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Subject:  
2022 First Semi-Annual Groundwater Monitoring Report

ENVIRONMENT

Dear Ms. Reams,

On behalf of Chevron Environmental Management Company (Chevron), Arcadis US, Inc. (Arcadis) has prepared the attached *2022 First Semi-Annual Groundwater Monitoring Report* for the first semi-annual groundwater sampling event of 2022 for the following facility:

Date:  
June 2, 2022

Contact:  
Gerald Robinson

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724.934.9507

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Our ref:  
30063667

<u>Chevron Facility No.</u>	<u>ADEC File No.</u>	<u>Hazard ID</u>	<u>Location</u>
97324	2100.26.008	23885	4417 Lake Otis Parkway Anchorage, Alaska

If you have any questions, please do not hesitate to contact me.

Sincerely,

Arcadis U.S., Inc.

Gerald A. Robinson  
Project Manager

Copies:  
Susan Erickson, Chevron (*electronic copy*)  
Mark Engelke, Cook Inlet Marketing Group, Inc. (*electronic copy*)  
Emma Giboney, Municipality of Anchorage (*electronic copy*)

Chevron Environmental Management Company

# **2022 FIRST SEMI-ANNUAL GROUNDWATER MONITORING REPORT**

Former Chevron-Branded  
Service Station No. 97324  
4417 Lake Otis Parkway  
Anchorage, Alaska  
ADEC File No. 2100.26.008

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June 02, 2022

## 2022 FIRST SEMI-ANNUAL GROUNDWATER MONITORING REPORT

### Former Chevron-Branded Service Station No. 97324



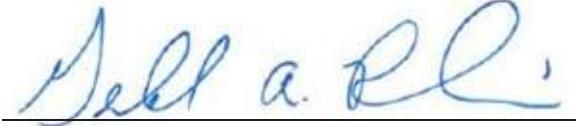
Sydney Clark, E.I.T.  
Environmental Engineer

4417 Lake Otis Parkway  
Anchorage, Alaska

ADEC File No: 2100.26.008  
HAZARD ID No: 23885

Prepared for:

Chevron Environmental Management Company



Gerald A. Robinson.  
Project Manager

Prepared by:  
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Our Ref.:

30063667 Date:  
June 2, 2022

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**SEMI-ANNUAL GROUNDWATER MONITORING REPORT**
**FIRST HALF 2022**
**June 2, 2022**

Facility No: Former Chevron-Branded  
Station No. 97324

Address: 4417 Lake Otis Parkway  
Anchorage, Alaska

Arcadis Contact Person / Phone No.:

Gerald Robinson / (724) 934-9507

Arcadis Project No.:

30063667

Primary Agency/Regulatory ID No.:

Alaska Department of Conservation (ADEC) / Rebekah Reams /ADEC File ID: 2100.26.008

**WORK CONDUCTED THIS PERIOD [First Half 2022]:**

1. Conducted semi-annual groundwater monitoring activities on April 4, 2022.
2. Prepared the *2022 First Semi-Annual Groundwater Monitoring Report*.

**WORK PROPOSED NEXT PERIOD [Second Half 2022]:**

1. Conduct semi-annual groundwater monitoring activities in the Second half of 2022
2. Conduct the soil assessment activities outlined in the *System Removal, Well Decommissioning, and Soil Assessment Work Plan* submitted to ADEC on July 14, 2021.
3. Prepare the *2022 Second Semi-Annual Groundwater Monitoring Report*.

Current Phase of Project:	Monitoring	
Frequency of Monitoring / Sampling:	Semi-annual	
Are Light Non-Aqueous Phase Liquid (LNAPL) Present On-site:	No	
Cumulative LNAPL Recovered to Date:	0.00	(gallons)
Approximate Depth to Groundwater:	15.50 to 24.58	(feet below top of casing)
Approximate Groundwater Elevation:	143.67 to 143.74	(feet relative to NAVD88)
Groundwater Flow Direction	North-Northwest (historically)	

Groundwater Gradient	Not Calculated	(feet per foot)
Current Remediation Techniques:	None	
Permits for Discharge:	None	
Summary of Unusual Activity:	None	
Agency Directive Requirements:	None	

## 1 INTRODUCTION

On behalf of Chevron Environmental Management Company (CEMC), Arcadis US, Inc. (Arcadis), has prepared this report to document the first semi-annual groundwater sampling event of 2022 for Chevron facility 97324, located at 4417 Lake Otis Parkway in Anchorage, Alaska (site). The site location map and site plan are presented on Figure 1 and Figure 2, respectively.

This work was conducted under the direction of a “Qualified Environmental Professional” (QEP) and “Qualified Sampler” (18 Alaska Administrative Code [AAC] 75.333). Site background and history summaries are attached as Appendix A.

## 2 GROUNDWATER MONITORING

### 2.1 Groundwater Gauging Methods

The 2022 First semi-annual groundwater gauging event was conducted on April 4, 2022. Monitoring wells MW-1R, MW-2R, MW-8RR, and MW-9 were gauged with an oil/water interface probe to determine depth-to-water and to ascertain if light non-aqueous phase liquid (LNAPL) was present.

In order to prevent the possibility of cross-contamination, wells were gauged in the order of lowest to highest historical petroleum hydrocarbon concentrations in groundwater. In addition, non-disposable groundwater gauging equipment was decontaminated prior to and after each use with a detergent solution and rinsed in potable water.

### 2.2 Groundwater Elevation and Flow Direction

During the 2022 First semi-annual event, monitoring wells MW-1R, MW-2R, MW-8RR, and MW-9 were gauged for groundwater elevations and the presence of LNAPL. The groundwater monitoring event field notes are presented in Appendix B.

The inferred groundwater flow direction for the second semi-annual 2022 monitoring event not calculated, however historic groundwater flow at the site has been observed to the north-northwest. Current and historical groundwater gauging and analytical results are summarized in Table 1 and Table 4, respectively. A groundwater elevation map with a rose diagram of historical flow directions is presented as Figure 3.

## 2.3 Groundwater Sampling Methods

The second semi-annual groundwater monitoring event was conducted on April 4, 2022. Groundwater samples were collected from monitoring wells MW-1R, MW-2R, MW-8RR, and MW-9 using a low flow purge sampling method.

Sampling procedures were conducted in accordance with ADEC *Field Sampling Guidance* (ADEC, 2019). Monitoring well caps were removed to allow groundwater levels to stabilize and equilibrate before using an oil/water interface probe meter capable of 0.01-foot accuracy to measure the depth to groundwater and total well depth. A bladder pump with compressor & control unit with clean/disposable Teflon lined tubing and bladders was used to purge groundwater from the wells and collect samples to minimize the risk of volatile contaminant absorption by the sampling equipment. Water table drawdown was continuously monitored during purging with a water level meter and the flow rate of the pump was adjusted to limit drawdown to 0.1 meter. The intake of the pump was set as close as possible to the soil groundwater interface. Water quality parameters were monitored during purging with a multi-parameter water quality meter equipped with a flow through cell and Turbidity meter. Parameters were recorded every 3 to 5 minutes until a minimum of three (minimum of four if using temperature as an indicator) of the parameters listed below stabilized. The flow rate was reduced to 100-150 ml/minute and samples were collected from the discharge line into laboratory sample bottles. Water quality parameters were considered stable when three successive readings were within the following ADEC limits:

- $\pm 3\%$  for temperature (minimum of  $\pm 0.2\text{ C}^\circ$ ),
- $\pm 0.1$  for pH,
- $\pm 3\%$  for conductivity,
- $\pm 10\text{ mv}$  for redox potential,
- $\pm 10\%$  for dissolved oxygen, and
- $\pm 10\%$  for turbidity.

Sample bottles were labeled, stored in a cooler packed with ice, and submitted to Pace Analytical (National Center for Testing & Innovation) Laboratory in Mount Juliet, Tennessee, under proper chain-of-custody procedures.

Groundwater samples collected from monitoring wells MW-1R, MW-2R, MW-8RR, and MW-9 were submitted to the analytical laboratory for the following analyses:

- Full-scan volatile organic compounds (VOCs) including benzene, toluene, ethylbenzene, total xylenes (collectively BTEX), methyl-t-butyl ether (MTBE), and naphthalene by United States Environmental Protection Agency (USEPA) Method 8260D
- Total petroleum hydrocarbons as gasoline range organics (TPH-g) by Alaska Method AK101
- Total petroleum hydrocarbons as diesel range organics (TPH-d) by Alaska Method AK102

Additionally, groundwater samples were collected from MW-2R are analyzed for polycyclic aromatic hydrocarbons (PAHs) by USEPA Method 8270E-SIM.

A groundwater duplicate sample was collected from monitoring well MW-2R. The duplicate sample was analyzed for full-scan VOCs, TPH-g, TPH-d, and PAHs. The duplicate sample was submitted blind with the sample set to Pace Analytical.

## 2.4 Groundwater Analytical Results

Routine analytical results for BTEX, MTBE, naphthalene, TPH-g, and TPH-d from the first semi-annual 2022 groundwater monitoring event are summarized in Table 1 and additional VOCs analyzed by USEPA Method 8260D are summarized in Table 2. Current and historic analytical data for PAHs are summarized in Table 3. Historical groundwater gauging and analytical data are summarized in Table 4. Historical additional VOCs analyzed by USEPA Method 8260D are summarized in Tables 5a, 5b, 5c and 5d.

Current analytical results for BTEX, MTBE, naphthalene, TPH-g, and TPH-d are summarized in Figure 4. Current analytical results for the solvents 1,2-dichloroethane, trichloroethene, tetrachloroethylene, cis-1,2-dichloroethene, and methylene chloride are summarized in Figure 5.

## 3 INVESTIGATION DERIVED WASTE

Purge water and decontamination water from groundwater was temporarily collected into 5-gallon buckets and treated onsite via a Granular Activated Carbon (GAC) bucket. The treatment of purge water and decontamination water was completed per the Technical Guidance Instructions for Investigation Derived waste Treatment Utilizing Granular Activated Carbon provided in Appendix C.

## 4 LABORATORY DATA QUALITY ASSURANCE SUMMARY

As required by ADEC (Technical Memorandum, October 2019), Arcadis completed a laboratory data review checklist for each of the laboratory report generated for the 2020 semi-annual event. The laboratory report is included as Appendix D and data review checklist is included as Appendix E. The following quality assurance (QA) summary describes six parameters, related to the quality and usability of the data presented in this report.

### 4.1 Precision

The RPD between field duplicated (FD) were within the control limits.

The RPD between the matrix spike and matrix spike duplicate (MS/MSD) exceeded control limits for compound TPH-g in sample location MW-1R for method AK101. The compound result was qualified as estimated.

The RPD between laboratory control sample/laboratory control sample duplicate (LCS/LCSD) exceeded the control limits for several compounds in sample locations MW-2R, MW-8RR, MW-1R, MW-9, the blind duplicate (BD-1), the equipment blank (EQB-1) and the trip blank for method USEPA 8260D. The compounds results were qualified as estimated.

The precision of the data, as measured by laboratory quality control (QC) indicators, suggest that the Data Quality Objectives (DQOs) were met.

### 4.2 Accuracy

The percent recovery for surrogates were within the control limits.

MS/MSD recovery exceeded control limits for compounds TPH-g for method AK101 in sample location MW-1R. The compound in the associated sample location was qualified as estimated.

Continuing calibration for compound acrolein exhibited a low bias recovery. Associated result in sample locations MW-2R, MW-8RR, MW-1R, MW-9, BD-1, EQB-1, and the trip blank were qualified as estimated.

The accuracy of the data, as measured by laboratory quality control (QC) indicators, suggest that the DQOs were met.

#### **4.3 Representativeness**

The data appear to be representative of site conditions and are generally consistent with historical groundwater monitoring results and expected impacts to groundwater.

#### **4.4 Comparability**

The laboratory results are presented in the same units as previous reports to allow comparison. The target compounds were not detected in trip blank, equipment blank and method blank.

#### **4.5 Completeness**

The results appear to be valid and usable, and thus, the laboratory results have 100% completeness.

#### **4.6 Sensitivity**

The concentrations of TPH-d, benzene, ethylbenzene and naphthalene exceeded the ADEC groundwater cleanup levels (GCLs) in the sample from MW-2R.

The concentration of 1,2-dichloroethane exceeded ADEC GCLs in samples from MW-1R and MW-2R.

The concentrations of trichloroethylene exceeded ADEC GCLs in the sample from MW-9.

The laboratory reported detection limit for compounds trichloroethene, 1,1,2,2-tetrachloroethane, 1,1,2-trichloroethane, 1,2-dibromoethane, chloroform, trans-1,3-dichloropropene and vinyl chloride exceeded the ADEC groundwater cleanup level; however, the laboratory method detection limit is below the ADEC groundwater cleanup level. As most samples were reported as not detected for the mentioned constituents, except where mentioned above.

The sensitivity of the analyses was adequate for the samples as the detection limits were less than the ADEC GCLs for compounds with above exceptions.

### **5 CONCLUSIONS AND RECOMMENDATIONS**

The groundwater data collected during the first semi-annual 2022 event indicate groundwater the groundwater table is relatively flat with no predominate flow direction. During the first semi-annual 2022 groundwater monitoring event, groundwater samples were collected for analysis from monitoring wells MW-1R, MW-2R, MW-8RR, and MW-9. Analytical results from the monitoring wells are generally consistent with historical data.

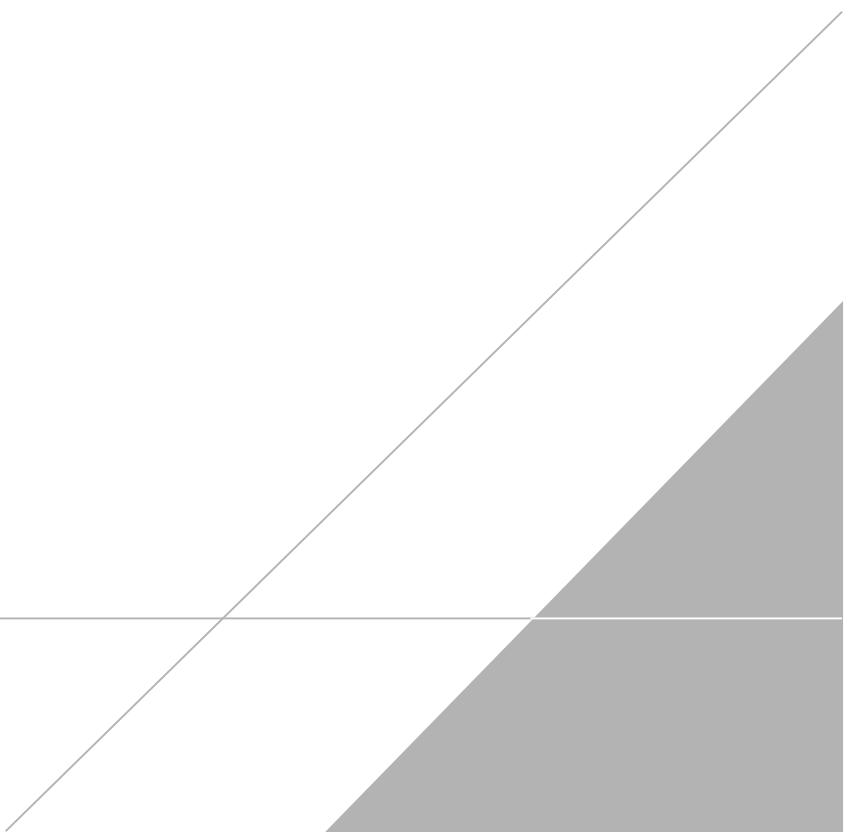
Groundwater monitoring will continue in accordance with the current semi-annual schedule. The next groundwater sampling event will be conducted in the fall of 2022.

## 6 REFERENCES

ADEC. *Field Sampling Guidance*. Division of Spill Prevention and Response Contaminated Sites Program. 2019.

ADEC Technical Memorandum, October 2019. *Minimum Quality Assurance Requirements for Sample Handling, Reports and Laboratory Data*. ADEC, Division of Spill Prevention and Response Contaminated Sites Program.

# TABLES



**Table 1. Current Groundwater Gauging and Analytical Results 1SA22**

Former Chevron-Branded Service Station 97324  
 4417 Lake Otis Parkway  
 Anchorage, Alaska

Well ID	Sample Date	TOC (ft)	Datum	DTW (ft bTOC)	Thickness (ft)	GW Elev (ft)	LNAPL							
							TPH-d (mg/L)	TPH-g (mg/L)	Benzene (mg/L)	Toluene (mg/L)	Ethylbenzene (mg/L)	Total Xylenes (mg/L)		
							1.5	2.2	0.0046	1.1	0.015	0.19		
												0.14		
												0.0017		
<b>ADEC Groundwater Cleanup Levels</b>														
MW-1R	4/4/2022	167.56	NAVD88	23.85	0.00	143.71	<0.8	<0.1 J	<0.00100	<0.00100	<0.00100	<0.00300	<0.00100 J	<0.00500 J
MW-2R	4/4/2022	168.25	NAVD88	24.58	0.00	143.67	<b>1.61 [1.61]</b>	<b>1.22 [1.36]</b>	<b>0.00629 [0.00654]</b>	<b>0.00132 [0.00161]</b>	<b>0.0723 [0.087]</b>	<b>0.0227 [0.0294]</b>	<b>&lt;0.00100 J [&lt;0.00100 J]</b>	<b>0.0310 J [ 0.0315 J]</b>
MW-8RR	4/4/2022	166.43	NAVD88	22.73	0.00	143.70	<0.8	<0.1	<0.00100	<0.00100	<0.00100	<0.00300	<0.00100 J	<0.00500 J
MW-9	4/4/2022	159.24	NAVD88	15.50	0.00	143.74	<0.8	<b>0.0445 J</b>	<0.00100	<0.00100	<0.00100	<0.00300	<0.00100 J	<0.00500 J
Trip Blank	4/4/2022	--	--	--	--	--	--	<0.1	<0.00100	<0.00100	<0.00100	<0.00300	<0.00100 J	<0.00500 J
Equipment Blank	4/4/2022	--	--	--	--	--	<0.8	<0.1	<0.00100	<0.00100	<0.00100	<0.00300	<0.00100 J	<0.00500 J

**Notes:**

ID = Identification

MW = Groundwater monitoring well

TOC = Top of casing

DTW = Depth to groundwater

ft bTOC = Feet below top of casing

ft = Feet relative to NAVD88

mg/L = Milligrams per liter

GW Elev = Groundwater elevation

&lt;0.00100 = Not detected at or above the reported detection limit (RDL)

**Bold and Shaded** = Value exceeds ADEC Groundwater Cleanup Level**Bold** = Detected above laboratory method detection limit (MDL)**Bold and *Italicized*** : Constituent considered non-detect, however

Laboratory RDL is greater than the ADEC Groundwater Cleanup Level

J = The compound was positively identified; however, the associated numerical value is an estimated concentration only.

TPH-g = Total petroleum hydrocarbons, gasoline range by LUFT GC/MS according to United States Environmental Protection Agency (USEPA) Method AK101

TPH-d = Total petroleum hydrocarbons, diesel range by LUFT GC/MS according to State of Alaska Method AK102.

Samples analytes by USEPA Method 8260D:

Benzene, Toluene, Ethylbenzene and Total Xylenes (collectively BTEX)

MTBE = Methyl-tert-butyl ether

Naphthalene

LUFT = Leaking Underground Fuel Tank

GC/MS = Gas chromatography/Mass Spectrometry

ADEC = Alaska Department of Environmental Conservation

NAVD88 = North American Vertical Datum of 1988

LNAPL = Light Non-Aqueous Phase Liquid

-- = Not Measured/Not analyzed

[ ] = Blind Duplicate Sample Result



**Table 3. Current and Historical Groundwater Analytical Results - PAHs**  
 Former Chevron-Branded Service Station 97324  
 4417 Lake Otis Parkway  
 Anchorage, Alaska

Well ID	Sample Date	1-Methylnaphthalene µg/L	2-Methylnaphthalene µg/L	Acenaphthene µg/L	Acenaphthylene µg/L	Anthracene µg/L	Benz(a)anthracene µg/L	Benz(a)pyrene µg/L	Benz(b)fluoranthene µg/L	Benz(g,h,i)perylene µg/L	Benz(k)fluoranthene µg/L	Chrysene µg/L	Dibenz(a,h)anthracene µg/L	Fluoranthene µg/L	Fluorene µg/L	Indeno(1,2,3-cd)pyrene µg/L	Naphthalene µg/L	Phenanthrene µg/L	Pyrene µg/L	
<b>ADEC Groundwater Cleanup Levels</b>		<b>11</b>	<b>36</b>	<b>530</b>	<b>260</b>	<b>43</b>	<b>0.3</b>	<b>0.25</b>	<b>2.5</b>	<b>0.26</b>	<b>0.8</b>	<b>2</b>	<b>0.25</b>	<b>260</b>	<b>290</b>	<b>0.19</b>	<b>1.7</b>	<b>170</b>	<b>120</b>	
MW-2R	9/11/2019	0.17	<b>0.058 J</b>	<0.11	<0.0503	<0.11	<0.053	<0.11	<0.053	<0.053	<0.11	<0.21	<0.11	<0.053	<b>1.8</b>	<0.11	<0.11	<0.11		
MW-2R	4/22/2020	0.360 J	<0.0510	<0.0510	<0.0510	<0.0510	<0.0510	<0.0510	<0.0510	<0.0510	<0.0510	<0.255	<0.0510	<0.0510	<0.0510	<0.0510	<0.0510	<0.0510		
MW-2R	10/9/2020	12.0 [11.4]	<b>0.922 [0.893]</b>	<b>0.0792 [0.0753]</b>	<0.0500 [<0.0500]	<0.0500 [<0.0500]	<0.0500 [<0.0500]	<0.0500 <b>[0.0260 J]</b>	<0.0500 <b>[0.0245 J]</b>	<0.0500 <b>[0.0413 J]</b>	<0.0500 <b>[0.0245 J]</b>	<0.250 [<0.250]	<0.0500 <b>[0.0305 J]</b>	<0.0500 [<0.0500]	<0.0500 <b>[0.0190 J]</b>	<0.0500 <b>[0.0184 J]</b>	<b>0.0273 [0.0261]</b>	<0.0500 <b>[0.0839]</b>	<0.0500 <b>[0.0668]</b>	
MW-2R	8/26/2021	<b>0.0117 [0.0111]</b>	<b>0.0074 [0.00679]</b>	<b>0.0000726 [0.0000692]</b>	<0.0000515 [<0.0000510]	<0.0000515 [<0.0000510]	<0.0000515 B [<0.0000510]	<b>0.0000381 J</b> [<0.0000510]	<b>0.0000402 J</b> [<0.0000510]	<b>0.0000425 J</b> [<0.0000510]	<b>0.0000347 J</b> [<0.0000255]	<b>0.0000315 J</b> [<0.0000510]	<b>0.0000384 J</b> [<0.0000510]	<0.0000515 B [<0.0000510]	<b>0.0000228 J</b> [<0.0000510]	<b>0.0000230 J</b> [<0.0000510]	<b>0.036 [0.0349]</b>	<b>0.0000243 J</b> [<0.0000510]	<0.0000515 B [<0.0000510]	<0.0000515 B [<0.0000510]
MW-2R	04/04/2022	<b>0.00811 [0.00959]</b>	--	<b>0.0000560 J</b> [ <b>0.0000643</b> ]	<0.0000625 [<0.0000625]	<0.0000625 [<0.0000625]	<0.0000625 [<0.0000625]	<0.0000625 [<0.0000625]	<0.0000625 [<0.0000625]	<0.0000625 [<0.0000625]	<0.0000625 [<0.0000625]	<0.0000625 [<0.0000625]	<0.0000625 [<0.0000625]	<0.0000625 [<0.0000625]	<b>0.0000223 J</b> [<0.0000625]	<0.0000625 [<0.0000625]	<b>0.0191 [0.0222]</b>	<0.0000625 [<0.0000625]	<0.0000625 [<0.0000625]	<0.0000625 [<0.0000625]
Equipment Blank	10/9/2020	<b>0.0208 J</b>	<0.500	<0.0500	<0.0500	<0.0500	<0.0500	<0.0500	<0.0500	<0.0500	<0.0500	<0.250	<0.0500	<0.0500	<0.0500	<0.500	<0.0500	<0.0500		
Equipment Blank	04/04/2022	<0.000625	--	<0.000625	<0.000625	<0.000625	<0.000625	<0.000625	<0.000625	<0.000625	<0.000625	<0.000313	<0.000625	<0.000625	<0.000625	<0.000625	<0.000625	<0.000625		

Notes:

PAHs = Polycyclic Aromatic Hydrocarbons by United States Environmental Protection Agency Method EPA 8270E-SIM.

