.0-124         |               |                | 1.73     | 20              |
| 1,1,2-Trichloroethane          | 0.0250               | 0.0224             | 0.0228              | 89.8          | 91.1           | 80.0-120         |               |                | 1.50     | 20              |
| Trichloroethene                | 0.0250               | 0.0228             | 0.0237              | 91.4          | 94.8           | 78.0-124         |               |                | 3.71     | 20              |
| Trichlorofluoromethane         | 0.0250               | 0.0294             | 0.0299              | 118           | 120            | 59.0-147         |               |                | 1.63     | 20              |
| 1,2,3-Trichloropropane         | 0.0250               | 0.0256             | 0.0274              | 102           | 109            | 73.0-130         |               |                | 6.82     | 20              |
| 1,2,3-Trimethylbenzene         | 0.0250               | 0.0229             | 0.0245              | 91.7          | 98.0           | 77.0-120         |               |                | 6.60     | 20              |
| 1,2,4-Trimethylbenzene         | 0.0250               | 0.0235             | 0.0245              | 94.0          | 98.1           | 76.0-121         |               |                | 4.23     | 20              |
| 1,3,5-Trimethylbenzene         | 0.0250               | 0.0241             | 0.0256              | 96.3          | 103            | 76.0-122         |               |                | 6.35     | 20              |
| Vinyl chloride                 | 0.0250               | 0.0252             | 0.0254              | 101           | 102            | 67.0-131         |               |                | 0.896    | 20              |
| Xylenes, Total                 | 0.0750               | 0.0663             | 0.0684              | 88.4          | 91.2           | 79.0-123         |               |                | 3.12     | 20              |
| (S) Toluene-d8                 |                      |                    |                     | 96.0          | 98.1           | 80.0-120         |               |                |          |                 |
| (S) Dibromofluoromethane       |                      |                    |                     | 108           | 108            | 75.0-120         |               |                |          |                 |
| (S) 4-Bromofluorobenzene       |                      |                    |                     | 102           | 110            | 77.0-126         |               |                |          |                 |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc



Method Blank (MB)

(MB) R3357755-3 11/07/18 13:23

| Analyte                | MB Result<br>mg/l | MB Qualifier | MB MDL<br>mg/l | MB RDL<br>mg/l |
|------------------------|-------------------|--------------|----------------|----------------|
| Anthracene             | U                 |              | 0.0000140      | 0.0000500      |
| Acenaphthene           | U                 |              | 0.0000100      | 0.0000500      |
| Acenaphthylene         | U                 |              | 0.0000120      | 0.0000500      |
| Benzo(a)anthracene     | U                 |              | 0.00000410     | 0.0000500      |
| Benzo(a)pyrene         | U                 |              | 0.0000116      | 0.0000500      |
| Benzo(b)fluoranthene   | U                 |              | 0.00000212     | 0.0000500      |
| Benzo(g,h,i)perylene   | U                 |              | 0.00000227     | 0.0000500      |
| Benzo(k)fluoranthene   | U                 |              | 0.0000136      | 0.0000500      |
| Chrysene               | U                 |              | 0.0000108      | 0.0000500      |
| Dibenz(a,h)anthracene  | U                 |              | 0.00000396     | 0.0000500      |
| Fluoranthene           | U                 |              | 0.0000157      | 0.0000500      |
| Fluorene               | U                 |              | 0.00000850     | 0.0000500      |
| Indeno(1,2,3-cd)pyrene | U                 |              | 0.0000148      | 0.0000500      |
| Naphthalene            | U                 |              | 0.0000198      | 0.000250       |
| Phenanthrene           | U                 |              | 0.00000820     | 0.0000500      |
| Pyrene                 | U                 |              | 0.0000117      | 0.0000500      |
| 1-Methylnaphthalene    | U                 |              | 0.00000821     | 0.000250       |
| 2-Methylnaphthalene    | U                 |              | 0.00000902     | 0.000250       |
| 2-Chloronaphthalene    | U                 |              | 0.00000647     | 0.000250       |
| (S) Nitrobenzene-d5    | 142               |              |                | 31.0-160       |
| (S) 2-Fluorobiphenyl   | 103               |              |                | 48.0-148       |
| (S) p-Terphenyl-d14    | 108               |              |                | 37.0-146       |