ADEC = Alaska Department of Environmental Conservation

µg/L = micrograms per liter

<0.000500 = Not detected at or above the reported detection limit (RDL)

**Bold** = Value exceeds Laboratory Method Detection Limit (MDL)

**Bold and Shaded** = Value exceeds ADEC Groundwater Cleanup Level

J = The compound was positively identified; however, the associated numerical value is an estimated concentration only

The laboratory for this site was changed from Eurofins Calscience to Pace Analytical prior to the second quarter 2020 groundwater monitoring event.



















**Table 4. Historical Groundwater Gauging and Analytical Results****First Quarter 1992 to Current**

Former Chevron-Branded Service Station 97324  
 4417 Lake Otis Parkway  
 Anchorage, Alaska

Well ID	Sample Date	TOC (ft amsl)	DTW (ft bTOC)	LNAPL Thickness (ft)	GW Elev (ft amsl)	TPH-d (mg/L)	TPH-g (mg/L)	Benzene (mg/L)	Toluene (mg/L)	Ethylbenzene (mg/L)	Total Xylenes (mg/L)	MTBE (mg/L)	Naphthalene (mg/L)	Comments
<b>ADEC Groundwater Cleanup Levels</b>														
Trip Blank	4/22/2020	--	--	--	--	1.5	2.2	<b>0.0046</b>	1.1	<b>0.015</b>	<b>0.19</b>	<b>0.14</b>	<b>0.0017</b>	
Trip Blank	10/9/2020	--	--	--	--	--	<0.100	<0.00100	<0.00100	<0.00100	<0.00300	<0.00100	<0.00500	
Trip Blank	8/26/2021	--	--	--	--	--	<b>0.0429 J</b>	<0.00100	<0.00100	<0.00100	<0.00300	<0.00100	<b>0.00124 J</b>	
Trip Blank	4/4/2022	--	--	--	--	--	<0.1	<0.00100	<0.00100	<0.00100	<0.00300	<0.00100 J	<0.00500 J	
Tudor Motel	9/21/2007	--	--	--	--	--	--	--	--	--	--	--	--	
Tudor Motel	5/1/2008	--	--	--	--	--	--	--	--	--	--	--	--	
Tudor Motel	7/15/2008	--	--	--	--	--	--	--	--	--	--	--	--	
Equipment Blank	9/11/2019	--	--	--	--	<0.076	<0.100	<b>0.000013 J</b>	<b>0.0011 J</b>	<0.00050	<0.00114	<0.00044	<b>0.000030 J*B</b>	
Equipment Blank	4/22/2020	--	--	--	--	<0.800	<0.100	<0.00100	<0.00100	<0.00100	<0.00300	<0.00100	<0.00500	
Equipment Blank	10/9/2020	--	--	--	--	<0.800	<0.100	<0.00100	<0.00100	<0.00100	<0.00300	<0.00100	<0.00500	
Equipment Blank	8/26/2021	--	--	--	--	0.624 J	<0.100	<0.00100	<0.00100	<0.00100	<0.00300	<0.00100	<0.00500	
Equipment Blank	4/4/2022	--	--	--	--	<0.8	<0.1	<0.00100	<0.00100	<0.00100	<0.00300	<0.00100 J	<0.00500 J	

**Notes:**

ID = Identification

MW = Groundwater monitoring well

TOC = Top of casing

DTW = Depth to groundwater

ft bTOC = Feet below top of casing

ft = Feet relative to NAVD88

mg/L = Milligrams per liter

GW Elev = Groundwater elevation

&lt;0.00100 = Not detected at or above the reported detection limit (RDL)

**Bold** = Detected above laboratory method detection limit (MDL)**Bold and Shaded** = Value exceeds ADEC Groundwater Cleanup Level**Bold and Italicized** : Constituent considered non-detect, however Laboratory RDL is greater than the ADEC Groundwater Cleanup Level

[ ] = Blind Duplicate Sample Result

\* = LCS or LCSD is outside acceptance limits.

ND = Constituent considered non detect at the MDL

TPH-g = Total petroleum hydrocarbons, gasoline range by LUFT GC/MS according to United States Environmental Protection Agency (USEPA) Method AK101

TPH-d = Total petroleum hydrocarbons, diesel range by LUFT GC/MS according to State of Alaska Method AK102.

Samples analytes by USEPA Method 8260D:

Benzene, Toluene, Ethylbenzene and Total Xylenes (collectively BTEX)

MTBE = Methyl-tert-butyl ether

Naphthalene

LUFT = Leaking Underground Fuel Tank

GC/MS = Gas chromatography/Mass Spectrometry

J = The compound was positively identified; however, the associated numerical value is an estimated concentration only.

B = Compound considered non-detect at the listed value due to associated blank contamination.

ADEC = Alaska Department of Environmental Conservation

NAVD 88 = North American Vertical Datum of 1988

LNAPL = Light Non-Aqueous Phase Liquid

-- = Not Measured/Not analysed

The laboratory for this site was changed from Eurofins Calscience to Pace Analytical prior to the second quarter 2020 groundwater monitoring event. Prior to this date, Eurofins Calscience was using the carbon ranges as follows: TPH-g as C6-C10; TPH-d as C13-C22. Pace Analytical reports the following carbon ranges: TPH-g as C5-C12; TPH-d as C12-C22.

Table 5a. Historical Groundwater Analytical Results - Additional VOCs

First Quarter 1992 to Current

Former Chevron-Branded Service Station 97324

4417 Lake Otis Parkway

Anchorage, Alaska

Well ID	Sample Date	EDC (mg/L)	TCE (mg/L)	PCE (mg/L)	cis-1,2-DCE (mg/L)	Methylene chloride (mg/L)	Isopropylbenzene mg/L	1,2-Dichlorobenzene (o-Dichlorobenzene) mg/L	trans-1,2-Dichloroethene mg/L	1,1,1-Trichloroethane mg/L	1,1,2,2-Tetrachloroethane mg/L	1,1,2-Trichloroethane mg/L	1,1,2-Trichlorotetrafluoroethane (Freon 113) mg/L	Comments	
ADEC Groundwater Cleanup Levels		0.0017	0.0028	0.041	0.036	0.11	—	0.3	0.36	8	0.00076	0.00041	10		
MW-1R	8/26/2021	<b>0.00311 J</b>	<0.00100	<0.00100 J	<0.00100	<0.00500	<0.00100 J	<0.00100	<0.00100	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100	
MW-1R	04/04/2022	<b>0.00191</b>	<0.00100	<b>0.000905 J</b>	<0.00100	<0.00500	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100 <0.0000500	
MW-2R	8/26/2021	<b>0.0105 J [0.0106 J]</b>	<0.00100 [<0.00100]	<0.00100 J [<0.00100 J]	<0.00100 [<0.00100]	<0.00500 [<0.00500]	<0.0433 J [<0.0426 J]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	
MW-2R	04/04/2022	<b>0.0064 [0.00642]</b>	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00500 [<0.00500]	<b>0.0377 [0.0435]</b>	<0.00100 J [<0.00100 J]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 J [<0.00100 J]	<0.00100 [<0.00100]	<0.00100 [<0.00100] <0.00125 [<0.000125]	
MW-8RR	8/26/2021	<0.00100	<0.00100 J	<b>0.00159 J</b>	<0.00100 J	<0.00500	<b>0.000295 J</b>	<0.00100	<0.00100	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100 J	0.00012
MW-8RR	04/04/2022	<b>0.000622 J</b>	<0.00100	<b>0.00183</b>	<0.00100	<0.00500	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100	
MW-9	8/26/2021	<0.00100	<b>0.0135</b>	<b>0.0452 J</b>	<b>0.0376</b>	<0.00500	<0.00100 J	<0.00100	<b>0.000275 J</b>	<0.00100	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100
MW-9	04/04/2022	<0.00100	<b>0.0101</b>	<b>0.0373</b>	<b>0.0263</b>	<0.00500	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.000125
Trip Blank	8/26/2021	<0.00100	<0.00100	<0.00100 J	<0.00100	<0.00500	<0.00100 J	<0.00100	<0.00100	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100	
Trip Blank	04/04/2022	<0.00100	<0.00100	<0.00100	<0.00100	<0.00500	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100	<0.0000500
Equipment Blank	8/26/2021	<0.00100	<0.00100	<0.00100 J	<0.00100	<0.00500	<0.00100 J	<0.00100	<0.00100	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100	
Equipment Blank	04/04/2022	<0.00100	<0.00100	<0.00100	<0.00100	<0.00500	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100 J	<0.00100	<0.00100	<0.00100	<0.0000500

**Notes:**

ID = Identification

MW = Groundwater monitoring well

mg/L = Milligrams per liter

&lt;0.00500 = Not detected at or above the Reported Detection Limit

**Bold** = Detected above laboratory method detection limit (MDL)**Bold and Shaded** = Value exceeds ADEC Groundwater Cleanup Level**Bold and Italicized** : Constituent considered non-detect, however Laboratory RDL is greater than the ADEC Groundwater Cleanup Level

[] = Blind Duplicate Sample Result

ADEC = Alaska Department of Environmental Conservation

Constituents analyzed by United States Environmental Protection Agency Method 8260D

EDB-1,2 Dibromoethane

EDC-1,2-Dichloroethane

TCE = Trichloroethylene

PCE = Tetrachloroethylene

Table 5b. Historical Groundwater Analytical Results - Additional VOCs

First Quarter 1992 to Current

Former Chevron-Branded Service Station 97324

4417 Lake Otis Parkway

Anchorage, Alaska

Well ID	Sample Date	1,1-Dichloroethane mg/L	1,1-Dichloroethylene (Dichloroethylene) mg/L	1,2,3-Trichlorobenzene mg/L	1,2,4-Trichlorobenzene mg/L	1,2,4-Trimethylbenzene mg/L	1,2-Dibromoethane mg/L	1,2-Dichloropropane mg/L	1,3-Dichlorobenzene mg/L	1,4-Dichlorobenzene mg/L	2-Butanone (Methyl ethyl ketone) mg/L	4-Methyl-2-pentanone mg/L	Acetone mg/L	Comments
<b>ADEC Groundwater Cleanup Levels</b>														
		<b>0.028</b>	<b>0.28</b>	<b>0.007</b>	<b>0.004</b>	<b>0.056</b>	<b>0.000075</b>	<b>0.0082</b>	<b>0.0047</b>	<b>0.0048</b>	--	<b>6.3</b>	<b>14</b>	
MW-1R	8/26/2021	<0.00100	<0.00100	<0.00100	<0.00100	<0.00100	<0.00000500	<0.00100	<0.00100	<0.00100	<0.0100	<0.0100	<0.0500	
MW-1R	04/04/2022	<0.00100	<0.00100	<0.00100 J	<0.00100 J	<0.00100 J	<0.00000500	<0.00100	<0.00100 J	<0.00100 J	<0.0100 J	<0.0100 J	<0.0500 J	
MW-2R	8/26/2021	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<b>0.092 [0.0853]</b>	<b>&lt;0.000125 [&lt;0.000125]</b>	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.0100 [<0.0100]	<0.0100 [<0.0100]	<0.0500 [<0.0500]	
MW-2R	04/04/2022	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 J [<0.00100 J]	<0.00100 J [<0.00100 J]	<b>0.0178 J [ 0.0195 J ]</b>	<b>&lt;0.000125 [&lt;0.000125]</b>	<0.00100 [<0.00100]	<0.00100 J [<0.00100 J]	<0.00100 J [<0.00100 J]	<0.0100 J [<0.0100 J]	<0.0100 J [<0.0100 J]	<0.0500 J [<0.0500 J]	
MW-8RR	8/26/2021	<0.00100 J	<0.00100 J	<0.00100	<0.00100	<b>0.000995 J</b>	<b>0.000006</b>	<0.00100	<0.00100	<0.00100	<0.0100	<0.0100	<0.0500	
MW-8RR	04/04/2022	<0.00100	<0.00100	<0.00100 J	<0.00100 J	<0.00100 J	<b>0.000012</b>	<0.00100	<0.00100 J	<0.00100 J	<0.0100 J	<0.0100 J	<0.0500 J	
MW-9	8/26/2021	<0.00100	<0.00100	<0.00100	<0.00100	<0.00100	<b>&lt;0.000125</b>	<0.00100	<0.00100	<0.00100	<0.0100	<0.0100	<0.0500	
MW-9	04/04/2022	<0.00100	<0.00100	<0.00100 J	<0.00100 J	<0.00100 J	<0.000125	<0.00100	<0.00100 J	<0.00100 J	<0.0100 J	<0.0100 J	<0.0500 J	
Trip Blank	8/26/2021	<0.00100	<0.00100	<0.00100	<0.00100	<0.00100	<0.00000500	<0.00100	<0.00100	<0.00100	<0.0100	<0.0100	<0.0500	
Trip Blank	04/04/2022	<0.00100	<0.00100	<0.00100 J	<0.00100 J	<0.00100 J	<0.00000500	<0.00100	<0.00100 J	<0.00100 J	<0.0100 J	<0.0100 J	<0.0500 J	
Equipment Blank	8/26/2021	<0.00100	<0.00100	<0.00100	<0.00100	<0.00100	<0.00000500	<0.00100	<0.00100	<0.00100	<0.0100	<0.0100	<0.0500	
Equipment Blank	04/04/2022	<0.00100	<0.00100	<0.00100	<0.00100 J	<0.00100 J	<0.00000500	<0.00100	<0.00100 J	<0.00100 J	<0.0100 J	<0.0100 J	<0.0500 J	

**Notes:**

ID = Identification

MW = Groundwater monitoring well

mg/L = Milligrams per liter

&lt;0.00500 = Not detected at or above the Reported Detection Limit

**Bold** = Detected above laboratory method detection limit (MDL)**Bold and Shaded** = Value exceeds ADEC Groundwater Cleanup LevelBold and *Italicized* : Constituent considered non-detect, however Laboratory RDL is greater than the ADEC Groundwater Cleanup Level

[] = Blind Duplicate Sample Result

ADEC = Alaska Department of Environmental Conservation

Constituents analyzed by United States Environmental Protection Agency Method 8260D

**Table 5c. Historical Groundwater Analytical Results - Additional VOCs****First Quarter 1992 to Current**

Former Chevron-Branded Service Station 97324

4417 Lake Otis Parkway

Anchorage, Alaska

Well ID	Sample Date	Bromoform mg/L	Bromochloromethane mg/L	Bromodichloromethane mg/L	Bromomethane (Methyl bromide) mg/L	Carbon Disulfide mg/L	Carbon Tetrachloride mg/L	Chlorobenzene mg/L	Chloroethane mg/L	Chloroform mg/L	Chloromethane (Methyl chloride) mg/L	cis-1,3-Dichloropropene mg/L	Dibromochloromethane mg/L	Comments
<b>ADEC Groundwater Cleanup Levels</b>	--	<b>0.0013</b>	<b>0.033</b>	<b>0.0075</b>	<b>0.81</b>	<b>0.0046</b>	<b>0.078</b>	--	<b>0.0022</b>	<b>0.19</b>	<b>0.0047</b>	<b>0.0087</b>		
MW-1R	8/26/2021	<0.00100	<0.00100	<0.00100 J	<0.00500 J	<0.00100	<0.00100	<0.00100	<0.00500	<0.00500	<0.00250	<0.00100	<0.00100	
MW-1R	04/04/2022	<0.00100	<0.00100	<0.00100 J	<0.00500	<0.00100	<0.00100	<0.00100	<0.00500	<0.00500	<0.00250	<0.00100	<0.00100	
MW-2R	8/26/2021	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 J [<0.00100 J]	<0.00500 J [<0.00500 J]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00500 [<0.00500]	<0.00500 [<0.00500]	<0.00250 [<0.00250]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	
MW-2R	04/04/2022	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 J [<0.00100 J]	<0.00500 [<0.00500]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00500 [<0.00500]	<0.00500 [<0.00500]	<0.00250 [<0.00250]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	
MW-8RR	8/26/2021	<0.00100	<0.00100	<0.00100 J	<0.00500 J	<0.00100 J	<0.00100 J	<0.00100	<0.00500 J	<0.00500 J	<0.00250 J	<0.00100	<0.00100	
MW-8RR	04/04/2022	<0.00100	<0.00100	<0.00100 J	<0.00500	<0.00100	<0.00100	<0.00100	<0.00500	<0.00500	<0.00250	<0.00100	<0.00100	
MW-9	8/26/2021	<0.00100	<0.00100	<0.00100 J	<0.00500 J	<0.00100	<0.00100	<0.00100	<0.00500	<0.00500	<0.00250	<0.00100	<0.00100	
MW-9	04/04/2022	<0.00100	<0.00100	<0.00100 J	<0.00500	<0.00100	<0.00100	<0.00100	<0.00500	<0.00500	<0.00250	<0.00100	<0.00100	
Trip Blank	8/26/2021	<0.00100	<0.00100	<0.00100 J	<0.00500 J	<0.00100	<0.00100	<0.00100	<0.00500	<0.00500	<0.00250	<0.00100	<0.00100	
Trip Blank	04/04/2022	<0.00100	<0.00100	<0.00100 J	<0.00500	<0.00100	<0.00100	<0.00100	<0.00500	<0.00500	<0.00250	<0.00100	<0.00100	
Equipment Blank	8/26/2021	<0.00100	<0.00100	<0.00100 J	<0.00500 J	<0.00100	<0.00100	<0.00100	<0.00500	<0.00500	<0.00250	<0.00100	<0.00100	
Equipment Blank	04/04/2022	<0.00100	<0.00100	<0.00100 J	<0.00500	<0.00100	<0.00100	<0.00100	<0.00500	<0.00500	<0.00250	<0.00100	<0.00100	

**Notes:**

ID = Identification

MW = Groundwater monitoring well

mg/L = Milligrams per liter

&lt;0.00500 = Not detected at or above the Reported Detection Limit

**Bold** = Detected above laboratory method detection limit (MDL)**Bold and Shaded** = Value exceeds ADEC Groundwater Cleanup LevelBold and *Italicized* : Constituent considered non-detect, however Laboratory RDL is greater than the ADEC Groundwater Cleanup Level

[] = Blind Duplicate Sample Result

ADEC = Alaska Department of Environmental Conservation

Constituents analyzed by United States Environmental Protection Agency Method 8260D

**Table 5d. Historical Groundwater Analytical Results - Additional VOCs****First Quarter 1992 to Current**

Former Chevron-Branded Service Station 97324

4417 Lake Otis Parkway

Anchorage, Alaska

Well ID	Sample Date	Dichlorodifluoromethane (Freon 12) mg/L	Styrene mg/L	trans-1,3-Dichloropropene mg/L	Trichlorofluoromethane (Freon 11) mg/L	Vinyl chloride (Chloroethene) mg/L	Comments
<b>ADEC Groundwater Cleanup Levels</b>							
<b>MW-1R</b>	8/26/2021	<0.00500	<0.00100	<0.00100	<0.00500	<b>&lt;0.00100</b>	
<b>MW-1R</b>	04/04/2022	<0.00500	<0.00100	<0.00100	<0.00500	<b>&lt;0.00100</b>	
<b>MW-2R</b>	8/26/2021	<0.00500 [<0.00500]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00500 [<0.00500]	<b>&lt;0.00100 [&lt;0.00100]</b>	
<b>MW-2R</b>	04/04/2022	<0.00500 [<0.00500]	<0.00100 [<0.00100]	<0.00100 [<0.00100]	<0.00500 [<0.00500]	<b>&lt;0.00100 [&lt;0.00100]</b>	
<b>MW-8RR</b>	8/26/2021	<0.00500 J	<0.00100	<0.00100	<0.00500	<b>&lt;0.00100 J</b>	
<b>MW-8RR</b>	04/04/2022	<0.00500	<0.00100	<0.00100	<0.00500	<b>&lt;0.00100</b>	
<b>MW-9</b>	8/26/2021	<0.00500	<0.00100	<0.00100	<0.00500	<b>&lt;0.00100</b>	
<b>MW-9</b>	04/04/2022	<0.00500	<0.00100	<0.00100	<0.00500	<b>&lt;0.00100</b>	
<b>Trip Blank</b>	8/26/2021	<0.00500	<0.00100	<0.00100	<0.00500	<b>&lt;0.00100</b>	
<b>Trip Blank</b>	04/04/2022	<0.00500	<0.00100	<0.00100	<0.00500	<b>&lt;0.00100</b>	
<b>Equipment Blank</b>	8/26/2021	<0.00500	<0.00100	<0.00100	<0.00500	<b>&lt;0.00100</b>	
<b>Equipment Blank</b>	04/04/2022	<0.00500	<0.00100	<0.00100	<0.00500	<b>&lt;0.00100</b>	

**Notes:**

ID = Identification

MW = Groundwater monitoring well

mg/L = Milligrams per liter

&lt;0.00500 = Not detected at or above the Reported Detection Limit

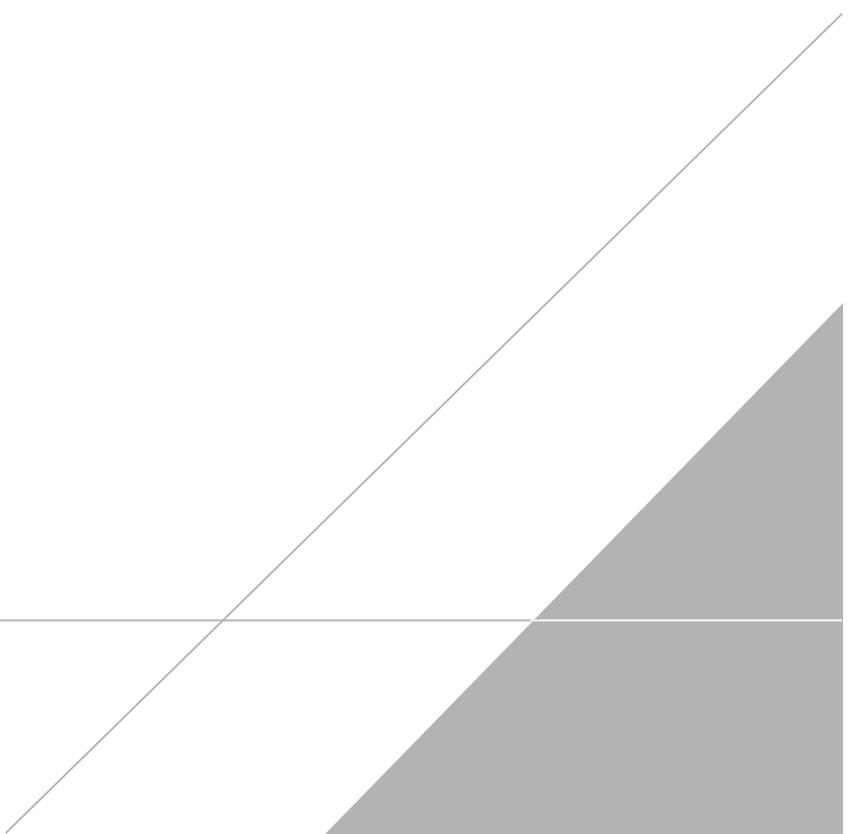
**Bold** = Detected above laboratory method detection limit (MDL)**Bold and Shaded** = Value exceeds ADEC Groundwater Cleanup Level**Bold and *Italicized*** : Constituent considered non-detect, however Laboratory RDL is greater than the ADEC Groundwater Cleanup Level

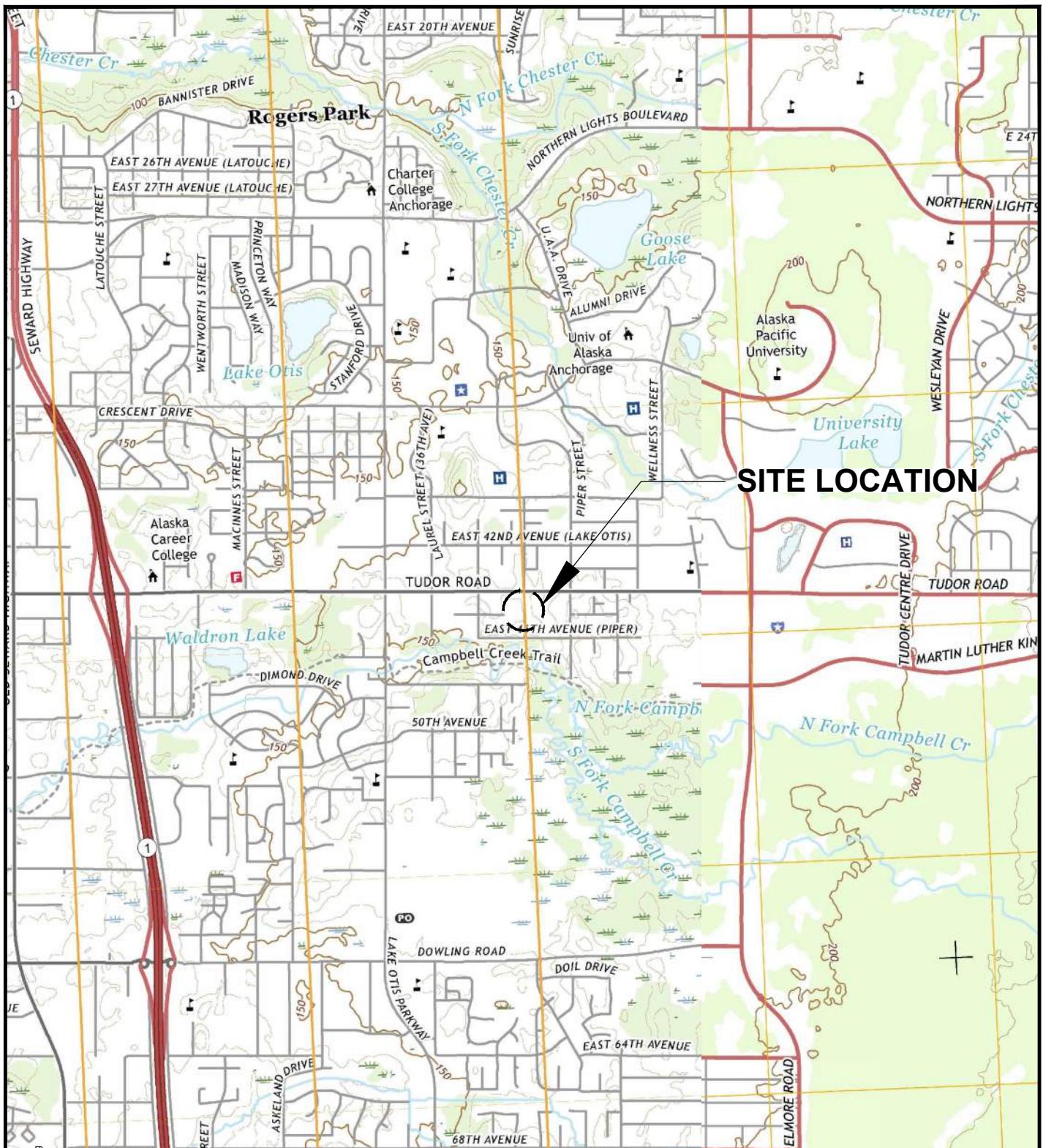
[ ] = Blind Duplicate Sample Result

ADEC = Alaska Department of Environmental Conservation

Constituents analyzed by United States Environmental Protection Agency Method 8260D

## FIGURES





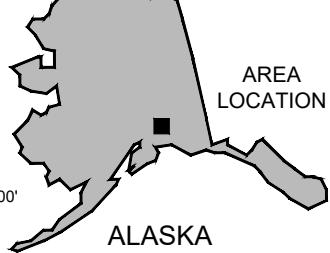
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IMAGES: ---  
REFS: ---  
PM BY: Y.M. BAEBU  
XREFS: ---  
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AK\_Anchorage\_A-8\_NW\_20190703\_TM.gdb

FORMER CHEVRON-BRANDED SERVICE STATION 97324  
4417 LAKE OTIS PARKWAY  
ANCHORAGE, ALASKA

## SITE LOCATION MAP

 ARCADIS

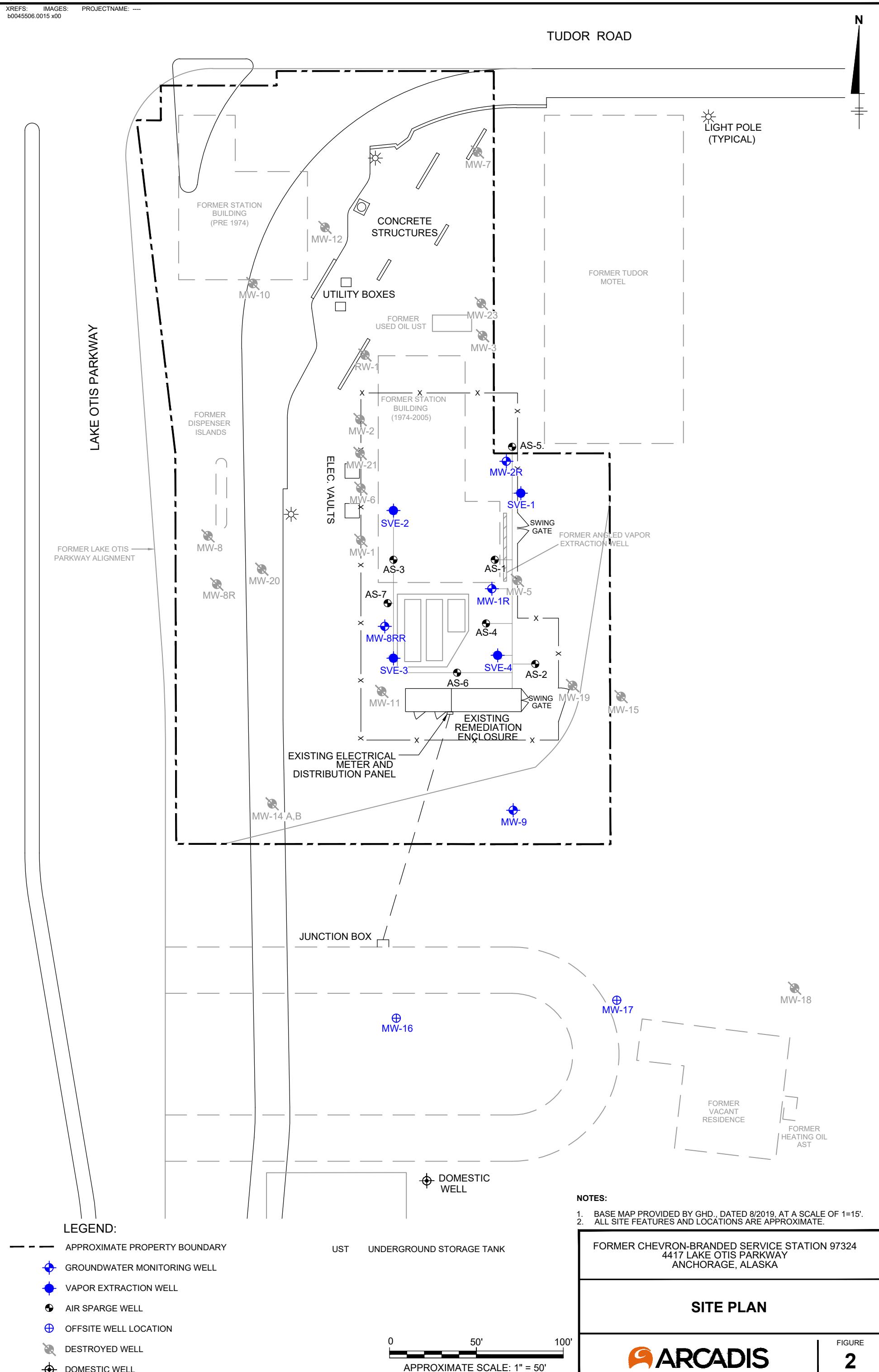
FIGURE  
1



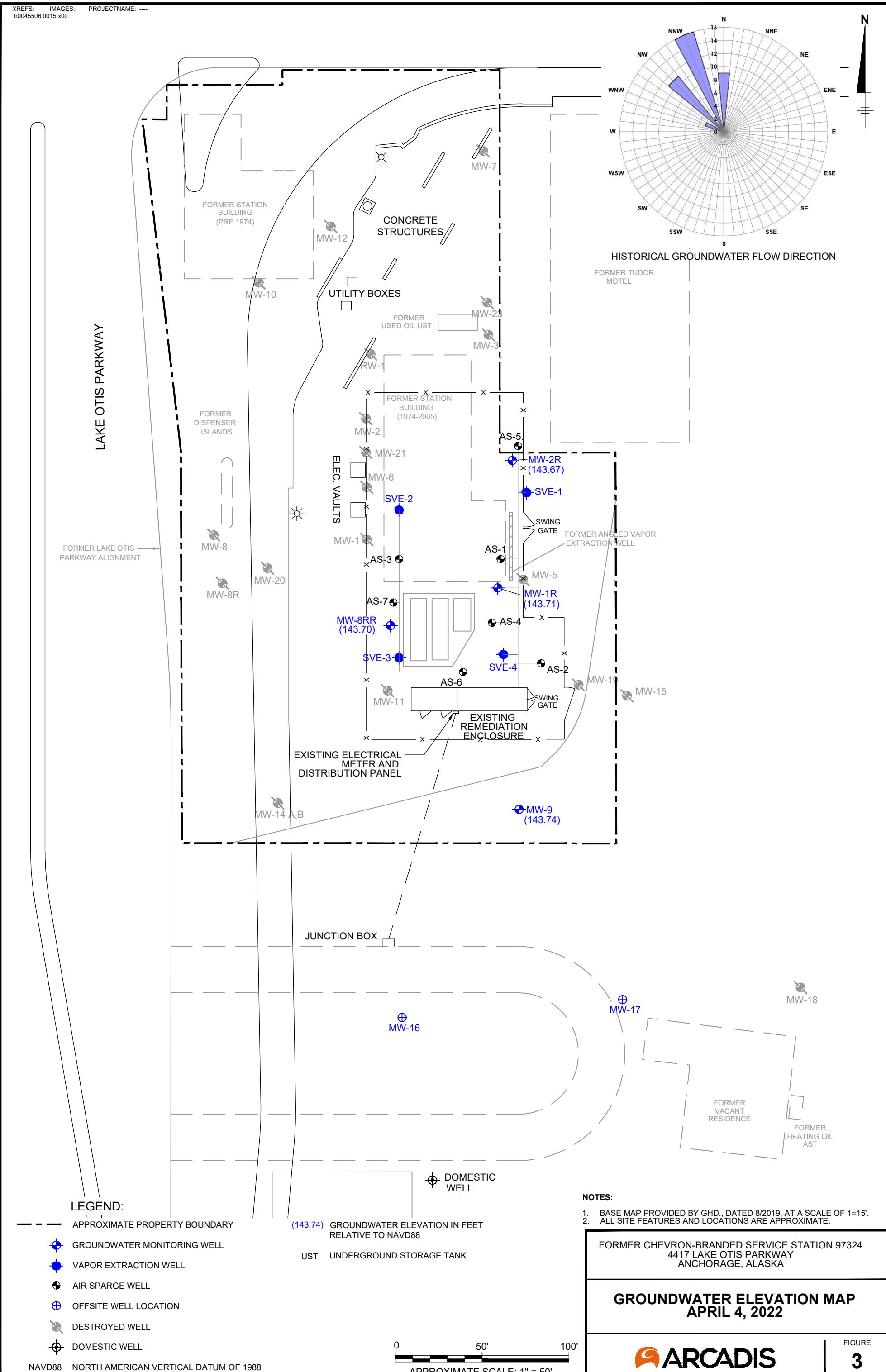
APPROXIMATE GRAPHIC SCALE

2000'

4000'



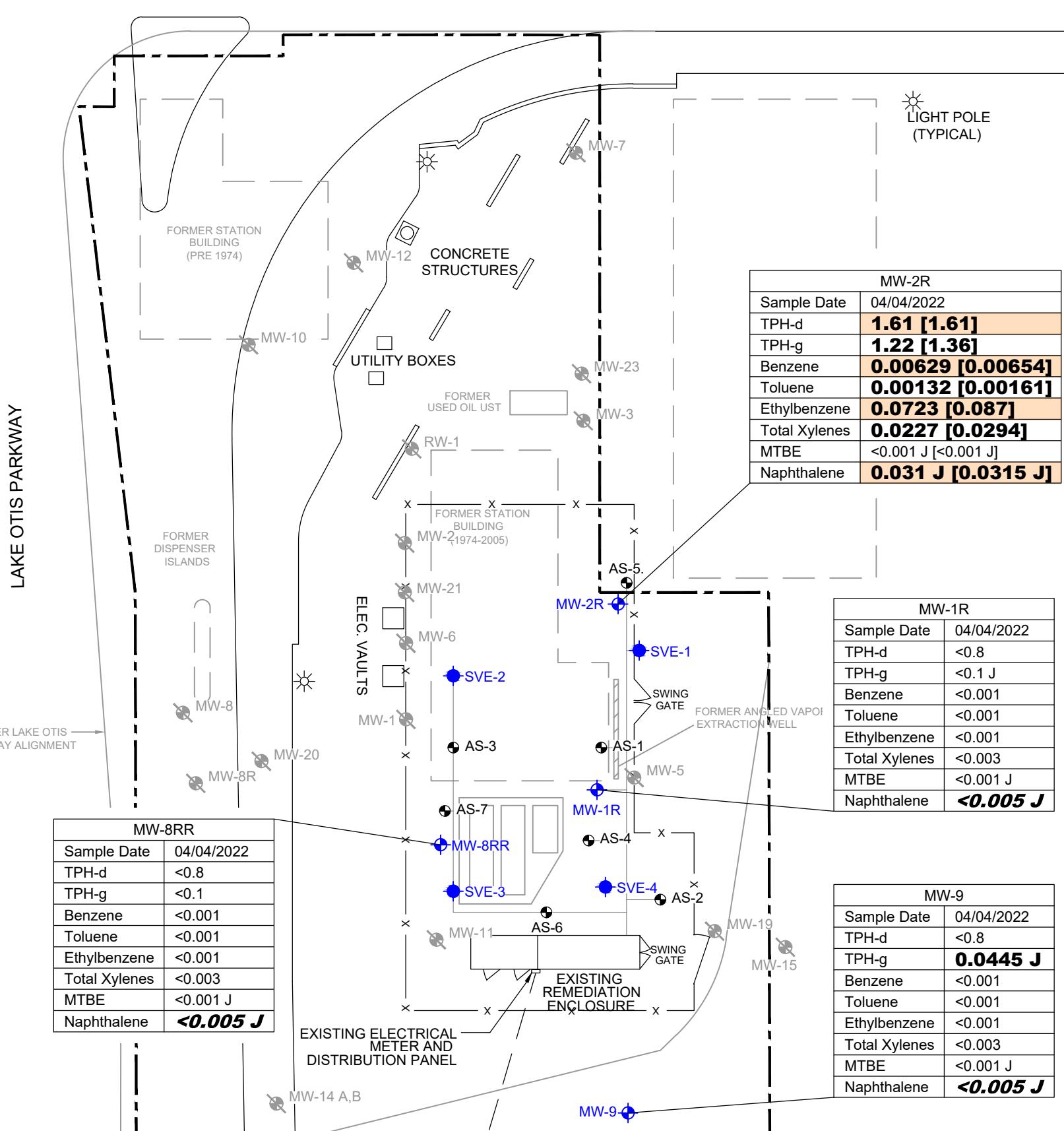
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XREFS: IMAGES: PROJECTNAME: ----  
b0045506.0015 x00

TUDOR ROAD

N



#### LEGEND:

— APPROXIMATE PROPERTY BOUNDARY

⊕ GROUNDWATER MONITORING WELL

● VAPOR EXTRACTION WELL

● AIR SPARGE WELL

⊕ OFFSITE WELL LOCATION

● DESTROYED WELL

● DOMESTIC WELL

UST UNDERGROUND STORAGE TANK

TPH-d TOTAL PETROLEUM HYDROCARBONS DIESEL RANGE ORGANICS

TPH-g TOTAL PETROLEUM HYDROCARBONS GASOLINE RANGE ORGANICS

MTBE METHYL-TERT-BUTYL ETHER

mg/L MILLIGRAMS PER LITER

<0.001 NOT DETECTED AT OR ABOVE THE REPORTED DETECTION LIMIT (RDL)

J THE COMPOUND WAS POSITIVELY IDENTIFIED; HOWEVER, THE ASSOCIATED NUMERICAL VALUE IS AN ESTIMATED CONCENTRATION ONLY

[ ] BLIND DUPLICATE SAMPLE RESULT

**BOLD** DETECTED ABOVE LABORATORY METHOD DETECTION LIMIT (MDL)

**BOLD** CONSTITUENT CONSIDERED NON-DETECT, HOWEVER LABORATORY RDL IS GREATER THAN THE ADEC GROUNDWATER CLEANUP LEVEL

**BOLD** VALUE EXCEEDS ADEC GROUNDWATER CLEANUP LEVEL

CONCENTRATION  
IN (mg/L)

Analyte	ADEC Groundwater Cleanup Levels
TPH-d	1.5
TPH-g	2.2
Benzene	0.0046
Toluene	1.1
Ethylbenzene	0.015
Total Xylenes	0.19
MTBE	0.14
Naphthalene	0.0017

0 50' 100'  
APPROXIMATE SCALE: 1" = 50'

#### NOTES:

1. BASE MAP PROVIDED BY GHD., DATED 8/2019, AT A SCALE OF 1=15'.
2. ALL SITE FEATURES AND LOCATIONS ARE APPROXIMATE.

CHEVRON-BRANDED SERVICE STATION 97324  
4417 LAKE OTIS PARKWAY  
ANCHORAGE, ALASKA

#### GROUNDWATER ANALYTICAL RESULTS MAP APRIL 4, 2022

## APPENDIX A



**Chevron Environmental  
Management Company**

**Appendix A:**  
**Site History and Background**

**Former Chevron Facility 97324**

4417 Lake Otis Parkway

Anchorage, Alaska

ADEC File No: 2100.26.008

HAZARD ID No: 23885

June 19, 2020

## Appendix A: 97324 Site Description and Background

# 1 97324 SITE BACKGROUND AND HISTORY

## 1.1 Site Description and Vicinity

Former Chevron Facility 97324 is located at 4417 Lake Otis Parkway in Anchorage, Alaska. The site was formerly operated as a Chevron-branded service station with three underground storage tanks (UST), two dispenser islands, and a station building with an auto service bay. The surrounding properties are mixed commercial and industrial; the site is bordered to the north, west, and south by former or current ADEC contaminated sites.

## 1.2 Site History

In 2004, the facility building, three petroleum underground storage tanks (USTs) equipped with dispenser pumps, and product lines were removed from the property. A remediation system consisting of seven air sparge (AS) wells and four soil vapor extraction (SVE) wells was operated seasonally until 2017, when it was shut down.

# 2 SITE CHARACTERIZATIONS

A soil and groundwater remediation system which included seven air sparge (AS) wells and four soil vapor extraction (SVE) wells was shut down in 2017. Currently, six groundwater monitoring wells remain in place, four of which are sampled and monitored semiannually.

# 3 CURRENT SITE MONITORING ACTIVITIES

The site currently has a network of six monitoring wells; four wells are monitored and sampled semiannually (MW-1R, MW-2R, MW-8RR, and MW-9). Historically, concentrations of volatile organic compounds (VOCs), gasoline range organics (GRO), and diesel range organics (DRO) have exceeded their respective ADEC Method 2 groundwater cleanup levels in several monitoring wells.

# 4 GEOLOGY AND HYDROGEOLOGY

## 4.1 Site Hydrogeology

The site is in south central Alaska, south of the Knik Arm and north of the Turnagain Arm of Cook Inlet. From 1992 until present, static groundwater depths at the site have ranged between 8.58 to 24.53 feet below top of casing (ft btoc). Historic ground water flow is to the northwest.

# 5 REFERENCES

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GHD Inc. 2018. Second Semiannual 2018 Groundwater Monitoring Report Former Chevron-Branded Service Station 97324, 4417 Lake Otis Parkway , Anchorage, AK. December 5

## APPENDIX B



**Project Name :** 97324      **Weather(°F) :** Clear  
**Project Number :** 30063667      **Prepared By:** Evan Wujcik  
**Purpose :** Gw sampling  
**PPE :** Level D  
**Equipment:** Water Quality Meter (i.e. YSI), Water Level Meter (WLM), Bladder Pump, Photoionization Detector (PID)

Date	Time	Description of Activities
04/04/2022	07:00	Arrive on site Open permit to work Locate Wells
04/04/2022	09:00	Sample MW-2R Decon equipment Blind duplicate Samples collected at this location See chain of custody for analytes
04/04/2022	10:00	Sample MW-8RR Decon equipment See chain of custody for analytes
04/04/2022	11:00	Sample MW-1R Decon equipment MS/MSD Samples collected at this location See chain of custody for analytes
04/04/2022	12:00	Sample MW-9 Decon equipment See chain of custody for analytes
04/04/2022	12:30	Load vehicle Mobilize offsite

**Signature:**

<b>Waste Management:</b>										
Drums On Site										
Date	Are there any waste drums on site?	Number of Drums upon Arrival	Size of Drums	Type of Drums	Condition of Drums	Waste Drummed Today?	Number of drums Created	Size of drums	Condition of Drums	General Waste Comments
04/04/2022	no					no				

**Daily Log****Equipment and Calibration Information:**

Supplier: Pine

Model:

Rental Number:

Calibrated:

Bump  
Checked:Calibration  
Passed:**Water Quality Meter SN:**

Date	Time	Calibrated Fluid and Value	Lot #	Expiration Date	Initial Reading	Final Reading
04/04/2022						

**Equipment and Calibration Information:**

Supplier: Pine

Model:

Rental Number:

Calibrated:

Bump  
Checked:Calibration  
Passed:**PIDSN:**

Date	Time	Calibrated Fluid and Value	Lot #	Expiration Date	Initial Reading	Final Reading
04/04/2022	--					



## Groundwater Gauging Log

Project Number	30063667							
Client:	Chevron							
Site ID:	97324							
Site Location:	Anchorage, Alaska							
Measuring Point:	Top of Casing							
Date(s):	04/04/2022							
Sampler(s):	Evan Wujcik							
Gauging Equipment:	Water Level Meter							
Well ID	Date	Gauging Time	Static Water Level (ft bmp)	Depth to Product (ft bmp)	Total Depth (ft bmp)	PID Reading (ppm)	LNAPL Removed (gal)	Comments
MW-1R	04/04/2022	09:12	23.85	ND	31.00	0	--	--
MW-2R	04/04/2022	08:12	24.58	ND	31.20	0	--	--
MW-8RR	04/04/2022	08:59	22.73	ND	32.50	0	--	--
MW-9	04/04/2022	10:09	15.5	ND	19.30	0	--	--

ft-bmp = feet below measuring point

ND = Not Detected

PID = Photoionization Detector Reading

ppm = parts per million

-- = Not Recorded

<b>Project Number</b>	30063667	<b>Well ID</b>	MW-1R	<b>Date</b>		4/4/2022				
<b>Site Location</b>	Anchorage, Alaska	<b>Site ID</b>	97324	<b>Weather (°F)</b>	Clear	<b>Sampled by</b>	Evan Wujcik			
<b>Measuring Point Description</b>	Top of Casing	<b>Screen Depth Interval (ft-bmp)</b>	-- to --	<b>Casing Diameter (in.)</b>	2	<b>Well Casing Material</b>	PVC			
<b>Static Water Level (ft-bmp)</b>	23.85	<b>Total Depth (ft-bmp)</b>	31	<b>Water Column (ft)</b>	7.15	<b>Gallons in Well</b>	1.16			
<b>Water Quality Meter Make/Model</b>	Horiba U-52	<b>Purge Method</b>	Low-Flow	<b>Sample Method</b>		Grab				
<b>Sample Time</b>	11:00	<b>Well Volumes Purged</b>	0.68	<b>Sample ID</b>	MW-1R-W-20220404	<b>Evacuation Equipment</b>	Bladder			
<b>Purge Start</b>	10:30	<b>Gallons Purged</b>	0.79	<b>Duplicate ID</b>	MS/MSD					
<b>Purge End</b>	10:50	<b>Total Purge Time (h:m)</b>	0:20							
Time	Rate (ml/min)	Depth to Water (ft)	pH (standard units)	Conductivity (mS/cm)	Turbidity (NTU)	Dissolved Oxygen (mg/L)	Temperature (°C)	Redox (mV)	Appearance	
									Color	Odor
10:33	200	23.85	6.56	0.431	60.0	3.43	4.13	122	--	--
10:36	200	23.85	6.48	0.352	49.7	4.59	4.13	132	--	--
10:39	200	23.85	6.39	0.300	40.6	5.39	4.12	140	--	--
10:42	200	23.85	6.35	0.292	35.7	5.55	4.12	144	--	--
10:45	200	23.85	6.32	0.282	33.4	5.51	4.12	146	--	--