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3357755-1 11/07/18 12:39 • (LCSD) R3357755-2 11/07/18 13:01

| Analyte               | Spike Amount<br>mg/l | LCS Result<br>mg/l | LCSD Result<br>mg/l | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | LCS Qualifier | LCSD Qualifier | RPD<br>% | RPD Limits<br>% |
|-----------------------|----------------------|--------------------|---------------------|---------------|----------------|------------------|---------------|----------------|----------|-----------------|
| Anthracene            | 0.00200              | 0.00203            | 0.00213             | 102           | 106            | 67.0-150         |               |                | 4.81     | 20              |
| Acenaphthene          | 0.00200              | 0.00186            | 0.00191             | 93.0          | 95.5           | 65.0-138         |               |                | 2.65     | 20              |
| Acenaphthylene        | 0.00200              | 0.00197            | 0.00202             | 98.5          | 101            | 66.0-140         |               |                | 2.51     | 20              |
| Benzo(a)anthracene    | 0.00200              | 0.00202            | 0.00208             | 101           | 104            | 61.0-140         |               |                | 2.93     | 20              |
| Benzo(a)pyrene        | 0.00200              | 0.00210            | 0.00199             | 105           | 99.5           | 60.0-143         |               |                | 5.38     | 20              |
| Benzo(b)fluoranthene  | 0.00200              | 0.00205            | 0.00190             | 102           | 95.0           | 58.0-141         |               |                | 7.59     | 20              |
| Benzo(g,h,i)perylene  | 0.00200              | 0.00214            | 0.00242             | 107           | 121            | 52.0-153         |               |                | 12.3     | 20              |
| Benzo(k)fluoranthene  | 0.00200              | 0.00202            | 0.00187             | 101           | 93.5           | 58.0-148         |               |                | 7.71     | 20              |
| Chrysene              | 0.00200              | 0.00205            | 0.00213             | 102           | 106            | 64.0-144         |               |                | 3.83     | 20              |
| Dibenz(a,h)anthracene | 0.00200              | 0.00205            | 0.00231             | 102           | 115            | 52.0-155         |               |                | 11.9     | 20              |
| Fluoranthene          | 0.00200              | 0.00216            | 0.00211             | 108           | 105            | 69.0-153         |               |                | 2.34     | 20              |



Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3357755-1 11/07/18 12:39 • (LCSD) R3357755-2 11/07/18 13:01

| Analyte                     | Spike Amount<br>mg/l | LCS Result<br>mg/l | LCSD Result<br>mg/l | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | <u>LCS Qualifier</u> | <u>LCSD Qualifier</u> | RPD<br>% | RPD Limits<br>% |
|-----------------------------|----------------------|--------------------|---------------------|---------------|----------------|------------------|----------------------|-----------------------|----------|-----------------|
| Fluorene                    | 0.00200              | 0.00193            | 0.00196             | 96.5          | 98.0           | 64.0-136         |                      |                       | 1.54     | 20              |
| Indeno(1,2,3-cd)pyrene      | 0.00200              | 0.00211            | 0.00238             | 105           | 119            | 54.0-153         |                      |                       | 12.0     | 20              |
| Naphthalene                 | 0.00200              | 0.00206            | 0.00209             | 103           | 105            | 61.0-137         |                      |                       | 1.45     | 20              |
| Phenanthrene                | 0.00200              | 0.00189            | 0.00199             | 94.5          | 99.5           | 62.0-137         |                      |                       | 5.15     | 20              |
| Pyrene                      | 0.00200              | 0.00189            | 0.00210             | 94.5          | 105            | 60.0-142         |                      |                       | 10.5     | 20              |
| 1-Methylnaphthalene         | 0.00200              | 0.00232            | 0.00228             | 116           | 114            | 66.0-142         |                      |                       | 1.74     | 20              |
| 2-Methylnaphthalene         | 0.00200              | 0.00210            | 0.00207             | 105           | 103            | 62.0-136         |                      |                       | 1.44     | 20              |
| 2-Chloronaphthalene         | 0.00200              | 0.00197            | 0.00204             | 98.5          | 102            | 64.0-140         |                      |                       | 3.49     | 20              |
| <i>(S) Nitrobenzene-d5</i>  |                      |                    |                     | 150           | 144            | 31.0-160         |                      |                       |          |                 |
| <i>(S) 2-Fluorobiphenyl</i> |                      |                    |                     | 105           | 103            | 48.0-148         |                      |                       |          |                 |
| <i>(S) p-Terphenyl-d14</i>  |                      |                    |                     | 111           | 106            | 37.0-146         |                      |                       |          |                 |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc



Guide to Reading and Understanding Your Laboratory Report

The information below is designed to better explain the various terms used in your report of analytical results from the Laboratory. This is not intended as a comprehensive explanation, and if you have additional questions please contact your project representative.