**Comments:** None

#### Well Casing Volume Conversion

Well diameter (in.) = 1 = 0.04 1.5 = 0.09 2.5 = 0.26 3.5 = 0.50 6 = 1.47  
 gallons per foot 1.25 = 0.06 2 = 0.16 3 = 0.37 4 = 0.65

#### Sample Information

Sample ID: MW-1R-W-20220404 Sample Time: 11:00 Sample Depth (ft-bmp): 24.5

Analytes and Methods: See Chain-of-Custody.

ft-bmp = feet below measuring point  
 in. = inches  
 ft = feet  
 mL/min = milliliters per minute

mS/cm = millSiemens per centimeter  
 NTU = Nephelometric Turbidity Unit  
 mg/L = milligrams per liter  
 PVC = Polyvinyl Chloride

mV = millivolts  
 °F = degrees Fahrenheit  
 °C = degrees Celsius  
 -- = Not Recorded

<b>Project Number</b>	30063667	<b>Well ID</b>	MW-2R	<b>Date</b>		4/4/2022				
<b>Site Location</b>	Anchorage, Alaska	<b>Site ID</b>	97324	<b>Weather (°F)</b>	Clear	<b>Sampled by</b>	Evan Wujcik			
<b>Measuring Point Description</b>	Top of Casing	<b>Screen Depth Interval (ft-bmp)</b>	-- to --	<b>Casing Diameter (in.)</b>	2	<b>Well Casing Material</b>	PVC			
<b>Static Water Level (ft-bmp)</b>	24.58	<b>Total Depth (ft-bmp)</b>	31.2	<b>Water Column (ft)</b>	6.62	<b>Gallons in Well</b>	1.08			
<b>Water Quality Meter Make/Model</b>	Horiba U-52	<b>Purge Method</b>	Low-Flow	<b>Sample Method</b>		Grab				
<b>Sample Time</b>	09:00	<b>Well Volumes Purged</b>	0.59	<b>Sample ID</b>	MW-2R-W-20220404	<b>Evacuation Equipment</b>	Bladder			
<b>Purge Start</b>	08:30	<b>Gallons Purged</b>	0.63	<b>Duplicate ID</b>	BD					
<b>Purge End</b>	08:50	<b>Total Purge Time (h:m)</b>	0:20							
Time	Rate (ml/min)	Depth to Water (ft)	pH (standard units)	Conductivity (mS/cm)	Turbidity (NTU)	Dissolved Oxygen (mg/L)	Temperature (°C)	Redox (mV)	Appearance	
									Color	Odor
08:33	200	24.60	7.26	1.13	16.8	0.95	3.55	31	--	--
08:36	200	24.60	7.28	1.10	16.0	0.80	3.50	27	--	--
08:39	200	24.60	7.30	1.09	15.6	0.76	3.48	25	--	--
08:42	200	24.60	7.29	1.08	15.0	0.73	3.45	24	--	--

**Comments:** None

#### Well Casing Volume Conversion

Well diameter (in.) = 1 = 0.04 1.5 = 0.09 2.5 = 0.26 3.5 = 0.50 6 = 1.47  
 gallons per foot 1.25 = 0.06 2 = 0.16 3 = 0.37 4 = 0.65

#### Sample Information

Sample ID:	MW-2R-W-20220404	Sample Time:	09:00	Sample Depth (ft-bmp):	25
Analytes and Methods:	See Chain-of-Custody.				

ft-bmp = feet below measuring point  
 in. = inches  
 ft = feet  
 mL/min = milliliters per minute

mS/cm = millSiemens per centimeter  
 NTU = Nephelometric Turbidity Unit  
 mg/L = milligrams per liter  
 PVC = Polyvinyl Chloride

mV = millivolts  
 °F = degrees Fahrenheit  
 °C = degrees Celsius  
 -- = Not Recorded

<b>Project Number</b>	30063667	<b>Well ID</b>	MW-8RR	<b>Date</b>		4/4/2022				
<b>Site Location</b>	Anchorage, Alaska	<b>Site ID</b>	97324	<b>Weather (°F)</b>	Clear	<b>Sampled by</b>	Evan Wujcik			
<b>Measuring Point Description</b>	Top of Casing	<b>Screen Depth Interval (ft-bmp)</b>	-- to --	<b>Casing Diameter (in.)</b>	2	<b>Well Casing Material</b>	PVC			
<b>Static Water Level (ft-bmp)</b>	22.73	<b>Total Depth (ft-bmp)</b>	32.5	<b>Water Column (ft)</b>	9.77	<b>Gallons in Well</b>	1.59			
<b>Water Quality Meter Make/Model</b>	Horiba U-52	<b>Purge Method</b>	Low-Flow	<b>Sample Method</b>			Grab			
<b>Sample Time</b>	10:00	<b>Well Volumes Purged</b>	0.50	<b>Sample ID</b>	MW-8RR-W-20220404	<b>Evacuation Equipment</b>	Bladder			
<b>Purge Start</b>	09:30	<b>Gallons Purged</b>	0.79	<b>Duplicate ID</b>	--					
<b>Purge End</b>	09:50	<b>Total Purge Time (h:m)</b>	0:20							
Time	Rate (ml/min)	Depth to Water (ft)	pH (standard units)	Conductivity (mS/cm)	Turbidity (NTU)	Dissolved Oxygen (mg/L)	Temperature (°C)	Redox (mV)	Appearance	
									Color	Odor
09:33	200	22.73	6.96	0.831	92.0	3.83	3.88	54	--	--
09:36	200	22.73	6.88	0.828	75.6	4.05	4.00	72	--	--
09:39	200	22.73	6.81	0.827	60.7	4.15	4.05	82	--	--
09:42	200	22.73	6.76	0.825	40.4	4.09	4.08	85	--	--
09:45	200	22.73	6.74	0.822	39.0	4.02	4.11	87	--	--

**Comments:** None

#### Well Casing Volume Conversion

Well diameter (in.) = 1 = 0.04 1.5 = 0.09 2.5 = 0.26 3.5 = 0.50 6 = 1.47  
 gallons per foot 1.25 = 0.06 2 = 0.16 3 = 0.37 4 = 0.65

#### Sample Information

Sample ID:	MW-8RR-W-20220404	Sample Time:	10:00	Sample Depth (ft-bmp):	23
Analytics and Methods:	See Chain-of-Custody.				

ft-bmp = feet below measuring point  
 in. = inches  
 ft = feet  
 mL/min = milliliters per minute

mS/cm = millSiemens per centimeter  
 NTU = Nephelometric Turbidity Unit  
 mg/L = milligrams per liter  
 PVC = Polyvinyl Chloride

mV = millivolts  
 °F = degrees Fahrenheit  
 °C = degrees Celsius  
 -- = Not Recorded

<b>Project Number</b>	30063667	<b>Well ID</b>	MW-9	<b>Date</b>		4/4/2022				
<b>Site Location</b>	Anchorage, Alaska	<b>Site ID</b>	97324	<b>Weather (°F)</b>	Clear	<b>Sampled by</b>	Evan Wujcik			
<b>Measuring Point Description</b>	Top of Casing	<b>Screen Depth Interval (ft-bmp)</b>	-- to --	<b>Casing Diameter (in.)</b>	2	<b>Well Casing Material</b>	PVC			
<b>Static Water Level (ft-bmp)</b>	15.5	<b>Total Depth (ft-bmp)</b>	19.3	<b>Water Column (ft)</b>	3.80	<b>Gallons in Well</b>	0.62			
<b>Water Quality Meter Make/Model</b>	Horiba U-52	<b>Purge Method</b>	Low-Flow	<b>Sample Method</b>		Grab				
<b>Sample Time</b>	12:00	<b>Well Volumes Purged</b>	1.02	<b>Sample ID</b>	MW-9-W-20220404	<b>Evacuation Equipment</b>	Bladder			
<b>Purge Start</b>	11:30	<b>Gallons Purged</b>	0.63	<b>Duplicate ID</b>	--					
<b>Purge End</b>	11:50	<b>Total Purge Time (h:m)</b>	0:20							
Time	Rate (ml/min)	Depth to Water (ft)	pH (standard units)	Conductivity (mS/cm)	Turbidity (NTU)	Dissolved Oxygen (mg/L)	Temperature (°C)	Redox (mV)	Appearance	
									Color	Odor
11:33	200	15.50	6.39	0.317	114	6.26	3.30	152	--	--
11:36	200	15.50	6.27	0.321	99.4	5.52	3.32	161	--	--
11:39	200	15.50	6.24	0.324	80.4	4.81	3.33	164	--	--
11:42	200	15.50	6.21	0.325	72.3	4.60	3.30	166	--	--

**Comments:** None

#### Well Casing Volume Conversion

Well diameter (in.) = 1 = 0.04 1.5 = 0.09 2.5 = 0.26 3.5 = 0.50 6 = 1.47  
 gallons per foot 1.25 = 0.06 2 = 0.16 3 = 0.37 4 = 0.65

#### Sample Information

Sample ID:	MW-9-W-20220404	Sample Time:	12:00	Sample Depth (ft-bmp):	16
Analytes and Methods:	See Chain-of-Custody.				

ft-bmp = feet below measuring point  
 in. = inches  
 ft = feet  
 mL/min = milliliters per minute

mS/cm = millSiemens per centimeter  
 NTU = Nephelometric Turbidity Unit  
 mg/L = milligrams per liter  
 PVC = Polyvinyl Chloride

mV = millivolts  
 °F = degrees Fahrenheit  
 °C = degrees Celsius  
 -- = Not Recorded

## APPENDIX C



# **SOP – Investigation Derived Waste (IDW) Treatment Utilizing Granular Activated Carbon (GAC)**

Rev: 0

Rev Date: May 24, 2022

## Version Control

Issue	Revision No.	Date Issued	Page No.	Description	Reviewed By

## Approval Signatures

Prepared by:

  
\_\_\_\_\_  
Elysha Nygaard

5/24/2022

Date

Reviewed by:

  
\_\_\_\_\_  
Nick Wood, P.E.

5/24/2022

Date

  
\_\_\_\_\_  
Gerald Robinson

5/24/2022

Date

## 1 Introduction

The objective of this Standard Operating Procedure (SOP) is to describe methods and procedures utilizing granular activated carbon (GAC) for treatment of purged groundwater produced during groundwater monitoring and sampling activities. This SOP is only applicable to impacted groundwater considered to be non-hazardous by the Resource Conservation and Recovery Act (RCRA) regulations. The United States Environmental Protection Agency (USEPA) Operating Procedure for the Management of Investigation Derived Waste (IDW), states that non-hazardous liquids may be discharged to the ground surface as long as doing so does not endanger human health, the environment, or violate state or federal regulations. Management and transportation of purged groundwater may have project, client and/or agency specific requirements and may require approval prior to initiating GAC treatment.

## 2 Intended Use and Responsibilities

This document describes general and/or specific procedures, methods, actions, steps, and considerations to be used and observed by Arcadis staff when performing work, tasks, or actions under the scope and relevancy of this document. This document may describe expectations, requirements, guidance, recommendations, and/or instructions pertinent to the service, work task, or activity it covers.

It is the responsibility of the Arcadis Certified Project Manager (CPM) to provide this document to the persons conducting services that fall under the scope and purpose of this procedure, instruction, and/or guidance. The Arcadis CPM will also ensure that the persons conducting the work falling under this document are appropriately trained and familiar with its content. The persons conducting the work under this document are required to meet the minimum competency requirements outlined herein, and inquire to the CPM regarding any questions, misunderstanding, or discrepancy related to the work under this document.

This document is not considered to be all inclusive nor does it apply to all projects. It is the CPM's responsibility to determine the proper scope and personnel required for each project. There may be project- and/or client- and/or state-specific requirements that may be more or less stringent than what is described herein. The CPM is responsible for informing Arcadis and/or Subcontractor personnel of omissions and/or deviations from this document that may be required for the project. In turn, project staff are required to inform the CPM if or when there is a deviation or omission from work performed as compared to what is described herein.

In following this document to execute the scope of work for a project, it may be necessary for staff to make professional judgment decisions to meet the project's scope of work based upon site conditions, staffing expertise, regulation-specific requirements, health and safety concerns, etc. Staff are required to consult with the CPM when or if a deviation or omission from this document is required that has not already been previously approved by the CPM. Upon approval by the CPM, the staff can perform the deviation or omission as confirmed by the CPM.

## 3 Scope and Application

Activated carbon is one of the most commonly used treatment methodologies for groundwater extraction and treatment remediation systems (USEPA, 2012) and is proven to be a safe engineering practice to remove organic constituents from waste waters. (USEPA, 1991). GAC media can remove petroleum and solvent volatile organic

compounds (VOCs) and other compounds such as per- and polyfluoroalkyl substances (PFAS), chloramines and trihalomethanes from water.

Groundwater wells are purged and treated with GAC media provided that no measurable non-aqueous phase liquid (NAPL) is present. As purge water is passed through the activated carbon contained in a portable vessel, constituents of potential concern (COPCs) adsorb on the surface area of the carbon granules and the treated water passes through the carbon matrix and is discharged to ground surface.

Estimation of the usage rate of GAC to predict breakthrough values relies on site and media specific data. For large scale GAC treatment systems, pilot testing may be conducted to obtain these values. For small scale portable GAC treatment systems across various sites, pilot testing may be cost and time prohibitive, as well as potentially inaccurate due to the variability in concentrations of COPCs. Therefore, Freundlich Isotherm modeling is commonly used to calculate breakthrough estimates of target COPCs.

Use of a portable GAC vessels effectively treats COPCs in generated purge water to desired concentrations applicable to regulatory criteria. Use of a GAC vessels also reduces the overall operational impact to the site and removes the need for drumming, storage, characterization and disposal of purged groundwater.

## 4 Personnel Qualifications

Arcadis field personnel will have completed site-specific training as well as have current health and safety training as required by Arcadis, client, or regulations such as the 40-hour HAZWOPER training and annual 8-hour HAZWOPER refresher. Arcadis personnel will also have up to date training as specified in the Health and Safety Plan (HASP) which may include first aid, fire extinguisher use or COPC specific trainings as needed. In addition, Arcadis personnel conducting work will be knowledgeable in the relevant processes, procedures, scope of work (SOW), standard operating procedures (SOPs) and Technical Guidance Instruction (TGIs), and possess the required skills and experience necessary to successfully and safely complete the work. The HASP and other site-specific SOW documentation will identify other training or work requirements.

## 5 Equipment List

The following field equipment is suggested for treating purged groundwater utilizing a portable GAC vessel:

- Appropriate personal protective equipment (PPE) as specified in the HASP
- Electronic oil-water interface probe and water level indicator with 0.01-foot increments
- Photoionization detector (PID) and/or other air quality measurement equipment as required by the HASP
- Non-phosphate laboratory soap (Alconox® or equivalent)
- Distilled, de-ionized, or potable water for equipment decontamination
- GAC media and portable vessel, including valves and fittings
- 5-gallon buckets
- Tubing
- Groundwater purge equipment such as submersible pumps or bailers
- Plastic drop cloth (e.g. Visqueen) to place beneath portable GAC vessel to reduce potential contamination from spills
- Tools and/or keys for accessing and opening wells
- FieldNow capable device or field logs

- GAC Volume Tracking Form

## 6 Cautions

Electronic water-level indicators and interface probes may sometimes produce false-positive readings. For accuracy, the probe should be raised and lowered several times to verify consistent, repeatable results. Ensure that the type of indicator or probe is compatible with the depth and diameter of the wells to be measured. If the presence of NAPL is suspected, do not use a water level indicator, use an oil-water interface probe. Purge water containing NAPL should not be treated using GAC methods.

Minimize the amount of sediment in purge water by pre-filtering, this will ensure GAC media filters impacted water efficiently.

## 7 Health and Safety Considerations

The site-specific HASP will be followed to ensure the safety of Arcadis field personnel. Access to groundwater wells may expose personnel to hazardous materials such as contaminated groundwater or petroleum compounds. Other potential hazards include pressurized wells, insects or animals that may inhabit the wells, other biological or environmental hazards in the vicinity of the well (e.g. dense vegetation and slope) and potentially the use of sharp tools (e.g. scissors or safety blade). Appropriate PPE will be worn and control measures taken while conducting these activities. Proper lifting and handling techniques will be used when moving portable GAC vessels and related materials.

## 8 Procedure

### Background

Portable vessels containing GAC media are utilized as on-site treatment of impacted purge water during groundwater monitoring and sampling activities. As purge water passes through the carbon granules, VOCs adsorb to the surface area, allowing treated water to pass through and allow for discharge and disposal to ground surface.

Maximum concentrations of site-specific COPCs are utilized to calculate carbon usage rates. To calculate the volume of water that can be treated during purging before breakthrough occurs, Freundlich Isotherm equations are used. The Freundlich Isotherm is an empirically derived adsorption equation relating the concentration of a COPC within the impacted purge water to the concentration of the COPC on the surface of the adsorption material (GAC) (USEPA, 1980). Using these parameters for a given COPC concentration, carbon usage rates can be calculated for the rate and volume of purge water that can be treated per weight of GAC material in a single portable vessel as well as breakthrough estimates. The USEPA provides a free download of the Environmental Technologies Design Option Tool (ETDOT), which includes the Adsorption Design Software (AdDesign<sup>SM</sup>) to model groundwater treatment with GAC using Freundlich Isotherms. Software can be downloaded from the link below:

<https://github.com/USEPA/Environmental-Technologies-Design-Option-Tool>

While this document is intended as general guidance for Arcadis use, a model using the Chevron Environmental Company (CEMC) Alaska Portfolio of Sites is provided in Attachment A as a specific example utilizing the

AdDesignS™ isotherm software to calculate GAC breakthrough to meet the Alaska Department of Environmental Preservation (AEP) criteria.

### **General Assumptions for GAC Usage**

- GAC adsorbs uniformly, entire capacity is utilized prior to breakthrough
- Virgin granular activated carbon produced from coconut shell
- Disregard minor losses due to environmental factors such as ambient temperature (not including extreme temperatures)
- Neutral pH of groundwater (increase carbon bed by 20% per every unit above a pH of 7.0)
- COPCs are dissolved phase; no free product
- Typical density for GAC media is approximately 0.50 – 0.80 g/mL
- Suspended solids are not inhibiting adsorption
- Assume general background concentrations of non-COPCs that may affect COPC adsorption (fluoride, nitrate, phosphate, etc.)
- GAC media can immobilize 4% – 10% of its total mass (DeSilva, 2000)
- Volume of GAC vessel sufficient to ensure adequate contact time between untreated water and GAC media

### **Field Procedures**

Prior to each use of the GAC vessel, the carbon material will be saturated with potable water and allowed to sit for 24 hours. Saturating the carbon opens pore space and increases surface area of the granules, reducing the potential formation of preferential pathways and resulting in optimal performance of the GAC material each time it is used.

Groundwater wells containing measurable NAPL will not be purged or treated through the portable GAC vessel. The presence of NAPL reduces the adsorption capacity and clogs pore space of the carbon media.

Purged groundwater will be pumped through the portable GAC vessel (Attachment C) at no more than the established flow rate (typically 200 to 300 milliliters per minute [mL/min]) to allow for adequate treatment time. As the untreated water begins to purge through the GAC vessel, the valve is opened to allow the treated water to discharge into a labeled waste container prior to discharge. Locations of discharged GAC treated water will be pre-determined and will not be in the immediate vicinity of any surface water, stormwater drains or any other sensitive receptors. After completing discharge of treated purge water, the portable GAC vessel valve will be closed.

Following the completion of groundwater monitoring and sampling activities, field personnel will record the total volume of purged water that has been treated on the GAC Tracking Log (Attachment B), which will remain with the portable GAC vessel until disposal. The following procedure should be followed when using a portable GAC vessel:

- Connect pump or evacuation equipment to well and a portable GAC vessel
- Connect portable GAC vessel outlet to a separate labeled waste container for treated purge water
- Purge impacted groundwater through the portable GAC vessel
- Discharge treated purge water from the GAC vessel into a identified waste container
- Document volume purged on GAC Tracking Log

- Discharge the treated purge water to pre-identified location onsite away from any sensitive receptors
- Replace GAC media prior to reaching calculated breakthrough values
- Store GAC in appropriate labeled container for staging and/or transportation and removal

## 9 Waste Management

Decontamination fluids, used PPE, and other disposable equipment will be properly stored on site in labeled containers and disposed of properly. Used GAC material will be containerized and labeled for transportation off site for re-use, disposal or regeneration. Ensure all waste is properly stored, labeled and documented in field logs. Review the *TGI – Investigation Derived Waste Handling and Storage* for additional information and refer to agency and/or client specific requirements.

## 10 Data Recording and Management

Digital data collection is the Arcadis standard using available FieldNow® applications that enable real-time, paperless data collection, entry, and automated reporting. Paper forms should only be used as backup to FieldNow® digital data collection and/or as necessary to collect data not captured by available FieldNow® applications. The Field Now® digital form applications follow a standardized approach, correlate to most TGIs and are available to all projects accessible with a PC or capable mobile device. Once the digital forms are saved within FieldNow®, the data is instantly available for review on a web interface. This facilitates review by project management team members and SMEs enabling error or anomalous data detection for correction while the staff are still in the field. Continual improvements of FieldNow® applications are ongoing, and revisions are made as necessary in response to feedback from users and subject matter experts.

## 11 Quality Assurance

Conducting sampling of GAC treated effluent purge water may be required to demonstrate that treatment is effective.

## 12 References

- DeSilva, Frank. 2000. Activated Carbon Filtration. Water Quality Products Magazine. January
- USEPA. 1980. Carbon Adsorption Isotherms for Toxic Organics. Wastewater Research Division. Municipal Environmental Research Laboratory, Cincinnati, Ohio. EPA-600/08-80-23. April.
- USEPA. 1991. Engineering Bulletin, Granular Activated Carbon Treatment. Superfund EPA-540/02-91/024. October.
- USEPA 2012. A Citizens Guide to Activated Carbon Treatment. Office of Solid Waste and Emergency Response. September.

### **Attachments**

- |              |  |
|--------------|--|
| Attachment A | Chevron Environmental Management Company – Alaska Portfolio Isotherm Model |
| Attachment B | GAC Tracking Log   |
| Attachment C | Portable GAC Vessel Diagram  |

# **Attachment A**

**Chevron Environmental Management Company – Alaska Portfolio  
Isotherms**

## Attachment A Chevron Environmental Management Company – Alaska Portfolio Isotherm Model

On behalf of Chevron Environmental Management Company (CEMC), Arcadis prepared conservative GAC breakthrough estimates for COPCs observed at Sites located across Anchorage and Fairbanks at the request of the Alaska Department of Environmental Preservation (ADEP). The USEPA recommended AdDesignS™ isotherm software was utilized to model the time until breakthrough is first observed.

### Site Specific Modeling Parameters

To demonstrate the effective treatment of petroleum constituents in groundwater via GAC without confirmation samples, Arcadis has applied conservative parameters to the calculations for GAC breakthrough. The following parameters and conditions have been applied:

- 26 Sites in Arcadis' Chevron Alaska Portfolio are sampled utilizing portable GAC vessels to purge treated groundwater to surface. There are 18 Sites across Anchorage that utilize one portable GAC vessel and 8 Sites across Fairbanks that utilize a second portable GAC vessel. For the purpose of applying the most conservative parameters, the model assumed one portable GAC unit was utilized for all 26 Sites.
- The COPCs that have been detected in approximately half of the 26 Sites were selected to be utilized in the AdDesignS™ software. The highest observed concentration at any of the 26 Sites for each constituent was applied for the duration of the model. Therefore, the model assumes the GAC is exposed to the highest concentration of each COPC for the duration of the modeled run time. The constituents selected were benzene, toluene, ethylbenzene, o-xylene, and naphthalene. Methyl tert-butyl ether (MTBE) was also included in the model scenario, despite detections in only 4 of the 26 Sites due to its impact on carbon adsorption.
- A maximum low-flow methodology flowrate of 300 mL/min per ADEP guidance was utilized in the model, despite the majority of the 26 Sites conducting low-flow purging at a flowrate between 200 and 250 mL/min. Lower flowrates increase the empty bed contact time (EBCT), which increases the time to breakthrough. Sufficient EBCT ensures adequate contact time between the influent untreated groundwater and the GAC media. The maximum low-flow methodology flowrate was selected to provide the most conservative breakthrough estimates.

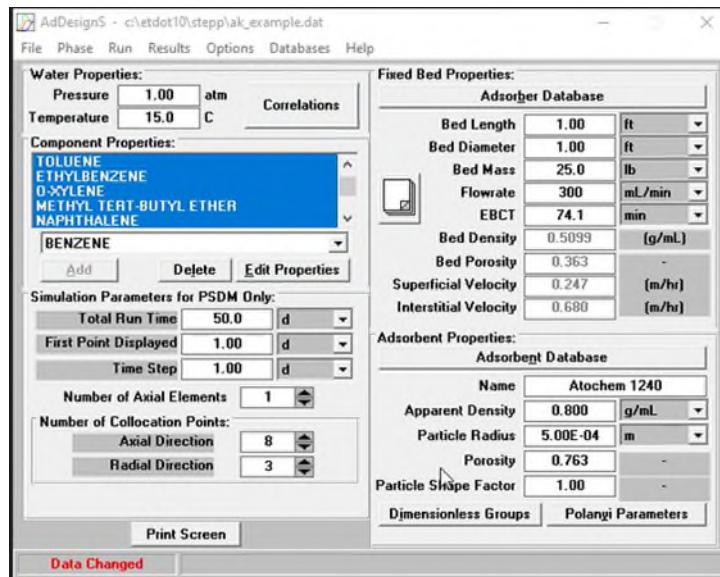
### AdDesignS™ Modeling Parameters

- Freundlich constants ( $1/n$  and  $K$ ) used for individual isotherm models were selected in the software's database using the maximum concentration range observed for that constituent across the 26 Sites.
- Models based on a single 5-gallon capacity portable GAC vessel (Figure A-1, Attachment B).
- To account for potential fouling of the GAC media, default groundwater conditions were selected.
- Pore-surface diffusion model (PSDM) was utilized.
- Adsorbent properties were selected using the software's database similar to the GAC media commonly used across the 26 sites (virgin GAC produced from coconut shell char).

## Attachment A Chevron Environmental Management Company – Alaska Portfolio Isotherm Model

### AdDesignS™ Models

Figure A-1 PSDM Parameters, Fixed Bed and Adsorbent Properties



### Alaska Chevron Site Modeling Results

- As the most conservative parameters were input in the model, virtually no breakthrough of any COPCs was observed until over 3 weeks of continuous run time. As expected, MTBE was the first constituent to show breakthrough, at approximately 24 days. While MTBE was included in the model due to its effect on carbon adsorption, it was only detected at 4 of the 26 Sites in which GAC is utilized.
- Arcadis' typical purge volume for all 26 Sites for one event is approximately 225 gallons of purge water, which at 300 mL/min is approximately 2.37 days. Even with the most conservative parameters applied to the model, this equates to more than 2,000 gallons of purge water at 300 mL/min before any breakthrough is observed and more than 2,300 gallons of purge water at 300 mL/min before 5% of the influent concentration of MTBE is observed breaking through.
- Additional breakthrough estimates are detailed in Figures A-2 through A-9.

Figure A-2 Benzene Breakthrough

Results for the PSDM (No Reactions Present)				
Results for:				
BENZENE	Length of the MTZ (cm):	8.935		
5% of influent conc.	43.85	8.52E+02	1.67	2.50
50% of influent conc.	46.24	8.98E+02	1.76	25.00
95% of influent conc.	57.41	1.12E+03	2.19	47.50
Treatment Objective	43.85	8.52E+02	1.67	2.50

## Attachment A Chevron Environmental Management Company – Alaska Portfolio Isotherm Model

Figure A-3 Toluene Breakthrough

Results for the PSDM (No Reactions Present)

Results for:

TOLUENE	Length of the MTZ (cm):	1.761	<input type="button" value="Close"/>	
	Time (days)	BVT(m³/m³)	VTM(m³/kg)	C (mg/L)
5% of influent conc.	42.28	8.21E+02	1.61	2.50
50% of influent conc.	44.00	8.55E+02	1.68	25.00
95% of influent conc.	44.82	8.71E+02	1.71	47.50
Treatment Objective	42.28	8.21E+02	1.61	2.50

Figure A-4 Ethylbenzene Breakthrough

Results for the PSDM (No Reactions Present)

Results for:

ETHYLBENZENE	Length of the MTZ (cm):	N/A	<input type="button" value="Close"/>	
	Time (days)	BVT(m³/m³)	VTM(m³/kg)	C (mg/L)
5% of influent conc.	1.70E+02	3.31E+03	6.48	1.00
50% of influent conc.	3.57E+02	6.93E+03	13.60	10.00
95% of influent conc.	N/A	N/A	N/A	N/A
Treatment Objective	1.70E+02	3.31E+03	6.48	1.00

Figure A-5 o-Xylene Breakthrough

Results for the PSDM (No Reactions Present)

Results for:

OXYLENE	Length of the MTZ (cm):	N/A	<input type="button" value="Close"/>	
	Time (days)	BVT(m³/m³)	VTM(m³/kg)	C (mg/L)
5% of influent conc.	N/A	N/A	N/A	N/A
50% of influent conc.	N/A	N/A	N/A	N/A
95% of influent conc.	N/A	N/A	N/A	N/A
Treatment Objective	N/A	N/A	N/A	N/A

Figure A-6 Naphthalene Breakthrough

Results for the PSDM (No Reactions Present)

Results for:

NAPHTHALENE	Length of the MTZ (cm):	N/A	<input type="button" value="Close"/>	
	Time (days)	BVT(m³/m³)	VTM(m³/kg)	C (mg/L)
5% of influent conc.	2.09E+02	4.05E+03	7.95	1.50E-02
50% of influent conc.	N/A	N/A	N/A	N/A
95% of influent conc.	N/A	N/A	N/A	N/A
Treatment Objective	2.09E+02	4.05E+03	7.95	1.50E-02

Figure A-7 MTBE Breakthrough

Results for the PSDM (No Reactions Present)

Results for:

METHYL TERT-BUTYL ETHER	Length of the MTZ (cm):	4.737	<input type="button" value="Close"/>	
	Time (days)	BVT(m³/m³)	VTM(m³/kg)	C (mg/L)
5% of influent conc.	24.56	4.77E+02	0.94	0.15
50% of influent conc.	28.49	5.53E+02	1.09	1.50
95% of influent conc.	28.99	5.63E+02	1.10	2.85
Treatment Objective	24.56	4.77E+02	0.94	0.15

## Attachment A Chevron Environmental Management Company – Alaska Portfolio Isotherm Model

Figure A-8 COPC Breakthrough (Annual)

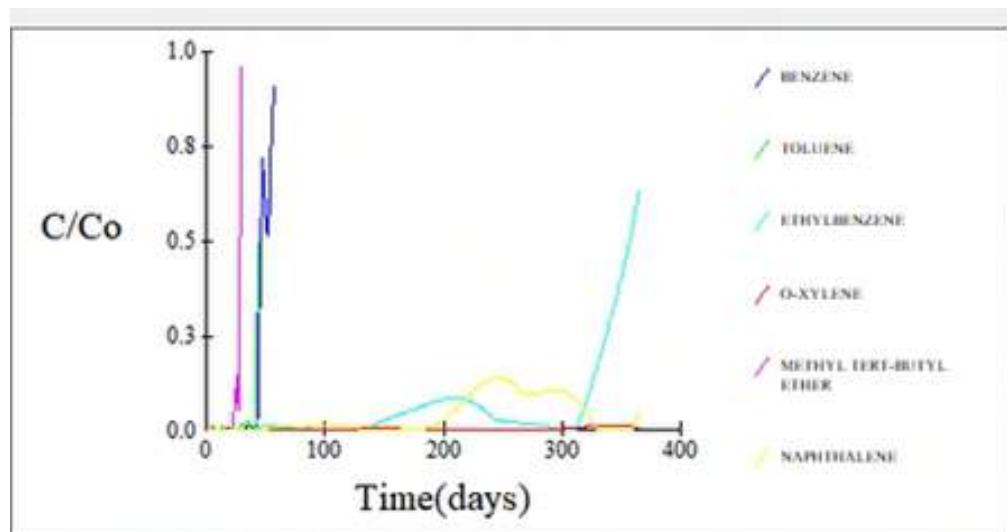
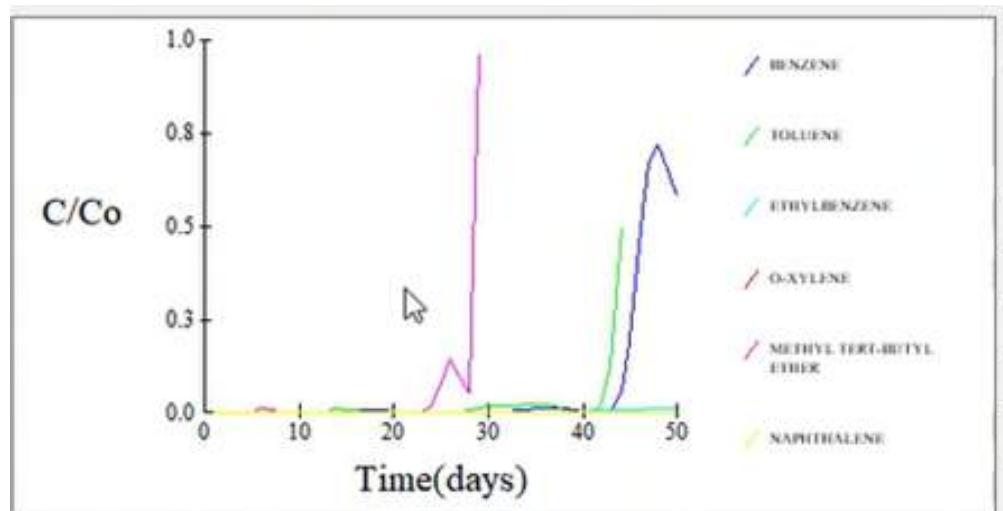


Figure A-9 COPC Breakthrough (Monthly)



### Alaska Chevron Site Proposed Path Forward

Based on the models above, the most conservative breakthrough estimates are much longer than the period of routine GAC utilization activities conducted across the 26 Sites. However, Arcadis plans to replace the GAC media in each portable vessel after each round of sampling to ensure fresh carbon is utilized for each reporting period to prevent any fouling from occurring between events. If site conditions change such as significant increases to observed COPC concentrations, the isotherm model will be re-run to determine new breakthrough estimates.

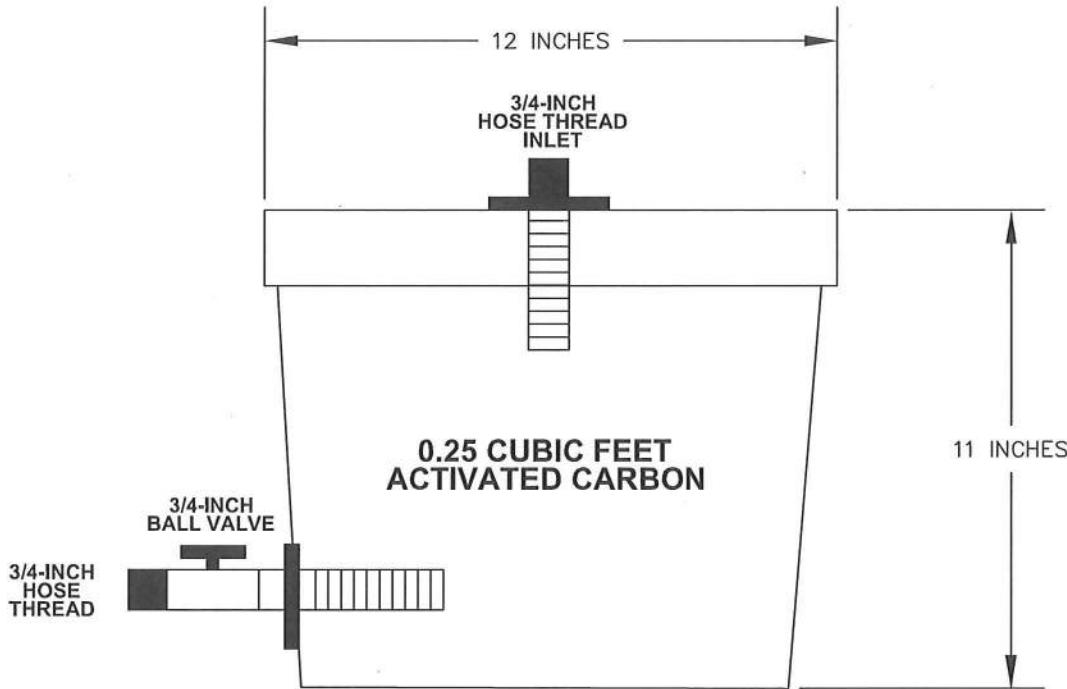
# **Attachment B**

## **GAC Tracking Log**

# Granular Activated Carbon (GAC) Volume Tracking Log

# **Attachment C**

**Portable GAC Vessel Diagram**



**PORTABLE LIQUID-PHASE GRANULAR ACTIVATED CARBON (LGAC) VESSEL**

ATLANTIC RICHFIELD COMPANY  
FORMER AND CURRENT ARCO STATION  
WASHINGTON STATE

**TYPICAL PORTABLE LGAC VESSEL**

 **ARCADIS**

FIGURE  
**1**

Arcadis U.S., Inc.  
630 Plaza Drive, Suite 200  
Highlands Ranch  
Colorado 80129  
Phone: 720 344 3500  
Fax: 720 344 3535  
[www.arcadis.com](http://www.arcadis.com)

## APPENDIX D





# ANALYTICAL REPORT

April 15, 2022

<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

## Arcadis - Chevron - AK

Sample Delivery Group: L1479046  
Samples Received: 04/06/2022  
Project Number: 30063667.19.21  
Description: 97324  
Site: 4417 LAKE OTIS PKWY, ANCHORAGE  
Report To: Sydney Clark/Erika Midkiff  
880 H St.  
Anchorage, AK 99501

Entire Report Reviewed By:

Jordan N Zito  
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 Analytical National is performed per guidance provided in laboratory standard operating procedures ENV-SOP-MTJL-0067 and ENV-SOP-MTJL-0068. Where sampling conducted by the customer, results relate to the accuracy of the information provided, and as the samples are received.

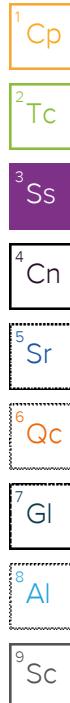
Pace Analytical National

12065 Lebanon Rd Mount Juliet, TN 37122 615-758-5858 800-767-5859 [www.pacenational.com](http://www.pacenational.com)

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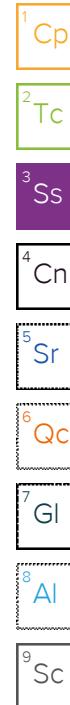
# SAMPLE SUMMARY



Sample ID: MW-2R-W-20220404 L1479046-01 GW			Collected by E Wujcik	Collected date/time 04/04/22 09:00	Received date/time 04/06/22 09:30
Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Volatile Organic Compounds (GC) by Method AK101	WG1846884	1	04/13/22 19:44	04/13/22 19:44	JHH
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844497	25	04/07/22 15:06	04/07/22 15:06	BRA
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844554	1	04/07/22 11:51	04/07/22 11:51	BMB
Semi-Volatile Organic Compounds (GC) by Method AK102	WG1847788	1	04/13/22 14:26	04/13/22 20:33	DMG
Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM	WG1844094	1.25	04/07/22 04:03	04/08/22 01:32	AGW
Sample ID: MW-8RR-W-20220404 L1479046-02 GW			Collected by E Wujcik	Collected date/time 04/04/22 10:00	Received date/time 04/06/22 09:30
Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Volatile Organic Compounds (GC) by Method AK101	WG1846884	1	04/13/22 20:10	04/13/22 20:10	JHH
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844497	1	04/07/22 14:42	04/07/22 14:42	BRA
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844554	1	04/07/22 12:10	04/07/22 12:10	BMB
Semi-Volatile Organic Compounds (GC) by Method AK102	WG1847788	1	04/13/22 14:26	04/13/22 20:54	DMG
Sample ID: MW-1R-W-20220404 L1479046-03 GW			Collected by E Wujcik	Collected date/time 04/04/22 11:00	Received date/time 04/06/22 09:30
Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Volatile Organic Compounds (GC) by Method AK101	WG1846884	1	04/13/22 20:37	04/13/22 20:37	JHH
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844497	1	04/07/22 14:18	04/07/22 14:18	BRA
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844554	1	04/07/22 12:29	04/07/22 12:29	BMB
Semi-Volatile Organic Compounds (GC) by Method AK102	WG1847788	1	04/13/22 14:26	04/13/22 21:14	DMG
Sample ID: MW-9-W-20220404 L1479046-04 GW			Collected by E Wujcik	Collected date/time 04/04/22 12:00	Received date/time 04/06/22 09:30
Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Volatile Organic Compounds (GC) by Method AK101	WG1846884	1	04/13/22 21:03	04/13/22 21:03	JHH
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844497	25	04/07/22 15:29	04/07/22 15:29	BRA
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844554	1	04/07/22 12:48	04/07/22 12:48	BMB
Semi-Volatile Organic Compounds (GC) by Method AK102	WG1847788	1	04/13/22 14:26	04/13/22 22:14	DMG
Sample ID: BD-1-W-20220404 L1479046-05 GW			Collected by E Wujcik	Collected date/time 04/04/22 00:00	Received date/time 04/06/22 09:30
Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Volatile Organic Compounds (GC) by Method AK101	WG1846884	1	04/13/22 21:30	04/13/22 21:30	JHH
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844497	25	04/07/22 15:53	04/07/22 15:53	BRA
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844554	1	04/07/22 13:07	04/07/22 13:07	BMB
Semi-Volatile Organic Compounds (GC) by Method AK102	WG1847788	1	04/13/22 14:26	04/13/22 22:35	DMG
Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM	WG1844094	1.25	04/07/22 04:03	04/08/22 01:52	SHG
Sample ID: EQB-1-W-20220404 L1479046-06 GW			Collected by E Wujcik	Collected date/time 04/04/22 13:00	Received date/time 04/06/22 09:30
Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Volatile Organic Compounds (GC) by Method AK101	WG1846884	1	04/13/22 18:51	04/13/22 18:51	JHH
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844497	1	04/07/22 13:54	04/07/22 13:54	BRA
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844554	1	04/07/22 10:34	04/07/22 10:34	BMB
Semi-Volatile Organic Compounds (GC) by Method AK102	WG1847788	1	04/13/22 14:26	04/13/22 22:55	DMG

# SAMPLE SUMMARY

EQB-1-W-20220404 L1479046-06 GW			Collected by E Wujcik	Collected date/time 04/04/22 13:00	Received date/time 04/06/22 09:30	
Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM	WG1844094	1.25	04/07/22 04:03	04/08/22 00:52	AGW	Mt. Juliet, TN
TRIP BLANK-W-20220404 L1479046-07 GW			Collected by E Wujcik	Collected date/time 04/04/22 00:00	Received date/time 04/06/22 09:30	
Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Volatile Organic Compounds (GC) by Method AK101	WG1846884	1	04/13/22 17:58	04/13/22 17:58	JHH	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844497	1	04/07/22 13:30	04/07/22 13:30	BRA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1844554	1	04/07/22 10:15	04/07/22 10:15	BMB	Mt. Juliet, TN



# CASE NARRATIVE

Unless qualified or noted within the narrative below, all sample aliquots were received at the correct temperature, in the proper containers, with the appropriate preservatives, and within method specified holding times. 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.



Jordan N Zito  
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> GI
- <sup>8</sup> AL
- <sup>9</sup> SC

## Sample Delivery Group (SDG) Narrative

Analyzed from headspace vial.

Batch	Method	Lab Sample ID
WG1846884	AK101	L1479046-07

## Volatile Organic Compounds (GC) by Method AK101

The sample matrix interfered with the ability to make any accurate determination; spike value is low.

Batch	Lab Sample ID	Analytes
WG1846884	(MS) R3781213-3, (MS) R3781213-5, (MSD) R3781213-6, (MSD) R3781213-4, L1479046-03	TPHGAK C6 to C10

The associated batch QC was outside the established quality control range for precision.

Batch	Lab Sample ID	Analytes
WG1846884	(MSD) R3781213-4, (MSD) R3781213-6, L1479046-03	TPHGAK C6 to C10

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

The reported concentration is an estimate. The continuing calibration standard associated with this data responded low. Method sensitivity check is acceptable.

Batch	Lab Sample ID	Analytes
WG1844554	L1479046-01	Acrolein
WG1844554	L1479046-02	Acrolein
WG1844554	L1479046-03	Acrolein
WG1844554	L1479046-04	Acrolein
WG1844554	L1479046-05	Acrolein
WG1844554	L1479046-06	Acrolein
WG1844554	L1479046-07	Acrolein

The associated batch QC was above the established quality control range for accuracy.

Batch	Lab Sample ID	Analytes
WG1844554	(LCSD) R3778810-2, L1479046-01, 02, 03, 04, 05, 06, 07	1,2-Dibromo-3-Chloropropane, Acetone and Methyl tert-butyl ether

# CASE NARRATIVE

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

The associated batch QC was outside the established quality control range for precision.

Batch	Lab Sample ID	Analytes
WG1844554	(LCSD) R3778810-2, L1479046-01, 02, 04, 05, 06, 07	27 analytes

The sample matrix interfered with the ability to make any accurate determination; spike value is high.

Batch	Lab Sample ID	Analytes
WG1844554	(MS) R3778810-4, (MSD) R3778810-5, L1479046-03	1,2-Dibromo-3-Chloropropane, 2-Butanone (MEK), 4-Methyl-2-pentanone (MIBK), Acetone and Methyl tert-butyl ether

- <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> GI
- <sup>8</sup> Al
- <sup>9</sup> Sc



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

Analyte	Result ug/l	<u>Qualifier</u>	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	<u>Batch</u>	1 Cp
Naphthalene	31.0	J3	1.00	5.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	2 Tc
n-Propylbenzene	81.0	J3	0.0993	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	3 Ss
Styrene	U		0.118	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	4 Cn
1,1,2-Tetrachloroethane	U		0.147	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	5 Sr
1,1,2,2-Tetrachloroethane	U	J3	0.133	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	6 Qc
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	7 Gl
Tetrachloroethylene	U		0.300	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	8 Al
Toluene	1.32		0.278	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	9 Sc
1,2,3-Trichlorobenzene	U	J3	0.230	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
1,2,4-Trichlorobenzene	U	J3	0.481	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
1,1,1-Trichloroethane	U		0.149	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
1,1,2-Trichloroethane	U		0.158	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
Trichloroethylene	U		0.190	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
Trichlorofluoromethane	U		0.160	5.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
1,2,4-Trimethylbenzene	17.8	J3	0.322	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
1,2,3-Trimethylbenzene	2.94	J3	0.104	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
1,3,5-Trimethylbenzene	1.08	J3	0.104	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
Vinyl chloride	U		0.234	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
Xylenes, Total	22.7		0.174	3.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
o-Xylene	1.76		0.174	1.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
m&p-Xylene	20.9		0.430	2.00	1	04/07/2022 11:51	<a href="#">WG1844554</a>	
(S) Toluene-d8	96.3			80.0-120		04/07/2022 11:51	<a href="#">WG1844554</a>	
(S) 4-Bromofluorobenzene	96.4			77.0-126		04/07/2022 11:51	<a href="#">WG1844554</a>	
(S) 1,2-Dichloroethane-d4	108			70.0-130		04/07/2022 11:51	<a href="#">WG1844554</a>	

## Sample Narrative:

L1479046-01 WG1844497: Non-target compounds too high to run at a lower dilution.

## Semi-Volatile Organic Compounds (GC) by Method AK102

Analyte	Result ug/l	<u>Qualifier</u>	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	<u>Batch</u>
AK102 DRO C10-C25	1610		229	800	1	04/13/2022 20:33	<a href="#">WG1847788</a>
(S) o-Terphenyl	71.2			50.0-150		04/13/2022 20:33	<a href="#">WG1847788</a>

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

Analyte	Result ug/l	<u>Qualifier</u>	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	<u>Batch</u>
Anthracene	U		0.0238	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Acenaphthene	0.0560	J	0.0238	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Acenaphthylene	U		0.0213	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Benzo(a)anthracene	U		0.0250	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Benzo(a)pyrene	U		0.0225	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Benzo(b)fluoranthene	U		0.0213	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Benzo(g,h,i)perylene	U		0.0225	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Benzo(k)fluoranthene	U		0.0250	0.313	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Chrysene	U		0.0225	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Dibenz(a,h)anthracene	U		0.0225	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Fluoranthene	0.0223	J	0.0137	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Fluorene	U		0.0213	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Indeno(1,2,3-cd)pyrene	U		0.0225	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Naphthalene	19.1		0.160	0.625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Phenanthrene	U		0.0225	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
Pyrene	U		0.0213	0.0625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
1-Methylnaphthalene	8.11		0.0250	0.625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
2-Methylnaphthalene	2.04		0.0350	0.625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>

MW-2R-W-20220404

Collected date/time: 04/04/22 09:00

## SAMPLE RESULTS - 01

L1479046

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

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
2-Chloronaphthalene	U		0.0150	0.625	1.25	04/08/2022 01:32	<a href="#">WG1844094</a>
(S) Nitrobenzene-d5	79.6			11.0-135		04/08/2022 01:32	<a href="#">WG1844094</a>
(S) 2-Fluorobiphenyl	85.2			32.0-120		04/08/2022 01:32	<a href="#">WG1844094</a>
(S) p-Terphenyl-d14	86.8			23.0-122		04/08/2022 01:32	<a href="#">WG1844094</a>

## Sample Narrative:

L1479046-01 WG1844094: Dilution due to sample volume.