Abbreviations and Definitions

|                              |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
|------------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| MDL                          | Method Detection Limit.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    |
| ND                           | Not detected at the Reporting Limit (or MDL where applicable).                                                                                                                                                                                                                                                                                                                                                                                                                                                             |
| RDL                          | Reported Detection Limit.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                  |
| Rec.                         | Recovery.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                  |
| RPD                          | Relative Percent Difference.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                               |
| SDG                          | Sample Delivery Group.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                     |
| (S)                          | Surrogate (Surrogate Standard) - Analytes added to every blank, sample, Laboratory Control Sample/Duplicate and Matrix Spike/Duplicate; used to evaluate analytical efficiency by measuring recovery. Surrogates are not expected to be detected in all environmental media.                                                                                                                                                                                                                                               |
| U                            | Not detected at the Reporting Limit (or MDL where applicable).                                                                                                                                                                                                                                                                                                                                                                                                                                                             |
| Analyte                      | The name of the particular compound or analysis performed. Some Analyses and Methods will have multiple analytes reported.                                                                                                                                                                                                                                                                                                                                                                                                 |
| Dilution                     | If the sample matrix contains an interfering material, the sample preparation volume or weight values differ from the standard, or if concentrations of analytes in the sample are higher than the highest limit of concentration that the laboratory can accurately report, the sample may be diluted for analysis. If a value different than 1 is used in this field, the result reported has already been corrected for this factor.                                                                                    |
| Limits                       | These are the target % recovery ranges or % difference value that the laboratory has historically determined as normal for the method and analyte being reported. Successful QC Sample analysis will target all analytes recovered or duplicated within these ranges.                                                                                                                                                                                                                                                      |
| Original Sample              | The non-spiked sample in the prep batch used to determine the Relative Percent Difference (RPD) from a quality control sample. The Original Sample may not be included within the reported SDG.                                                                                                                                                                                                                                                                                                                            |
| Qualifier                    | This column provides a letter and/or number designation that corresponds to additional information concerning the result reported. If a Qualifier is present, a definition per Qualifier is provided within the Glossary and Definitions page and potentially a discussion of possible implications of the Qualifier in the Case Narrative if applicable.                                                                                                                                                                  |
| Result                       | The actual analytical final result (corrected for any sample specific characteristics) reported for your sample. If there was no measurable result returned for a specific analyte, the result in this column may state "ND" (Not Detected) or "BDL" (Below Detectable Levels). The information in the results column should always be accompanied by either an MDL (Method Detection Limit) or RDL (Reporting Detection Limit) that defines the lowest value that the laboratory could detect or report for this analyte. |
| Case Narrative (Cn)          | A brief discussion about the included sample results, including a discussion of any non-conformances to protocol observed either at sample receipt by the laboratory from the field or during the analytical process. If present, there will be a section in the Case Narrative to discuss the meaning of any data qualifiers used in the report.                                                                                                                                                                          |
| Quality Control Summary (Qc) | This section of the report includes the results of the laboratory quality control analyses required by procedure or analytical methods to assist in evaluating the validity of the results reported for your samples. These analyses are not being performed on your samples typically, but on laboratory generated material.                                                                                                                                                                                              |
| Sample Chain of Custody (Sc) | This is the document created in the field when your samples were initially collected. This is used to verify the time and date of collection, the person collecting the samples, and the analyses that the laboratory is requested to perform. This chain of custody also documents all persons (excluding commercial shippers) that have had control or possession of the samples from the time of collection until delivery to the laboratory for analysis.                                                              |
| Sample Results (Sr)          | This section of your report will provide the results of all testing performed on your samples. These results are provided by sample ID and are separated by the analyses performed on each sample. The header line of each analysis section for each sample will provide the name and method number for the analysis reported.                                                                                                                                                                                             |
| Sample Summary (Ss)          | This section of the Analytical Report defines the specific analyses performed for each sample ID, including the dates and times of preparation and/or analysis.                                                                                                                                                                                                                                                                                                                                                            |