<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

## Volatile Organic Compounds (GC) by Method AK101

Analyte	Result ug/l	<u>Qualifier</u>	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	<u>Batch</u>
TPHGAK C6 to C10	U		28.7	100	1	04/13/2022 20:10	<a href="#">WG1846884</a>
(S) a,a,a-Trifluorotoluene(FID)	94.8			50.0-150		04/13/2022 20:10	<a href="#">WG1846884</a>

<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> GI<sup>8</sup> Al<sup>9</sup> Sc

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

Analyte	Result ug/l	<u>Qualifier</u>	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	<u>Batch</u>
1,2,3-Trichloropropane	U		0.00200	0.00500	1	04/07/2022 14:42	<a href="#">WG1844497</a>
Acetone	U	<u>J3 J4</u>	11.3	50.0	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,2-Dibromoethane	0.0120		0.00410	0.00500	1	04/07/2022 14:42	<a href="#">WG1844497</a>
Acrolein	U	<u>C3 J3</u>	2.54	50.0	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Acrylonitrile	U	<u>J3</u>	0.671	10.0	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Benzene	U		0.0941	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Bromobenzene	U	<u>J3</u>	0.118	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Bromochloromethane	U		0.128	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Bromodichloromethane	U		0.136	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Bromoform	U	<u>J3</u>	0.129	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Bromomethane	U		0.605	5.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
n-Butylbenzene	U	<u>J3</u>	0.157	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
sec-Butylbenzene	U	<u>J3</u>	0.125	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
tert-Butylbenzene	U	<u>J3</u>	0.127	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Carbon disulfide	U		0.0962	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Carbon tetrachloride	U		0.128	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Chlorobenzene	U		0.116	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Chlorodibromomethane	U		0.140	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Chloroethane	U		0.192	5.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Chloroform	U		0.111	5.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Chloromethane	U		0.960	2.50	1	04/07/2022 12:10	<a href="#">WG1844554</a>
2-Chlorotoluene	U	<u>J3</u>	0.106	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
4-Chlorotoluene	U	<u>J3</u>	0.114	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,2-Dibromo-3-Chloropropane	U	<u>J3 J4</u>	0.276	5.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Dibromomethane	U		0.122	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,2-Dichlorobenzene	U	<u>J3</u>	0.107	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,3-Dichlorobenzene	U	<u>J3</u>	0.110	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,4-Dichlorobenzene	U	<u>J3</u>	0.120	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Dichlorodifluoromethane	U		0.374	5.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,1-Dichloroethane	U		0.100	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,2-Dichloroethane	0.622	<u>J</u>	0.0819	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,1-Dichloroethene	U		0.188	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
cis-1,2-Dichloroethene	U		0.126	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
trans-1,2-Dichloroethene	U		0.149	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,2-Dichloropropane	U		0.149	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,1-Dichloropropene	U		0.142	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,3-Dichloropropene	U		0.110	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
cis-1,3-Dichloropropene	U		0.111	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
trans-1,3-Dichloropropene	U		0.118	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
2,2-Dichloropropane	U		0.161	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Di-isopropyl ether	U		0.105	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Ethylbenzene	U		0.137	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Hexachloro-1,3-butadiene	U	<u>J3</u>	0.337	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Isopropylbenzene	U		0.105	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
p-Isopropyltoluene	U	<u>J3</u>	0.120	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
2-Butanone (MEK)	U	<u>J3</u>	1.19	10.0	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Methylene Chloride	U		0.430	5.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
4-Methyl-2-pentanone (MIBK)	U	<u>J3</u>	0.478	10.0	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Methyl tert-butyl ether	U	<u>J3 J4</u>	0.101	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>

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

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
Naphthalene	U	J3	1.00	5.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
n-Propylbenzene	0.124	J J3	0.0993	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Styrene	U		0.118	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,1,2-Tetrachloroethane	U		0.147	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,1,2,2-Tetrachloroethane	U	J3	0.133	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Tetrachloroethylene	1.83		0.300	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Toluene	U		0.278	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,2,3-Trichlorobenzene	U	J3	0.230	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,2,4-Trichlorobenzene	U	J3	0.481	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,1,1-Trichloroethane	U		0.149	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,1,2-Trichloroethane	U		0.158	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Trichloroethylene	U		0.190	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Trichlorofluoromethane	U		0.160	5.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,2,4-Trimethylbenzene	U	J3	0.322	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,2,3-Trimethylbenzene	U	J3	0.104	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
1,3,5-Trimethylbenzene	U	J3	0.104	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Vinyl chloride	U		0.234	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
Xylenes, Total	U		0.174	3.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
o-Xylene	U		0.174	1.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
m&p-Xylene	U		0.430	2.00	1	04/07/2022 12:10	<a href="#">WG1844554</a>
(S) Toluene-d8	105			80.0-120		04/07/2022 12:10	<a href="#">WG1844554</a>
(S) 4-Bromofluorobenzene	110			77.0-126		04/07/2022 12:10	<a href="#">WG1844554</a>
(S) 1,2-Dichloroethane-d4	109			70.0-130		04/07/2022 12:10	<a href="#">WG1844554</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 AK102

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
AK102 DRO C10-C25	U		229	800	1	04/13/2022 20:54	<a href="#">WG1847788</a>
(S) o-Terphenyl	56.2			50.0-150		04/13/2022 20:54	<a href="#">WG1847788</a>

## Volatile Organic Compounds (GC) by Method AK101

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
TPHGAK C6 to C10	U	J3 J6	28.7	100	1	04/13/2022 20:37	<a href="#">WG1846884</a>
(S) a,a,a-Trifluorotoluene(FID)	94.0			50.0-150		04/13/2022 20:37	<a href="#">WG1846884</a>

<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

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

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
1,2,3-Trichloropropane	U		0.00200	0.00500	1	04/07/2022 14:18	<a href="#">WG1844497</a>
Acetone	U	J4 J5	11.3	50.0	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,2-Dibromoethane	U		0.00410	0.00500	1	04/07/2022 14:18	<a href="#">WG1844497</a>
Acrolein	U	C3	2.54	50.0	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Acrylonitrile	U		0.671	10.0	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Benzene	U		0.0941	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Bromobenzene	U		0.118	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Bromochloromethane	U		0.128	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Bromodichloromethane	U		0.136	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Bromoform	U		0.129	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Bromomethane	U		0.605	5.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
n-Butylbenzene	U		0.157	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
sec-Butylbenzene	U		0.125	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
tert-Butylbenzene	U		0.127	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Carbon disulfide	U		0.0962	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Carbon tetrachloride	U		0.128	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Chlorobenzene	U		0.116	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Chlorodibromomethane	U		0.140	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Chloroethane	U		0.192	5.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Chloroform	U		0.111	5.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Chloromethane	U		0.960	2.50	1	04/07/2022 12:29	<a href="#">WG1844554</a>
2-Chlorotoluene	U		0.106	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
4-Chlorotoluene	U		0.114	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,2-Dibromo-3-Chloropropane	U	J4 J5	0.276	5.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Dibromomethane	U		0.122	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,2-Dichlorobenzene	U		0.107	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,3-Dichlorobenzene	U		0.110	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,4-Dichlorobenzene	U		0.120	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Dichlorodifluoromethane	U		0.374	5.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,1-Dichloroethane	U		0.100	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,2-Dichloroethane	1.91		0.0819	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,1-Dichloroethylene	U		0.188	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
cis-1,2-Dichloroethene	U		0.126	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
trans-1,2-Dichloroethene	U		0.149	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,2-Dichloropropane	U		0.149	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,1-Dichloropropene	U		0.142	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,3-Dichloropropane	U		0.110	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
cis-1,3-Dichloropropene	U		0.111	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
trans-1,3-Dichloropropene	U		0.118	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
2,2-Dichloropropane	U		0.161	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Di-isopropyl ether	U		0.105	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Ethylbenzene	U		0.137	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Hexachloro-1,3-butadiene	U		0.337	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Isopropylbenzene	U		0.105	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
p-Isopropyltoluene	U		0.120	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
2-Butanone (MEK)	U	J5	1.19	10.0	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Methylene Chloride	U		0.430	5.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
4-Methyl-2-pentanone (MIBK)	U	J5	0.478	10.0	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Methyl tert-butyl ether	U	J4 J5	0.101	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>

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

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
Naphthalene	U		1.00	5.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
n-Propylbenzene	U		0.0993	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Styrene	U		0.118	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Tetrachloroethylene	0.905	J	0.300	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Toluene	U		0.278	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,2,3-Trichlorobenzene	U		0.230	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,2,4-Trichlorobenzene	U		0.481	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,1,1-Trichloroethane	U		0.149	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,1,2-Trichloroethane	U		0.158	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Trichloroethylene	U		0.190	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Trichlorofluoromethane	U		0.160	5.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,2,4-Trimethylbenzene	U		0.322	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,2,3-Trimethylbenzene	U		0.104	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
1,3,5-Trimethylbenzene	U		0.104	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Vinyl chloride	U		0.234	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
Xylenes, Total	U		0.174	3.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
o-Xylene	U		0.174	1.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
m&p-Xylene	U		0.430	2.00	1	04/07/2022 12:29	<a href="#">WG1844554</a>
(S) Toluene-d8	104			80.0-120		04/07/2022 12:29	<a href="#">WG1844554</a>
(S) 4-Bromofluorobenzene	109			77.0-126		04/07/2022 12:29	<a href="#">WG1844554</a>
(S) 1,2-Dichloroethane-d4	111			70.0-130		04/07/2022 12:29	<a href="#">WG1844554</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 AK102

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
AK102 DRO C10-C25	U		229	800	1	04/13/2022 21:14	<a href="#">WG1847788</a>
(S) o-Terphenyl	60.9			50.0-150		04/13/2022 21:14	<a href="#">WG1847788</a>

## Volatile Organic Compounds (GC) by Method AK101

Analyte	Result ug/l	<u>Qualifier</u>	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	<u>Batch</u>
TPHGAK C6 to C10 <i>(S)</i> a,a,a-Trifluorotoluene(FID)	44.5 95.3	J <i>J</i>	28.7 50.0-150	100 1		04/13/2022 21:03 04/13/2022 21:03	<a href="#">WG1846884</a> <a href="#">WG1846884</a>

<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

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

Analyte	Result ug/l	<u>Qualifier</u>	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	<u>Batch</u>
1,2,3-Trichloropropane	U		0.0500	0.125	25	04/07/2022 15:29	<a href="#">WG1844497</a>
Acetone	U	<u>J3 J4</u>	11.3	50.0	1	04/07/2022 12:48	<a href="#">WG1844554</a>
1,2-Dibromoethane	U		0.103	0.125	25	04/07/2022 15:29	<a href="#">WG1844497</a>
Acrolein	U	<u>C3 J3</u>	2.54	50.0	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Acrylonitrile	U	<u>J3</u>	0.671	10.0	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Benzene	U		0.0941	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Bromobenzene	U	<u>J3</u>	0.118	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Bromochloromethane	U		0.128	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Bromodichloromethane	U		0.136	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Bromoform	U	<u>J3</u>	0.129	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Bromomethane	U		0.605	5.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
n-Butylbenzene	U	<u>J3</u>	0.157	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
sec-Butylbenzene	U	<u>J3</u>	0.125	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
tert-Butylbenzene	U	<u>J3</u>	0.127	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Carbon disulfide	U		0.0962	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Carbon tetrachloride	U		0.128	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Chlorobenzene	U		0.116	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Chlorodibromomethane	U		0.140	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Chloroethane	U		0.192	5.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Chloroform	U		0.111	5.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Chloromethane	U		0.960	2.50	1	04/07/2022 12:48	<a href="#">WG1844554</a>
2-Chlorotoluene	U	<u>J3</u>	0.106	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
4-Chlorotoluene	U	<u>J3</u>	0.114	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
1,2-Dibromo-3-Chloropropane	U	<u>J3 J4</u>	0.276	5.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Dibromomethane	U		0.122	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
1,2-Dichlorobenzene	U	<u>J3</u>	0.107	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
1,3-Dichlorobenzene	U	<u>J3</u>	0.110	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
1,4-Dichlorobenzene	U	<u>J3</u>	0.120	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Dichlorodifluoromethane	U		0.374	5.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
1,1-Dichloroethane	U		0.100	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
1,2-Dichloroethane	U		0.0819	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
1,1-Dichloroethylene	U		0.188	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
cis-1,2-Dichloroethene	26.3		0.126	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
trans-1,2-Dichloroethene	U		0.149	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
1,2-Dichloropropane	U		0.149	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
1,1-Dichloropropene	U		0.142	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
1,3-Dichloropropane	U		0.110	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
cis-1,3-Dichloropropene	U		0.111	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
trans-1,3-Dichloropropene	U		0.118	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
2,2-Dichloropropane	U		0.161	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Di-isopropyl ether	U		0.105	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Ethylbenzene	U		0.137	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Hexachloro-1,3-butadiene	U	<u>J3</u>	0.337	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Isopropylbenzene	U		0.105	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
p-Isopropyltoluene	U	<u>J3</u>	0.120	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
2-Butanone (MEK)	U	<u>J3</u>	1.19	10.0	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Methylene Chloride	U		0.430	5.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>
4-Methyl-2-pentanone (MIBK)	U	<u>J3</u>	0.478	10.0	1	04/07/2022 12:48	<a href="#">WG1844554</a>
Methyl tert-butyl ether	U	<u>J3 J4</u>	0.101	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>

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

Analyte	Result ug/l	<u>Qualifier</u>	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	<u>Batch</u>	
Naphthalene	U	J3	1.00	5.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	<sup>1</sup> Cp
n-Propylbenzene	U	J3	0.0993	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	<sup>2</sup> Tc
Styrene	U		0.118	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	<sup>3</sup> Ss
1,1,2-Tetrachloroethane	U		0.147	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
1,1,2,2-Tetrachloroethane	U	J3	0.133	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
Tetrachloroethylene	37.3		0.300	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	<sup>4</sup> Cn
Toluene	U		0.278	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	<sup>5</sup> Sr
1,2,3-Trichlorobenzene	U	J3	0.230	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
1,2,4-Trichlorobenzene	U	J3	0.481	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
1,1,1-Trichloroethane	U		0.149	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
1,1,2-Trichloroethane	U		0.158	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
Trichloroethylene	10.1		0.190	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
Trichlorofluoromethane	U		0.160	5.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
1,2,4-Trimethylbenzene	U	J3	0.322	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
1,2,3-Trimethylbenzene	U	J3	0.104	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
1,3,5-Trimethylbenzene	U	J3	0.104	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
Vinyl chloride	U		0.234	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
Xylenes, Total	U		0.174	3.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
o-Xylene	U		0.174	1.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
m&p-Xylene	U		0.430	2.00	1	04/07/2022 12:48	<a href="#">WG1844554</a>	
(S) Toluene-d8	102			80.0-120		04/07/2022 12:48	<a href="#">WG1844554</a>	
(S) 4-Bromofluorobenzene	106			77.0-126		04/07/2022 12:48	<a href="#">WG1844554</a>	
(S) 1,2-Dichloroethane-d4	109			70.0-130		04/07/2022 12:48	<a href="#">WG1844554</a>	

## Sample Narrative:

L1479046-04 WG1844497: Non-target compounds too high to run at a lower dilution.

## Semi-Volatile Organic Compounds (GC) by Method AK102

Analyte	Result ug/l	<u>Qualifier</u>	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	<u>Batch</u>
AK102 DRO C10-C25	U		229	800	1	04/13/2022 22:14	<a href="#">WG1847788</a>
(S) o-Terphenyl	57.9			50.0-150		04/13/2022 22:14	<a href="#">WG1847788</a>





BD-1-W-20220404

Collected date/time: 04/04/22 00:00

## SAMPLE RESULTS - 05

L1479046

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

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch	
2-Chloronaphthalene	U		0.0150	0.625	1.25	04/08/2022 01:52	<a href="#">WG1844094</a>	<sup>1</sup> Cp
(S) Nitrobenzene-d5	73.2			11.0-135		04/08/2022 01:52	<a href="#">WG1844094</a>	<sup>2</sup> Tc
(S) 2-Fluorobiphenyl	82.0			32.0-120		04/08/2022 01:52	<a href="#">WG1844094</a>	<sup>3</sup> Ss
(S) p-Terphenyl-d14	95.2			23.0-122		04/08/2022 01:52	<a href="#">WG1844094</a>	<sup>4</sup> Cn
								<sup>5</sup> Sr
								<sup>6</sup> Qc
								<sup>7</sup> Gl
								<sup>8</sup> Al
								<sup>9</sup> Sc





EQB-1-W-20220404

Collected date/time: 04/04/22 13:00

## SAMPLE RESULTS - 06

L1479046

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

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	<u>Batch</u>
(S) p-Terphenyl-d14	114			23.0-122		04/08/2022 00:52	<a href="#">WG1844094</a>

## Sample Narrative:

L1479046-06 WG1844094: Dilution due to sample volume.

<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

## Volatile Organic Compounds (GC) by Method AK101

Analyte	Result ug/l	<u>Qualifier</u>	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	<u>Batch</u>
TPHGAK C6 to C10	U		28.7	100	1	04/13/2022 17:58	<a href="#">WG1846884</a>
(S) a,a,a-Trifluorotoluene(FID)	91.6			50.0-150		04/13/2022 17:58	<a href="#">WG1846884</a>

<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

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

Analyte	Result ug/l	<u>Qualifier</u>	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	<u>Batch</u>
1,2,3-Trichloropropane	U		0.00200	0.00500	1	04/07/2022 13:30	<a href="#">WG1844497</a>
Acetone	U	<a href="#">J3 J4</a>	11.3	50.0	1	04/07/2022 10:15	<a href="#">WG1844554</a>
1,2-Dibromoethane	U		0.00410	0.00500	1	04/07/2022 13:30	<a href="#">WG1844497</a>
Acrolein	U	<a href="#">C3 J3</a>	2.54	50.0	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Acrylonitrile	U	<a href="#">J3</a>	0.671	10.0	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Benzene	U		0.0941	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Bromobenzene	U	<a href="#">J3</a>	0.118	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Bromochloromethane	U		0.128	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Bromodichloromethane	U		0.136	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Bromoform	U	<a href="#">J3</a>	0.129	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Bromomethane	U		0.605	5.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
n-Butylbenzene	U	<a href="#">J3</a>	0.157	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
sec-Butylbenzene	U	<a href="#">J3</a>	0.125	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
tert-Butylbenzene	U	<a href="#">J3</a>	0.127	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Carbon disulfide	U		0.0962	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Carbon tetrachloride	U		0.128	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Chlorobenzene	U		0.116	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Chlorodibromomethane	U		0.140	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Chloroethane	U		0.192	5.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Chloroform	U		0.111	5.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Chloromethane	U		0.960	2.50	1	04/07/2022 10:15	<a href="#">WG1844554</a>
2-Chlorotoluene	U	<a href="#">J3</a>	0.106	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
4-Chlorotoluene	U	<a href="#">J3</a>	0.114	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
1,2-Dibromo-3-Chloropropane	U	<a href="#">J3 J4</a>	0.276	5.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Dibromomethane	U		0.122	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
1,2-Dichlorobenzene	U	<a href="#">J3</a>	0.107	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
1,3-Dichlorobenzene	U	<a href="#">J3</a>	0.110	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
1,4-Dichlorobenzene	U	<a href="#">J3</a>	0.120	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Dichlorodifluoromethane	U		0.374	5.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
1,1-Dichloroethane	U		0.100	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
1,2-Dichloroethane	U		0.0819	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
1,1-Dichloroethylene	U		0.188	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
cis-1,2-Dichloroethene	U		0.126	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
trans-1,2-Dichloroethene	U		0.149	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
1,2-Dichloropropane	U		0.149	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
1,1-Dichloropropene	U		0.142	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
1,3-Dichloropropene	U		0.110	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
cis-1,3-Dichloropropene	U		0.111	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
trans-1,3-Dichloropropene	U		0.118	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
2,2-Dichloropropane	U		0.161	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Di-isopropyl ether	U		0.105	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Ethylbenzene	U		0.137	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Hexachloro-1,3-butadiene	U	<a href="#">J3</a>	0.337	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Isopropylbenzene	U		0.105	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
p-Isopropyltoluene	U	<a href="#">J3</a>	0.120	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
2-Butanone (MEK)	U	<a href="#">J3</a>	1.19	10.0	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Methylene Chloride	U		0.430	5.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>
4-Methyl-2-pentanone (MIBK)	U	<a href="#">J3</a>	0.478	10.0	1	04/07/2022 10:15	<a href="#">WG1844554</a>
Methyl tert-butyl ether	U	<a href="#">J3 J4</a>	0.101	1.00	1	04/07/2022 10:15	<a href="#">WG1844554</a>

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

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch	
Naphthalene	U	J3	1.00	5.00	1	04/07/2022 10:15	WG1844554	<sup>1</sup> Cp
n-Propylbenzene	U	J3	0.0993	1.00	1	04/07/2022 10:15	WG1844554	<sup>2</sup> Tc
Styrene	U		0.118	1.00	1	04/07/2022 10:15	WG1844554	<sup>3</sup> Ss
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	04/07/2022 10:15	WG1844554	
1,1,2,2-Tetrachloroethane	U	J3	0.133	1.00	1	04/07/2022 10:15	WG1844554	
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	04/07/2022 10:15	WG1844554	
Tetrachloroethylene	U		0.300	1.00	1	04/07/2022 10:15	WG1844554	<sup>4</sup> Cn
Toluene	U		0.278	1.00	1	04/07/2022 10:15	WG1844554	<sup>5</sup> Sr
1,2,3-Trichlorobenzene	U	J3	0.230	1.00	1	04/07/2022 10:15	WG1844554	
1,2,4-Trichlorobenzene	U	J3	0.481	1.00	1	04/07/2022 10:15	WG1844554	
1,1,1-Trichloroethane	U		0.149	1.00	1	04/07/2022 10:15	WG1844554	
1,1,2-Trichloroethane	U		0.158	1.00	1	04/07/2022 10:15	WG1844554	
Trichloroethylene	U		0.190	1.00	1	04/07/2022 10:15	WG1844554	
Trichlorofluoromethane	U		0.160	5.00	1	04/07/2022 10:15	WG1844554	
1,2,4-Trimethylbenzene	U	J3	0.322	1.00	1	04/07/2022 10:15	WG1844554	
1,2,3-Trimethylbenzene	U	J3	0.104	1.00	1	04/07/2022 10:15	WG1844554	
1,3,5-Trimethylbenzene	U	J3	0.104	1.00	1	04/07/2022 10:15	WG1844554	
Vinyl chloride	U		0.234	1.00	1	04/07/2022 10:15	WG1844554	
Xylenes, Total	U		0.174	3.00	1	04/07/2022 10:15	WG1844554	
o-Xylene	U		0.174	1.00	1	04/07/2022 10:15	WG1844554	
m&p-Xylene	U		0.430	2.00	1	04/07/2022 10:15	WG1844554	
(S) Toluene-d8	106			80.0-120		04/07/2022 10:15	WG1844554	
(S) 4-Bromofluorobenzene	109			77.0-126		04/07/2022 10:15	WG1844554	
(S) 1,2-Dichloroethane-d4	110			70.0-130		04/07/2022 10:15	WG1844554	

## QUALITY CONTROL SUMMARY

[L1479046-01,02,03,04,05,06,07](#)

## Method Blank (MB)

(MB) R3781213-2 04/13/22 17:05

Analyte	MB Result ug/l	<u>MB Qualifier</u>	MB MDL ug/l	MB RDL ug/l
TPHGAK C6 to C10	U		28.7	100
(S) a,a,a-Trifluorotoluene(FID)	95.5		60.0-120	