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

| Qualifier | Description                                                                                           |
|-----------|-------------------------------------------------------------------------------------------------------|
| J         | The identification of the analyte is acceptable; the reported value is an estimate.                   |
| J6        | The sample matrix interfered with the ability to make any accurate determination; spike value is low. |
| T8        | Sample(s) received past/too close to holding time expiration.                                         |



Pace National is the only environmental laboratory accredited/certified to support your work nationwide from one location. One phone call, one point of contact, one laboratory. No other lab is as accessible or prepared to handle your needs throughout the country. Our capacity and capability from our single location laboratory is comparable to the collective totals of the network laboratories in our industry. The most significant benefit to our one location design is the design of our laboratory campus. The model is conducive to accelerated productivity, decreasing turn-around time, and preventing cross contamination, thus protecting sample integrity. Our focus on premium quality and prompt service allows us to be YOUR LAB OF CHOICE.

\* Not all certifications held by the laboratory are applicable to the results reported in the attached report.  
 \* Accreditation is only applicable to the test methods specified on each scope of accreditation held by Pace National.

## State Accreditations

|                         |             |                             |                   |
|-------------------------|-------------|-----------------------------|-------------------|
| Alabama                 | 40660       | Nebraska                    | NE-OS-15-05       |
| Alaska                  | 17-026      | Nevada                      | TN-03-2002-34     |
| Arizona                 | AZ0612      | New Hampshire               | 2975              |
| Arkansas                | 88-0469     | New Jersey-NELAP            | TN002             |
| California              | 2932        | New Mexico <sup>1</sup>     | n/a               |
| Colorado                | TN00003     | New York                    | 11742             |
| Connecticut             | PH-0197     | North Carolina              | Env375            |
| Florida                 | E87487      | North Carolina <sup>1</sup> | DW21704           |
| Georgia                 | NELAP       | North Carolina <sup>3</sup> | 41                |
| Georgia <sup>1</sup>    | 923         | North Dakota                | R-140             |
| Idaho                   | TN00003     | Ohio-VAP                    | CL0069            |
| Illinois                | 200008      | Oklahoma                    | 9915              |
| Indiana                 | C-TN-01     | Oregon                      | TN200002          |
| Iowa                    | 364         | Pennsylvania                | 68-02979          |
| Kansas                  | E-10277     | Rhode Island                | LA000356          |
| Kentucky <sup>1,6</sup> | 90010       | South Carolina              | 84004             |
| Kentucky <sup>2</sup>   | 16          | South Dakota                | n/a               |
| Louisiana               | AI30792     | Tennessee <sup>1,4</sup>    | 2006              |
| Louisiana <sup>1</sup>  | LA180010    | Texas                       | T 104704245-17-14 |
| Maine                   | TN0002      | Texas <sup>5</sup>          | LAB0152           |
| Maryland                | 324         | Utah                        | TN00003           |
| Massachusetts           | M-TN003     | Vermont                     | VT2006            |
| Michigan                | 9958        | Virginia                    | 460132            |
| Minnesota               | 047-999-395 | Washington                  | C847              |
| Mississippi             | TN00003     | West Virginia               | 233               |
| Missouri                | 340         | Wisconsin                   | 9980939910        |
| Montana                 | CERT0086    | Wyoming                     | A2LA              |