1 Cp

2 Tc

3 Ss

4 Cn

15 Sr

6 Qc

7 Gl

8 Al

9 Sc

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

(LCS) R3781213-1 04/13/22 15:50 • (LCSD) R3781213-7 04/14/22 16:13

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	<u>LCS Qualifier</u>	<u>LCSD Qualifier</u>	RPD %	RPD Limits %
TPHGAK C6 to C10	5000	4740	4370	94.8	87.4	60.0-120			8.12	20
(S) a,a,a-Trifluorotoluene(FID)			106	107	107	60.0-120				

## L1480598-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1480598-01 04/13/22 21:56 • (MS) R3781213-3 04/14/22 14:01 • (MSD) R3781213-4 04/14/22 14:27

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	<u>MS Qualifier</u>	<u>MSD Qualifier</u>	RPD %	RPD Limits %
TPHGAK C6 to C10	5000	U	2350	3050	47.0	61.0	1	70.0-130	J6	J3 J6	25.9	20
(S) a,a,a-Trifluorotoluene(FID)				103	102	102	1	50.0-150				

## L1479046-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1479046-03 04/13/22 20:37 • (MS) R3781213-5 04/14/22 14:54 • (MSD) R3781213-6 04/14/22 15:20

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	<u>MS Qualifier</u>	<u>MSD Qualifier</u>	RPD %	RPD Limits %
TPHGAK C6 to C10	5000	U	3250	2340	65.0	46.8	1	70.0-130	J6	J3 J6	32.6	20
(S) a,a,a-Trifluorotoluene(FID)				104	102	102	1	50.0-150				

WG1844497

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

## QUALITY CONTROL SUMMARY

[L1479046-01,02,03,04,05,06,07](#)

## Method Blank (MB)

(MB) R3778873-2 04/07/22 12:43

Analyte	MB Result ug/l	<u>MB Qualifier</u>	MB MDL ug/l	MB RDL ug/l
1,2,3-Trichloropropane	U		0.00200	0.00500
1,2-Dibromoethane	U		0.00410	0.00500

<sup>1</sup>Cp<sup>2</sup>Tc<sup>3</sup>Ss<sup>4</sup>Cn<sup>15</sup>Sr<sup>6</sup>Qc<sup>7</sup>Gl<sup>8</sup>Al<sup>9</sup>Sc

## Laboratory Control Sample (LCS)

(LCS) R3778873-1 04/07/22 12:19

Analyte	Spike Amount ug/l	LCS Result ug/l	LCS Rec. %	Rec. Limits %	<u>LCS Qualifier</u>
1,2,3-Trichloropropane	0.0500	0.0450	90.0	70.0-130	
1,2-Dibromoethane	0.0500	0.0410	82.0	70.0-130	

## L1479046-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1479046-03 04/07/22 14:18 • (MS) R3778873-3 04/07/22 16:17 • (MSD) R3778873-4 04/07/22 16:41

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	<u>MS Qualifier</u>	<u>MSD Qualifier</u>	RPD	RPD Limits
1,2,3-Trichloropropane	0.0500	U	0.0520	0.0530	104	106	1	70.0-130			1.90	20
1,2-Dibromoethane	0.0500	U	0.0500	0.0520	100	104	1	70.0-130			3.92	20

WG1844554

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

## QUALITY CONTROL SUMMARY

[L1479046-01,02,03,04,05,06,07](#)

## Method Blank (MB)

(MB) R3778810-3 04/07/22 09:30

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l	
Acetone	U		11.3	50.0	<sup>1</sup> Cp
Acrolein	U		2.54	50.0	<sup>2</sup> Tc
Acrylonitrile	U		0.671	10.0	<sup>3</sup> Ss
Benzene	U		0.0941	1.00	<sup>4</sup> Cn
Bromobenzene	U		0.118	1.00	<sup>15</sup> Sr
Bromochloromethane	U		0.128	1.00	<sup>6</sup> Qc
Bromodichloromethane	U		0.136	1.00	<sup>7</sup> Gl
Bromoform	U		0.129	1.00	<sup>8</sup> Al
Bromomethane	U		0.605	5.00	<sup>9</sup> Sc
n-Butylbenzene	U		0.157	1.00	
sec-Butylbenzene	U		0.125	1.00	
tert-Butylbenzene	U		0.127	1.00	
Carbon disulfide	U		0.0962	1.00	
Carbon tetrachloride	U		0.128	1.00	
Chlorobenzene	U		0.116	1.00	
Chlorodibromomethane	U		0.140	1.00	
Chloroethane	U		0.192	5.00	
Chloroform	U		0.111	5.00	
Chloromethane	U		0.960	2.50	
2-Chlorotoluene	U		0.106	1.00	
4-Chlorotoluene	U		0.114	1.00	
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	
Dibromomethane	U		0.122	1.00	
1,2-Dichlorobenzene	U		0.107	1.00	
1,3-Dichlorobenzene	U		0.110	1.00	
1,4-Dichlorobenzene	U		0.120	1.00	
Dichlorodifluoromethane	U		0.374	5.00	
1,1-Dichloroethane	U		0.100	1.00	
1,2-Dichloroethane	U		0.0819	1.00	
1,1-Dichloroethene	U		0.188	1.00	
cis-1,2-Dichloroethene	U		0.126	1.00	
trans-1,2-Dichloroethene	U		0.149	1.00	
1,2-Dichloropropane	U		0.149	1.00	
1,1-Dichloropropene	U		0.142	1.00	
1,3-Dichloropropane	U		0.110	1.00	
cis-1,3-Dichloropropene	U		0.111	1.00	
trans-1,3-Dichloropropene	U		0.118	1.00	
2,2-Dichloropropane	U		0.161	1.00	
Di-isopropyl ether	U		0.105	1.00	
Ethylbenzene	U		0.137	1.00	

## QUALITY CONTROL SUMMARY

[L1479046-01,02,03,04,05,06,07](#)

## Method Blank (MB)

(MB) R3778810-3 04/07/22 09:30

Analyte	MB Result ug/l	<u>MB Qualifier</u>	MB MDL ug/l	MB RDL ug/l	1 Cp
Hexachloro-1,3-butadiene	U		0.337	1.00	
Isopropylbenzene	U		0.105	1.00	
p-Isopropyltoluene	U		0.120	1.00	
2-Butanone (MEK)	U		1.19	10.0	
Methylene Chloride	U		0.430	5.00	
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	
Methyl tert-butyl ether	U		0.101	1.00	
Naphthalene	U		1.00	5.00	
n-Propylbenzene	U		0.0993	1.00	
Styrene	U		0.118	1.00	
1,1,2-Tetrachloroethane	U		0.147	1.00	
1,1,2,2-Tetrachloroethane	U		0.133	1.00	
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	
Tetrachloroethene	U		0.300	1.00	
Toluene	U		0.278	1.00	
1,2,3-Trichlorobenzene	U		0.230	1.00	
1,2,4-Trichlorobenzene	U		0.481	1.00	
1,1,1-Trichloroethane	U		0.149	1.00	
1,1,2-Trichloroethane	U		0.158	1.00	
Trichloroethene	U		0.190	1.00	
Trichlorofluoromethane	U		0.160	5.00	
1,2,4-Trimethylbenzene	U		0.322	1.00	
1,2,3-Trimethylbenzene	U		0.104	1.00	
1,3,5-Trimethylbenzene	U		0.104	1.00	
Vinyl chloride	U		0.234	1.00	
Xylenes, Total	U		0.174	3.00	
o-Xylene	U		0.174	1.00	
m&p-Xylenes	U		0.430	2.00	
(S) Toluene-d8	104		80.0-120		
(S) 4-Bromofluorobenzene	104		77.0-126		
(S) 1,2-Dichloroethane-d4	106		70.0-130		

1 Cp

2 Tc

3 Ss

4 Cn

15 Sr

6 Qc

7 Gl

8 Al

9 Sc

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

(LCS) R3778810-1 04/07/22 07:49 • (LCSD) R3778810-2 04/07/22 08:08

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	<u>LCS Qualifier</u>	<u>LCSD Qualifier</u>	RPD %	RPD Limits %
Acetone	25.0	30.5	45.9	122	184	19.0-160	J3 J4		40.3	27
Acrolein	25.0	12.4	17.5	49.6	70.0	10.0-160	J3		34.1	26

## QUALITY CONTROL SUMMARY

[L1479046-01,02,03,04,05,06,07](#)

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

(LCS) R3778810-1 04/07/22 07:49 • (LCSD) R3778810-2 04/07/22 08:08

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	<u>LCS Qualifier</u>	<u>LCSD Qualifier</u>	RPD %	RPD Limits %
Acrylonitrile	25.0	21.2	29.8	84.8	119	55.0-149	J3		33.7	20
Benzene	5.00	4.48	5.03	89.6	101	70.0-123			11.6	20
Bromobenzene	5.00	4.55	5.62	91.0	112	73.0-121	J3		21.0	20
Bromochloromethane	5.00	4.61	5.60	92.2	112	76.0-122			19.4	20
Bromodichloromethane	5.00	4.45	5.12	89.0	102	75.0-120			14.0	20
Bromoform	5.00	4.30	5.27	86.0	105	68.0-132	J3		20.3	20
Bromomethane	5.00	4.22	4.43	84.4	88.6	10.0-160			4.86	25
n-Butylbenzene	5.00	4.35	5.42	87.0	108	73.0-125	J3		21.9	20
sec-Butylbenzene	5.00	4.53	5.89	90.6	118	75.0-125	J3		26.1	20
tert-Butylbenzene	5.00	4.69	5.81	93.8	116	76.0-124	J3		21.3	20
Carbon disulfide	5.00	4.66	5.32	93.2	106	61.0-128			13.2	20
Carbon tetrachloride	5.00	4.62	5.11	92.4	102	68.0-126			10.1	20
Chlorobenzene	5.00	4.49	5.23	89.8	105	80.0-121			15.2	20
Chlorodibromomethane	5.00	4.43	5.25	88.6	105	77.0-125			16.9	20
Chloroethane	5.00	4.14	4.62	82.8	92.4	47.0-150			11.0	20
Chloroform	5.00	4.34	4.94	86.8	98.8	73.0-120			12.9	20
Chloromethane	5.00	4.28	4.40	85.6	88.0	41.0-142			2.76	20
2-Chlorotoluene	5.00	4.58	5.71	91.6	114	76.0-123	J3		22.0	20
4-Chlorotoluene	5.00	4.45	5.76	89.0	115	75.0-122	J3		25.7	20
1,2-Dibromo-3-Chloropropane	5.00	4.60	6.73	92.0	135	58.0-134	J3 J4		37.6	20
Dibromomethane	5.00	4.72	5.38	94.4	108	80.0-120			13.1	20
1,2-Dichlorobenzene	5.00	4.39	5.71	87.8	114	79.0-121	J3		26.1	20
1,3-Dichlorobenzene	5.00	4.36	5.52	87.2	110	79.0-120	J3		23.5	20
1,4-Dichlorobenzene	5.00	4.20	5.44	84.0	109	79.0-120	J3		25.7	20
Dichlorodifluoromethane	5.00	4.85	4.81	97.0	96.2	51.0-149			0.828	20
1,1-Dichloroethane	5.00	4.38	5.00	87.6	100	70.0-126			13.2	20
1,2-Dichloroethane	5.00	4.57	5.27	91.4	105	70.0-128			14.2	20
1,1-Dichloroethene	5.00	4.60	4.93	92.0	98.6	71.0-124			6.93	20
cis-1,2-Dichloroethene	5.00	4.55	5.21	91.0	104	73.0-120			13.5	20
trans-1,2-Dichloroethene	5.00	4.33	4.89	86.6	97.8	73.0-120			12.1	20
1,2-Dichloropropane	5.00	4.60	5.37	92.0	107	77.0-125			15.4	20
1,1-Dichloropropene	5.00	4.29	4.90	85.8	98.0	74.0-126			13.3	20
1,3-Dichloropropane	5.00	4.37	5.25	87.4	105	80.0-120			18.3	20
cis-1,3-Dichloropropene	5.00	4.55	5.17	91.0	103	80.0-123			12.8	20
trans-1,3-Dichloropropene	5.00	4.33	5.25	86.6	105	78.0-124			19.2	20
2,2-Dichloropropane	5.00	4.88	5.54	97.6	111	58.0-130			12.7	20
Di-isopropyl ether	5.00	4.59	5.38	91.8	108	58.0-138			15.8	20
Ethylbenzene	5.00	4.65	5.42	93.0	108	79.0-123			15.3	20
Hexachloro-1,3-butadiene	5.00	4.02	5.23	80.4	105	54.0-138	J3		26.2	20
Isopropylbenzene	5.00	4.72	5.41	94.4	108	76.0-127			13.6	20

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

## QUALITY CONTROL SUMMARY

[L1479046-01,02,03,04,05,06,07](#)

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

(LCS) R3778810-1 04/07/22 07:49 • (LCSD) R3778810-2 04/07/22 08:08

1 Cp

2 Tc

3 Ss

4 Cn

15 Sr

6 Qc

7 Gl

8 Al

9 Sc

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
p-Isopropyltoluene	5.00	4.56	5.65	91.2	113	76.0-125	J3	J3	21.4	20
2-Butanone (MEK)	25.0	27.1	35.5	108	142	44.0-160	J3	J3	26.8	20
Methylene Chloride	5.00	4.65	5.47	93.0	109	67.0-120			16.2	20
4-Methyl-2-pentanone (MIBK)	25.0	25.2	32.7	101	131	68.0-142	J3	J3	25.9	20
Methyl tert-butyl ether	5.00	5.10	6.42	102	128	68.0-125	J3 J4	J3 J4	22.9	20
Naphthalene	5.00	4.59	6.40	91.8	128	54.0-135	J3	J3	32.9	20
n-Propylbenzene	5.00	4.85	6.01	97.0	120	77.0-124	J3	J3	21.4	20
Styrene	5.00	4.26	4.96	85.2	99.2	73.0-130			15.2	20
1,1,2-Tetrachloroethane	5.00	4.54	5.34	90.8	107	75.0-125			16.2	20
1,1,2,2-Tetrachloroethane	5.00	4.56	6.06	91.2	121	65.0-130	J3	J3	28.2	20
1,1,2-Trichlorotrifluoroethane	5.00	4.25	4.74	85.0	94.8	69.0-132			10.9	20
Tetrachloroethene	5.00	4.40	5.13	88.0	103	72.0-132			15.3	20
Toluene	5.00	4.60	5.20	92.0	104	79.0-120			12.2	20
1,2,3-Trichlorobenzene	5.00	4.23	5.65	84.6	113	50.0-138	J3	J3	28.7	20
1,2,4-Trichlorobenzene	5.00	4.24	5.62	84.8	112	57.0-137	J3	J3	28.0	20
1,1,1-Trichloroethane	5.00	4.65	5.31	93.0	106	73.0-124			13.3	20
1,1,2-Trichloroethane	5.00	4.39	5.18	87.8	104	80.0-120			16.5	20
Trichloroethene	5.00	4.52	5.16	90.4	103	78.0-124			13.2	20
Trichlorofluoromethane	5.00	4.73	5.33	94.6	107	59.0-147			11.9	20
1,2,4-Trimethylbenzene	5.00	4.71	5.90	94.2	118	76.0-121	J3	J3	22.4	20
1,2,3-Trimethylbenzene	5.00	4.48	5.71	89.6	114	77.0-120	J3	J3	24.1	20
1,3,5-Trimethylbenzene	5.00	4.85	5.96	97.0	119	76.0-122	J3	J3	20.5	20
Vinyl chloride	5.00	4.38	4.79	87.6	95.8	67.0-131			8.94	20
Xylenes, Total	15.0	14.3	16.4	95.3	109	79.0-123			13.7	20
o-Xylene	5.00	4.78	5.52	95.6	110	80.0-122			14.4	20
m&p-Xylenes	10.0	9.47	10.9	94.7	109	80.0-122			14.0	20
(S) Toluene-d8				102	103	80.0-120				
(S) 4-Bromofluorobenzene				107	101	77.0-126				
(S) 1,2-Dichloroethane-d4				105	105	70.0-130				

## L1479046-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1479046-03 04/07/22 12:29 • (MS) R3778810-4 04/07/22 18:32 • (MSD) R3778810-5 04/07/22 18:51

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MS Rec. %	MSD Result ug/l	MS Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Acetone	25.0	U	51.1	54.3	204	217	1	10.0-160	J5	J5	6.07	35
Acrolein	25.0	U	28.2	30.0	113	120	1	10.0-160			6.19	39
Acrylonitrile	25.0	U	39.4	40.0	158	160	1	21.0-160			1.51	32
Benzene	5.00	U	6.46	6.80	129	136	1	17.0-158			5.13	27

## QUALITY CONTROL SUMMARY

[L1479046-01,02,03,04,05,06,07](#)

## L1479046-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1479046-03 04/07/22 12:29 • (MS) R3778810-4 04/07/22 18:32 • (MSD) R3778810-5 04/07/22 18:51

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
Bromobenzene	5.00	U	6.67	6.82	133	136	1	30.0-149			2.22	28
Bromochloromethane	5.00	U	6.69	6.96	134	139	1	38.0-142			3.96	26
Bromodichloromethane	5.00	U	6.44	6.71	129	134	1	31.0-150			4.11	27
Bromoform	5.00	U	6.29	6.58	126	132	1	29.0-150			4.51	29
Bromomethane	5.00	U	5.63	5.89	113	118	1	10.0-160			4.51	38
n-Butylbenzene	5.00	U	6.61	6.77	132	135	1	31.0-150			2.39	30
sec-Butylbenzene	5.00	U	7.07	7.26	141	145	1	33.0-155			2.65	29
tert-Butylbenzene	5.00	U	7.06	6.99	141	140	1	34.0-153			0.996	28
Carbon disulfide	5.00	U	6.54	7.08	131	142	1	10.0-156			7.93	28
Carbon tetrachloride	5.00	U	7.14	7.43	143	149	1	23.0-159			3.98	28
Chlorobenzene	5.00	U	6.43	6.58	129	132	1	33.0-152			2.31	27
Chlorodibromomethane	5.00	U	6.09	6.49	122	130	1	37.0-149			6.36	27
Chloroethane	5.00	U	5.99	6.42	120	128	1	10.0-160			6.93	30
Chloroform	5.00	U	6.30	6.64	126	133	1	29.0-154			5.26	28
Chloromethane	5.00	U	5.95	6.17	119	123	1	10.0-160			3.63	29
2-Chlorotoluene	5.00	U	6.82	7.11	136	142	1	32.0-153			4.16	28
4-Chlorotoluene	5.00	U	6.65	6.82	133	136	1	32.0-150			2.52	28
1,2-Dibromo-3-Chloropropane	5.00	U	7.76	8.15	155	163	1	22.0-151	J5	J5	4.90	34
Dibromomethane	5.00	U	6.87	7.00	137	140	1	30.0-151			1.87	27
1,2-Dichlorobenzene	5.00	U	6.43	6.62	129	132	1	34.0-149			2.91	28
1,3-Dichlorobenzene	5.00	U	6.33	6.43	127	129	1	36.0-146			1.57	27
1,4-Dichlorobenzene	5.00	U	6.18	6.45	124	129	1	35.0-142			4.28	27
Dichlorodifluoromethane	5.00	U	6.54	7.15	131	143	1	10.0-160			8.91	29
1,1-Dichloroethane	5.00	U	6.34	6.81	127	136	1	25.0-158			7.15	27
1,2-Dichloroethane	5.00	1.91	8.09	8.50	124	132	1	29.0-151			4.94	27
1,1-Dichloroethene	5.00	U	6.55	7.04	131	141	1	11.0-160			7.21	29
cis-1,2-Dichloroethene	5.00	U	6.32	6.81	126	136	1	10.0-160			7.46	27
trans-1,2-Dichloroethene	5.00	U	6.37	6.73	127	135	1	17.0-153			5.50	27
1,2-Dichloropropane	5.00	U	6.95	7.25	139	145	1	30.0-156			4.23	27
1,1-Dichloropropene	5.00	U	6.32	6.69	126	134	1	25.0-158			5.69	27
1,3-Dichloropropane	5.00	U	6.16	6.59	123	132	1	38.0-147			6.75	27
cis-1,3-Dichloropropene	5.00	U	6.11	6.59	122	132	1	34.0-149			7.56	28
trans-1,3-Dichloropropene	5.00	U	6.17	6.39	123	128	1	32.0-149			3.50	28
2,2-Dichloropropane	5.00	U	7.22	7.54	144	151	1	24.0-152			4.34	29
Di-isopropyl ether	5.00	U	6.83	7.25	137	145	1	21.0-160			5.97	28
Ethylbenzene	5.00	U	6.61	6.71	132	134	1	30.0-155			1.50	27
Hexachloro-1,3-butadiene	5.00	U	6.33	6.59	127	132	1	20.0-154			4.02	34
Isopropylbenzene	5.00	U	6.94	7.31	139	146	1	28.0-157			5.19	27
p-Isopropyltoluene	5.00	U	6.96	7.12	139	142	1	30.0-154			2.27	29
2-Butanone (MEK)	25.0	U	44.9	45.9	180	184	1	10.0-160	J5	J5	2.20	32

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

## QUALITY CONTROL SUMMARY

[L1479046-01,02,03,04,05,06,07](#)

## L1479046-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1479046-03 04/07/22 12:29 • (MS) R3778810-4 04/07/22 18:32 • (MSD) R3778810-5 04/07/22 18:51

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD	RPD Limits
Methylene Chloride	5.00	U	6.48	7.06	130	141	1	23.0-144			8.57	28
4-Methyl-2-pentanone (MIBK)	25.0	U	39.4	41.4	158	166	1	29.0-160	J5	J5	4.95	29
Methyl tert-butyl ether	5.00	U	7.67	8.31	153	166	1	28.0-150	J5	J5	8.01	29
Naphthalene	5.00	U	6.85	7.28	137	146	1	12.0-156			6.09	35
n-Propylbenzene	5.00	U	7.41	7.61	148	152	1	31.0-154			2.66	28
Styrene	5.00	U	5.90	6.26	118	125	1	33.0-155			5.92	28
1,1,1,2-Tetrachloroethane	5.00	U	6.54	6.89	131	138	1	36.0-151			5.21	29
1,1,2,2-Tetrachloroethane	5.00	U	7.22	7.39	144	148	1	33.0-150			2.33	28
1,1,2-Trichlorotrifluoroethane	5.00	U	6.49	6.71	130	134	1	23.0-160			3.33	30
Tetrachloroethylene	5.00	0.905	7.12	7.46	124	131	1	10.0-160			4.66	27
Toluene	5.00	U	6.33	6.62	127	132	1	26.0-154			4.48	28
1,2,3-Trichlorobenzene	5.00	U	6.15	6.39	123	128	1	17.0-150			3.83	36
1,2,4-Trichlorobenzene	5.00	U	6.14	6.68	123	134	1	24.0-150			8.42	33
1,1,1-Trichloroethane	5.00	U	6.97	7.50	139	150	1	23.0-160			7.33	28
1,1,2-Trichloroethane	5.00	U	6.28	6.65	126	133	1	35.0-147			5.72	27
Trichloroethylene	5.00	U	6.28	6.79	126	136	1	10.0-160			7.80	25
Trichlorofluoromethane	5.00	U	3.67	4.75	73.4	95.0	1	17.0-160			25.7	31
1,2,4-Trimethylbenzene	5.00	U	6.80	7.02	136	140	1	26.0-154			3.18	27
1,2,3-Trimethylbenzene	5.00	U	6.63	6.80	133	136	1	32.0-149			2.53	28
1,3,5-Trimethylbenzene	5.00	U	6.99	7.41	140	148	1	28.0-153			5.83	27
Vinyl chloride	5.00	U	6.48	6.68	130	134	1	10.0-160			3.04	27
Xylenes, Total	15.0	U	20.3	21.1	135	141	1	29.0-154			3.86	28
o-Xylene	5.00	U	6.73	7.11	135	142	1	45.0-144			5.49	26
m&p-Xylenes	10.0	U	13.6	14.0	136	140	1	43.0-146			2.90	26
(S) Toluene-d8					102	102		80.0-120				
(S) 4-Bromofluorobenzene					104	106		77.0-126				
(S) 1,2-Dichloroethane-d4					110	111		70.0-130				

1 Cp

2 Tc

3 Ss

4 Cn

15 Sr

6 Qc

7 Gl

8 Al

9 Sc

WG1847788

Semi-Volatile Organic Compounds (GC) by Method AK102

## QUALITY CONTROL SUMMARY

[L1479046-01,02,03,04,05,06](#)

## Method Blank (MB)

(MB) R3780795-1 04/13/22 19:32

Analyte	MB Result ug/l	<u>MB Qualifier</u>	MB MDL ug/l	MB RDL ug/l
AK102 DRO C10-C25	U		229	800
(S) o-Terphenyl	63.0			60.0-120