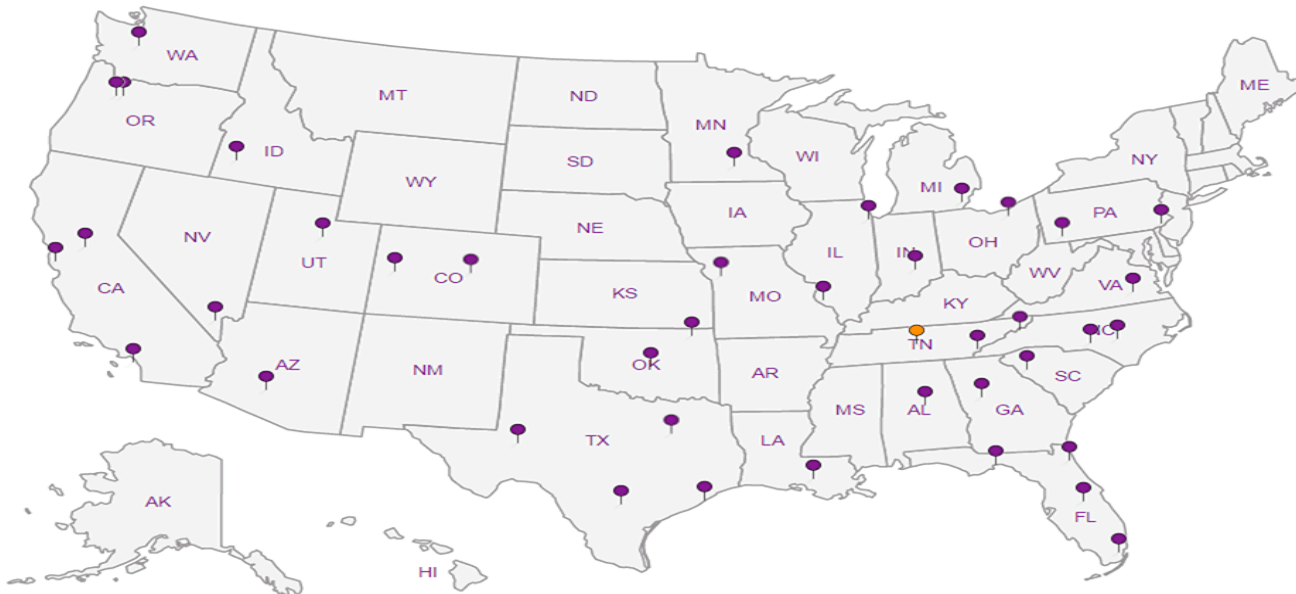
## Third Party Federal Accreditations

|                               |         |                    |               |
|-------------------------------|---------|--------------------|---------------|
| A2LA – ISO 17025              | 1461.01 | AIHA-LAP,LLC EMLAP | 100789        |
| A2LA – ISO 17025 <sup>5</sup> | 1461.02 | DOD                | 1461.01       |
| Canada                        | 1461.01 | USDA               | P330-15-00234 |
| EPA-Crypto                    | TN00003 |                    |               |

<sup>1</sup> Drinking Water <sup>2</sup> Underground Storage Tanks <sup>3</sup> Aquatic Toxicity <sup>4</sup> Chemical/Microbiological <sup>5</sup> Mold <sup>6</sup> Wastewater n/a Accreditation not applicable

## Our Locations

Pace National has sixty-four client support centers that provide sample pickup and/or the delivery of sampling supplies. If you would like assistance from one of our support offices, please contact our main office. Pace National performs all testing at our central laboratory.



1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

**Terracon - Longmont**  
**12831 Lefthand Circle, Suite C**  
**Longmont, CO 80501**

Billing information:  
**Same as Address**

Analysis / Container / Preservative

Chain of Custody Page 1 of 1

Report to:  
**Mike Skridulis**

Email To:  
**mike.skridulis@terracon.com**

Project Description:  
**Union Reservoir**

City/State Collected:  
**Longmont, CO**

Phone: **303-454-5249**  
 Fax: **970-484-0454**

Client Project #  
**22187053**

Lab Project #

Collected by (print):  
**Charles Covington**

Site/Facility ID #

P.O. #

Collected by (signature):  
*Charles Covington*

**Rush?** (Lab MUST Be Notified)  
 Same Day  Five Day  
 Next Day  5 Day (Rad Only)  
 Two Day  10 Day (Rad Only)  
 Three Day

Quote #  
 Date Results Needed

Immediately Packed on ice N    Y X

| Sample ID | Comp/Grab | Matrix * | Depth | Date     | Time | No. of Cntrs | VOC8260 (2) 40ml Amber wHCl | Diss. Methane, Ethane, Ethylene (2) 40ml | CO2 - 125ml HDPE No Pres | Diss. Metals - 250ml HDPE No Pres | PAHSIM (2) 40ml Amber No Pres | Cl, S04, TDS - 250ml HDPE No Pres |
|-----------|-----------|----------|-------|----------|------|--------------|-----------------------------|------------------------------------------|--------------------------|-----------------------------------|-------------------------------|-----------------------------------|
| MW-01     | Grab      | GW       |       | 10/31/18 | 1155 | 9            | X                           | X                                        | X                        | X                                 | X                             | X                                 |
| MW-02     | Grab      | GW       |       | 10/31/18 | 1235 | 9            | X                           | X                                        | X                        | X                                 | X                             | X                                 |
| MW-03     | Grab      | GW       |       | 10/31/18 | 1325 | 9            | X                           | X                                        | X                        | X                                 | X                             | X                                 |

|                             |                                          |                          |                                   |                               |                                   |
|-----------------------------|------------------------------------------|--------------------------|-----------------------------------|-------------------------------|-----------------------------------|
| VOC8260 (2) 40ml Amber wHCl | Diss. Methane, Ethane, Ethylene (2) 40ml | CO2 - 125ml HDPE No Pres | Diss. Metals - 250ml HDPE No Pres | PAHSIM (2) 40ml Amber No Pres | Cl, S04, TDS - 250ml HDPE No Pres |
|-----------------------------|------------------------------------------|--------------------------|-----------------------------------|-------------------------------|-----------------------------------|

12065 Lebanon Rd  
 Mount Juliet, TN 37122  
 Phone: 615-758-5858  
 Phone: 800-767-5859  
 Fax: 615-758-5859

L# **L1040168**  
**F005**

Acctnum: **TERRALCO**  
 Template:  
 Prelogin:  
 TSR: **Daphne Richards**  
 PB:

Shipped Via:

\* Matrix:  
 SS - Soil AIR - Air F - Filter  
 GW - Groundwater B - Bioassay  
 WW - WasteWater  
 DW - Drinking Water  
 OT - Other

Remarks:  
**PAD SCREEN: <0.5 mR/hr**  
 Tracking # **4510 1654 8135**

**Sample Receipt Checklist**  
 COC Seal Present/Intact:    NP    Y    N  
 CDC Signed/Accurate:    Y    N  
 Bottles arrive intact:    Y    N  
 Correct bottles used:    Y    N  
 Sufficient volume sent:    Y    N  
 If Applicable  
 VOA Zero Headpace:    Y    N  
 Preservation Correct/Checked:    Y    N

Relinquished by: (Signature)  
 Date: Time:

Received by: (Signature)  
 Date: Time:

Received by: (Signature)  
 Date: Time:

Trip Blank Received: Yes (No)  
 HCL/MeOH  
 TBR  
 Temp: **20.1 °C**  
**2019/11/1**  
**27**

If preservation required by Login: Date/Time  
 Hold:  
 Condition: **NCF / OK**

**Andy Vann**



|                  |                  |               |                        |
|------------------|------------------|---------------|------------------------|
| Login #:L1040168 | Client: TERRALCO | Date:11/01/18 | Evaluated by:Andy Vann |
|------------------|------------------|---------------|------------------------|

**Non-Conformance (check applicable items)**

| Sample Integrity               | Chain of Custody Clarification                   | If Broken Container:                                 |
|--------------------------------|--------------------------------------------------|------------------------------------------------------|
| Parameter(s) past holding time | Login Clarification Needed                       | Insufficient packing material around container       |
| Improper temperature           | Chain of custody is incomplete                   | Insufficient packing material inside cooler          |
| Improper container type        | Please specify Metals requested.                 | Improper handling by carrier (FedEx / UPS / Courier) |
| Improper preservation          | Please specify TCLP requested.                   | Sample was frozen                                    |
| Insufficient sample volume.    | Received additional samples not listed on coc.   | Container lid not intact                             |
| Sample is biphasic.            | Sample ids on containers do not match ids on coc | <b>If no Chain of Custody:</b>                       |
| Vials received with headspace. | Trip Blank not received.                         | Received by:                                         |
| Broken container               | Client did not "X" analysis.                     | Date/Time:                                           |
| Broken container:              | Chain of Custody is missing                      | Temp./Cont. Rec./pH:                                 |
| Sufficient sample remains      |                                                  | Carrier:                                             |
|                                |                                                  | Tracking#                                            |

**Login Comments: What metals?**

|                     |                 |       |            |       |       |
|---------------------|-----------------|-------|------------|-------|-------|
| Client informed by: | Call            | Email | Voice Mail | Date: | Time: |
| TSR Initials:       | Client Contact: |       |            |       |       |

**Login Instructions:**

Please log dissolved metals for AGDICP, ASDICP, BADICP, BDICP, CDDICP, CRDICP, CUDICP, PBDICP, HGD, NIDICP, SEDICP, and ZNDICP.  
CRICP(total) and CR3 and CR6