<sup>1</sup>Cp<sup>2</sup>Tc<sup>3</sup>Ss<sup>4</sup>Cn<sup>15</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) R3780795-2 04/13/22 19:53 • (LCSD) R3780795-3 04/13/22 20:13

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	<u>LCS Qualifier</u>	<u>LCSD Qualifier</u>	RPD %	RPD Limits %
AK102 DRO C10-C25	6000	5630	5510	93.8	91.8	75.0-125			2.15	20
(S) o-Terphenyl				96.1	92.9	60.0-120				

## L1479046-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1479046-03 04/13/22 21:14 • (MS) R3780795-4 04/13/22 21:34 • (MSD) R3780795-5 04/13/22 21:54

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	<u>MS Qualifier</u>	<u>MSD Qualifier</u>	RPD %	RPD Limits %
AK102 DRO C10-C25	6000	U	5590	5750	93.2	95.8	1	75.0-125			2.82	20
(S) o-Terphenyl					94.3	94.8		50.0-150				

## QUALITY CONTROL SUMMARY

L1479046-01,05,06

## Method Blank (MB)

(MB) R3779023-3 04/07/22 17:56

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l	1 Cp
Anthracene	U		0.0190	0.0500	
Acenaphthene	U		0.0190	0.0500	
Acenaphthylene	U		0.0170	0.0500	
Benzo(a)anthracene	U		0.0200	0.0500	
Benzo(a)pyrene	U		0.0180	0.0500	
Benzo(b)fluoranthene	U		0.0170	0.0500	
Benzo(g,h,i)perylene	U		0.0180	0.0500	
Benzo(k)fluoranthene	U		0.0200	0.250	
Chrysene	U		0.0180	0.0500	
Dibenz(a,h)anthracene	U		0.0180	0.0500	
Fluoranthene	U		0.0110	0.0500	
Fluorene	U		0.0170	0.0500	
Indeno(1,2,3-cd)pyrene	U		0.0180	0.0500	
Naphthalene	U		0.128	0.500	
Phenanthrene	U		0.0180	0.0500	
Pyrene	U		0.0170	0.0500	
1-Methylnaphthalene	U		0.0200	0.500	
2-Methylnaphthalene	U		0.0280	0.500	
2-Chloronaphthalene	U		0.0120	0.500	
(S) Nitrobenzene-d5	96.0			11.0-135	
(S) 2-Fluorobiphenyl	91.5			32.0-120	
(S) p-Terphenyl-d14	110			23.0-122	

1 Cp

2 Tc

3 Ss

4 Cn

15 Sr

6 Qc

7 Gl

8 Al

9 Sc

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

(LCS) R3779023-1 04/07/22 17:16 • (LCSD) R3779023-2 04/07/22 17:36

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Anthracene	2.00	1.54	1.59	77.0	79.5	43.0-127			3.19	20
Acenaphthene	2.00	1.61	1.67	80.5	83.5	42.0-120			3.66	20
Acenaphthylene	2.00	1.68	1.77	84.0	88.5	43.0-120			5.22	20
Benzo(a)anthracene	2.00	1.72	1.80	86.0	90.0	46.0-120			4.55	20
Benzo(a)pyrene	2.00	1.46	1.58	73.0	79.0	44.0-122			7.89	20
Benzo(b)fluoranthene	2.00	1.69	1.80	84.5	90.0	43.0-122			6.30	20
Benzo(g,h,i)perylene	2.00	1.61	1.61	80.5	80.5	25.0-137			0.000	23
Benzo(k)fluoranthene	2.00	1.59	1.65	79.5	82.5	39.0-128			3.70	22
Chrysene	2.00	1.77	1.85	88.5	92.5	42.0-129			4.42	20
Dibenz(a,h)anthracene	2.00	1.55	1.59	77.5	79.5	25.0-139			2.55	22
Fluoranthene	2.00	1.61	1.67	80.5	83.5	48.0-131			3.66	20

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

(LCS) R3779023-1 04/07/22 17:16 • (LCSD) R3779023-2 04/07/22 17:36

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	<u>LCS Qualifier</u>	<u>LCSD Qualifier</u>	RPD %	RPD Limits %
Fluorene	2.00	1.61	1.68	80.5	84.0	42.0-120			4.26	20
Indeno(1,2,3-cd)pyrene	2.00	1.54	1.61	77.0	80.5	37.0-133			4.44	20
Naphthalene	2.00	1.70	1.77	85.0	88.5	30.0-120			4.03	22
Phenanthrene	2.00	1.59	1.66	79.5	83.0	42.0-120			4.31	20
Pyrene	2.00	1.82	1.91	91.0	95.5	38.0-124			4.83	20
1-Methylnaphthalene	2.00	1.62	1.69	81.0	84.5	43.0-120			4.23	20
2-Methylnaphthalene	2.00	1.55	1.63	77.5	81.5	40.0-120			5.03	20
2-Chloronaphthalene	2.00	1.57	1.64	78.5	82.0	39.0-120			4.36	20
(S) Nitrobenzene-d5				90.0	92.5	11.0-135				
(S) 2-Fluorobiphenyl				85.5	89.0	32.0-120				
(S) p-Terphenyl-d14				108	111	23.0-122				

<sup>1</sup>Cp<sup>2</sup>Tc<sup>3</sup>Ss<sup>4</sup>Cn<sup>15</sup>Sr<sup>6</sup>Qc<sup>7</sup>Gl<sup>8</sup>Al<sup>9</sup>Sc

# GLOSSARY OF TERMS

## 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.

**Results Disclaimer -** Information that may be provided by the customer, and contained within this report, include Permit Limits, Project Name, Sample ID, Sample Matrix, Sample Preservation, Field Blanks, Field Spikes, Field Duplicates, On-Site Data, Sampling Collection Dates/Times, and Sampling Location. Results relate to the accuracy of this information provided, and as the samples are received.

### Abbreviations and Definitions

MDL	Method Detection Limit.
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.
Uncertainty (Radiochemistry)	Confidence level of 2 sigma.
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.

### Qualifier      Description

C3	The reported concentration is an estimate. The continuing calibration standard associated with this data responded low. Method sensitivity check is acceptable.
J	The identification of the analyte is acceptable; the reported value is an estimate.
J3	The associated batch QC was outside the established quality control range for precision.
J4	The associated batch QC was outside the established quality control range for accuracy.
J5	The sample matrix interfered with the ability to make any accurate determination; spike value is high.
J6	The sample matrix interfered with the ability to make any accurate determination; spike value is low.

<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> Gi

<sup>8</sup> Al

<sup>9</sup> Sc

# ACCREDITATIONS & LOCATIONS

Pace Analytical National 12065 Lebanon Rd Mount Juliet, TN 37122

Alabama	40660
Alaska	17-026
Arizona	AZ0612
Arkansas	88-0469
California	2932
Colorado	TN00003
Connecticut	PH-0197
Florida	E87487
Georgia	NELAP
Georgia <sup>1</sup>	923
Idaho	TN00003
Illinois	200008
Indiana	C-TN-01
Iowa	364
Kansas	E-10277
Kentucky <sup>16</sup>	KY90010
Kentucky <sup>2</sup>	16
Louisiana	AI30792
Louisiana	LA018
Maine	TN00003
Maryland	324
Massachusetts	M-TN003
Michigan	9958
Minnesota	047-999-395
Mississippi	TN00003
Missouri	340
Montana	CERT0086
A2LA – ISO 17025	1461.01
A2LA – ISO 17025 <sup>5</sup>	1461.02
Canada	1461.01
EPA-Crypto	TN00003

Nebraska	NE-OS-15-05
Nevada	TN000032021-1
New Hampshire	2975
New Jersey–NELAP	TN002
New Mexico <sup>1</sup>	TN00003
New York	11742
North Carolina	Env375
North Carolina <sup>1</sup>	DW21704
North Carolina <sup>3</sup>	41
North Dakota	R-140
Ohio–VAP	CL0069
Oklahoma	9915
Oregon	TN200002
Pennsylvania	68-02979
Rhode Island	LA000356
South Carolina	84004002
South Dakota	n/a
Tennessee <sup>14</sup>	2006
Texas	T104704245-20-18
Texas <sup>5</sup>	LAB0152
Utah	TN000032021-11
Vermont	VT2006
Virginia	110033
Washington	C847
West Virginia	233
Wisconsin	998093910
Wyoming	A2LA
AIHA-LAP,LLC EMLAP	100789
DOD	1461.01
USDA	P330-15-00234

<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

\* 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 Analytical.

<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

Company Name/Address: <b>Arcadis - Chevron - AK</b> 880 H St. Anchorage, AK 99501			Billing Information: <b>Attn: Accounts Payable 630 Plaza Dr Ste 600 Highlands Ranch, CO 80129</b>			Pres Chk	Analysis / Container / Preservative						Chain of Custody	Page <u>1</u> of <u>1</u>
Report to: <b>Sydney Clark/Erika Midkiff</b>			Email To: <b>Sydney.Clark@arcadis.com;Nicole.Monroe@arc</b>											
Project Description: <b>97324</b>		City/State Collected:	<i>Anchorage, AK</i>		Please Circle: <b>PT MT CT ET</b>									
Phone: <b>907-276-8095</b>		Client Project # <b>30063667.19.21</b>		Lab Project # <b>CHEVARCAK-97324</b>										
Collected by (print): <i>E. Wujcik</i>		Site/Facility ID # <b>4417 LAKE OTIS PKWY,</b>		P.O. #										
Collected by (signature): <i>E. Wujcik</i>		Rush? (Lab MUST Be Notified)		Quote #										
Immediately Packed on Ice N <input checked="" type="checkbox"/> Y <input type="checkbox"/>		<input type="checkbox"/> Same Day <input type="checkbox"/> Five Day <input type="checkbox"/> Next Day <input type="checkbox"/> 5 Day (Rad Only) <input type="checkbox"/> Two Day <input type="checkbox"/> 10 Day (Rad Only) <input checked="" type="checkbox"/> Three Day <input type="checkbox"/> Standard		Date Results Needed		No. of Cntrs								
Sample ID		Comp/Grab	Matrix *	Depth	Date	Time								
MW-2R-W-20220404	Gmb	GW	-	4.4.22	0900	13	X	X	X	X	X			
MW-8RR-W-20220404	Gmb	GW	-	4.4.22	1000	11	X	X	X	X	X			
MW-1R-W-20220404	Gmb	GW	-	4.4.22	1100	33	X	X	X	X	X			
MW-9-W-20220404	Gmb	GW	-	4.4.22	1200	11	X	X	X	X	X			
BD-1-W-20220404	Gmb	GW	-	4.4.22	-	13	X	X	X	X	X			
EGB-1-W-20220404	Gmb	GW	-	4.4.22	1300	13	X	X	X	X	X			
Trap Blank	-	GW	-	4.4.22	-	1	X	X	X	X	X			
		GW												
		GW												
		GW												
* Matrix: SS - Soil   AIR - Air   F - Filter GW - Groundwater   B - Bioassay WW - WasteWater DW - Drinking Water OT - Other _____	Remarks:						pH _____	Temp _____	Sample Receipt Checklist					
							Flow _____	Other _____	COC Seal Present/Intact: <input type="checkbox"/> NP <input checked="" type="checkbox"/> Y <input type="checkbox"/> N COC Signed/Accurate: <input checked="" type="checkbox"/> <input type="checkbox"/> N Bottles arrive intact: <input checked="" type="checkbox"/> <input type="checkbox"/> N Correct bottles used: <input checked="" type="checkbox"/> <input type="checkbox"/> N Sufficient volume sent: <input checked="" type="checkbox"/> <input type="checkbox"/> N <small>If Applicable</small> VOA Zero Headspace: <input checked="" type="checkbox"/> <input type="checkbox"/> N Preservation Correct/Checked: <input checked="" type="checkbox"/> <input type="checkbox"/> N RAD Screen < 0.5 mR/hr: <input checked="" type="checkbox"/> <input type="checkbox"/> N					
Samples returned via: <input checked="" type="checkbox"/> UPS <input type="checkbox"/> FedEx <input type="checkbox"/> Courier			Tracking # <i>SL71 5376 3015</i>											
Relinquished by : (Signature) <i>E. Wujcik</i>		Date: <i>4.5.22</i>	Time: <i>0800</i>	Received by: (Signature)		Trip Blank Received: <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <i>1</i>		NHC / MeOH TBR		If preservation required by Login: Date/Time				
Relinquished by : (Signature)		Date:	Time:	Received by: (Signature)		Temp <i>DLA7C</i> <i>2910-29</i>		Bottles Received: <i>94</i>						
Relinquished by : (Signature)		Date:	Time:	Received for lab by: (Signature) <i>Just L</i>		Date: <i>4/6/22</i>	Time: <i>900</i>	Hold:	Condition: <i>NCF / OK</i>					

**Pace**  
PEOPLE ADVANCING SCIENCE

**MT JULIET, TN**

12065 Lebanon Rd Mount Juliet, TN 37122  
Submitting a sample via this chain of custody constitutes acknowledgment and acceptance of the Pace Terms and Conditions found at:  
<https://info.pacelabs.com/hubfs/pas-standard-terms.pdf>

SDG # *11479046*  
**F190**

**Acctnum: CHEVARCAK**

**Template:T205504**

**Prelogin: P911868**

**PM: 110 - Brian Ford**

**PB: CR 3-16-22**

**Shipped Via:**

Remarks	Sample # (lab only)
---------	---------------------

## APPENDIX E



## **Laboratory Data Review Checklist**

Completed By:

Bhagyashree A Fulzele

Title:

Project Chemist

Date:

April 21, 2022

Consultant Firm:

ARCADIS U.S., Inc

Laboratory Name:

Pace Analytical

Laboratory Report Number:

L1479046

Laboratory Report Date:

04/15/2022

CS Site Name:

First Semi Annual 2022 Groundwater Monitoring Report

ADEC File Number:

2100.26.008

Hazard Identification Number:

23885

**Note: Any N/A or No box checked must have an explanation in the comments box.**

1. Laboratory

- a. Did an ADEC CS approved laboratory receive and perform all of the submitted sample analyses?

Yes  No  N/A  Comments:

Yes.

- b. If the samples were transferred to another “network” laboratory or sub-contracted to an alternate laboratory, was the laboratory performing the analyses ADEC CS approved?

Yes  No  N/A  Comments:

Not applicable.

2. Chain of Custody (CoC)

- a. CoC information completed, signed, and dated (including released/received by)?

Yes  No  N/A  Comments:

Yes.

- b. Correct analyses requested?

Yes  No  N/A  Comments:

Yes.

3. Laboratory Sample Receipt Documentation

- a. Sample/cooler temperature documented and within range at receipt (0° to 6° C)?

Yes  No  N/A  Comments:

Yes.

- b. Sample preservation acceptable – acidified waters, Methanol preserved VOC soil (GRO, BTEX, Volatile Chlorinated Solvents, etc.)?

Yes  No  N/A  Comments:

Yes.

- c. Sample condition documented – broken, leaking (Methanol), zero headspace (VOC vials)?

Yes  No  N/A  Comments:

No.

- d. If there were any discrepancies, were they documented? For example, incorrect sample containers/preservation, sample temperature outside of acceptable range, insufficient or missing samples, etc.?

Yes  No  N/A  Comments:

Yes.

- e. Data quality or usability affected?

Comments:

Data quality/usability was not affected.

4. Case Narrative

- a. Present and understandable?

Yes  No  N/A  Comments:

Yes.

- b. Discrepancies, errors, or QC failures identified by the lab?

Yes  No  N/A  Comments:

Yes.

- c. Were all corrective actions documented?

Yes  No  N/A  Comments:

Yes.

- d. What is the effect on data quality/usability according to the case narrative?

Comments:

Data quality/usability was not affected.

5. Samples Results

- a. Correct analyses performed/reported as requested on COC?

Yes  No  N/A  Comments:

Yes.

- b. All applicable holding times met?

Yes  No  N/A  Comments:

Yes.

- c. All soils reported on a dry weight basis?

Yes  No  N/A  Comments:

Not applicable.

- d. Are the reported LOQs less than the Cleanup Level or the minimum required detection level for the project?

Yes  No  N/A  Comments:

Yes.

e. Data quality or usability affected?

Data quality/usability was not affected.

6. QC Samples

a. Method Blank

i. One method blank reported per matrix, analysis and 20 samples?

Yes  No  N/A  Comments:

Yes.

ii. All method blank results less than limit of quantitation (LOQ) or project specified objectives?

Yes  No  N/A  Comments:

Yes.

iii. If above LOQ or project specified objectives, what samples are affected?

Comments:

None of the samples were affected.

iv. Do the affected sample(s) have data flags? If so, are the data flags clearly defined?

Yes  No  N/A  Comments:

Yes.

v. Data quality or usability affected?

Comments:

Data quality or usability was not affected.

b. Laboratory Control Sample/Duplicate (LCS/LCSD)

i. Organics – One LCS/LCSD reported per matrix, analysis and 20 samples? (LCS/LCSD required per AK methods, LCS required per SW846)

Yes  No  N/A  Comments:

Yes.

ii. Metals/Inorganics – one LCS and one sample duplicate reported per matrix, analysis and 20 samples?

Yes  No  N/A  Comments:

Not applicable.

- iii. Accuracy – All percent recoveries (%R) reported and within method or laboratory limits and project specified objectives, if applicable? (AK Petroleum methods: AK101 60%-120%, AK102 75%-125%, AK103 60%-120%; all other analyses see the laboratory QC pages)

Yes  No  N/A  Comments:

Method SW846 8260D: LCSD recovery was greater than the control limit for compounds acetone, 1,2-dibromo-3-chloropropane and methyl tert-butyl ether in preparation batch WG1844554. The compounds were non-detected in any of the associated samples; therefore, no other qualification of the data was required.

- iv. Precision – All relative percent differences (RPD) reported and less than method or laboratory limits and project specified objectives, if applicable? RPD reported from LCS/LCSD, and or sample/sample duplicate. (AK Petroleum methods 20%; all other analyses see the laboratory QC pages)

Yes  No  N/A  Comments:

Method SW846 8260D: LCS/LCSD RPD exceedances were observed for several compounds acetone, acrolein, acrylonitrile, bromobenzene, bromoform, n-butylbenzene, sec-butylbenzene, tert-butylbenzene, 2-chlorotoluene, 4-chlorotoluene, 1,2-dibromo-3-chloropropane, 1,2-dichlorobenzene, 1,3-dichlorobenzene, 1,4-dichlorobenzene, hexachloro-1,3-butadiene, p-isopropyltoluene, 2-butanone (mek), 4-methyl-2-pentanone (mibk), methyl tert-butyl ether, naphthalene, n-propylbenzene, 1,1,2,2-tetrachloroethane, 1,2,3-trichlorobenzene, 1,2,4-trichlorobenzene, 1,2,4-trimethylbenzene, 1,2,3-trimethylbenzene and 1,3,5-trimethylbenzene in preparation batch WG1844554. Compounds result in the associated samples were qualified as estimated (J/UJ).

- v. If %R or RPD is outside of acceptable limits, what samples are affected?

Comments:

Precision: Compounds result in sample IDs MW-2R-W-20220404, MW-8RR-W-20220404, MW-1R-W-20220404, MW-9-W-20220404, BD-1-W-20220404, EQB-1-W-20220404 and TRIP BLANK-W-20220404 was qualified as estimated (J/UJ).

- vi. Do the affected sample(s) have data flags? If so, are the data flags clearly defined?

Yes  No  N/A  Comments:

Yes.

- vii. Data quality or usability affected? (Use comment box to explain.)

Comments:

The LCS/LCSD recovery and RPD exceedances are considered minor and would result in the estimation of the associated data. The reported data should still consider as usable.

- c. Matrix Spike/Matrix Spike Duplicate (MS/MSD)

**Note: Leave blank if not required for project**

- i. Organics – One MS/MSD reported per matrix, analysis and 20 samples?

Yes  No  N/A  Comments:

The MS/MSD analysis was performed on sample ID MW-1R-W-20220404 for Method AK101, AK102 and SW846 8260D.

ii. Metals/Inorganics – one MS and one MSD reported per matrix, analysis and 20 samples?

Yes  No  N/A  Comments:

Not applicable.

iii. Accuracy – All percent recoveries (%R) reported and within method or laboratory limits and project specified objectives, if applicable? (AK Petroleum methods: AK101 60%-120%, AK102 75%-125%, AK103 60%-120%; all other analyses see the laboratory QC pages)

Yes  No  N/A  Comments:

Method SW846 8260D: MS and/or MSD recovery for compounds acetone, 1,2-dibromo-3-chloropropane, 2-butanone (MEK), 4-methyl-2-pentanone (MIBK) and methyl tert-butyl ether exceeded in sample ID MW-1R-W-20220404. Compounds were non-detected in the associated sample; hence no other qualification of the data was required.

Method AK101: MS/MSD recovery for compound TPHGAK C6 to C10 was less than the control limit in sample ID MW-1R-W-20220404. Compound result in the associated sample was qualified as estimated (UJ).

iv. Precision – All relative percent differences (RPD) reported and less than method or laboratory limits and project specified objectives, if applicable? RPD reported from MS/MSD, and or sample/sample duplicate. (AK Petroleum methods 20%; all other analyses see the laboratory QC pages)

Yes  No  N/A  Comments:

Method AK101: MS/MSD RPD for compound TPHGAK C6 to C10 was exceeded the control limit in sample MW-1R-W-20220404. The compound result in associated sample was qualified as estimated (UJ).

v. If %R or RPD is outside of acceptable limits, what samples are affected?

Comments:

The MS/MSD RPD exceedance was observed for compound acetone in sample ID MW-1R-W-20220404 and qualified as estimated (UJ/J).

vi. Do the affected sample(s) have data flags? If so, are the data flags clearly defined?

Yes  No  N/A  Comments:

Yes.

vii. Data quality or usability affected? (Use comment box to explain.)

Comments:

MS/MSD recovery and RPD exceedance is considered minor and would result in the estimation of associated data. The reported data should still consider as usable.

d. Surrogates – Organics Only or Isotope Dilution Analytes (IDA) – Isotope Dilution Methods Only

i. Are surrogate/IDA recoveries reported for organic analyses – field, QC and laboratory samples?

Yes  No  N/A  Comments:

Yes.

- ii. Accuracy – All percent recoveries (%R) reported and within method or laboratory limits and project specified objectives, if applicable? (AK Petroleum methods 50-150 %R; all other analyses see the laboratory report pages)

Yes  No  N/A  Comments:

Yes.

- iii. Do the sample results with failed surrogate/IDA recoveries have data flags? If so, are the data flags clearly defined?

Yes  No  N/A  Comments:

Not applicable.

- iv. Data quality or usability affected?

Comments:

Data quality or usability was not affected.

e. Trip Blanks

- i. One trip blank reported per matrix, analysis and for each cooler containing volatile samples?  
(If not, enter explanation below.)

Yes  No  N/A  Comments:

Trip blank sample was collected as TRIP BLANK-W-20220404.

- ii. Is the cooler used to transport the trip blank and VOA samples clearly indicated on the COC?  
(If not, a comment explaining why must be entered below)

Yes  No  N/A  Comments:

Yes.

- iii. All results less than LOQ and project specified objectives?

Yes  No  N/A  Comments:

Yes.

- iv. If above LOQ or project specified objectives, what samples are affected?

Comments:

None of the samples were affected.

- v. Data quality or usability affected?

Comments:

Data quality or usability was not affected.

f. Field Duplicate

- i. One field duplicate submitted per matrix, analysis and 10 project samples?

Yes  No  N/A  Comments:

Yes.

ii. Submitted blind to lab?

Yes  No  N/A  Comments:

Field duplicate BD-1-W-20220404 was collected from sample MW-2R-W-20220404.

iii. Precision – All relative percent differences (RPD) less than specified project objectives?  
(Recommended: 30% water, 50% soil)

$$\text{RPD (\%)} = \text{Absolute value of: } \frac{(R_1 - R_2)}{((R_1 + R_2)/2)} \times 100$$

Where  $R_1$  = Sample Concentration  
 $R_2$  = Field Duplicate Concentration

Yes  No  N/A  Comments:

Yes.

iv. Data quality or usability affected? (Use the comment box to explain why or why not.)

Comments:

Data quality/usability was not affected.

g. Decontamination or Equipment Blank (If not applicable, a comment stating why must be entered below)?

Yes  No  N/A  Comments:

Equipment blank sample was collected as EQB-1-W-20220404.

i. All results less than LOQ and project specified objectives?

Yes  No  N/A  Comments:

Yes.

ii. If above LOQ or project specified objectives, what samples are affected?

Comments:

None of the samples were affected.

iii. Data quality or usability affected?

Comments:

Data quality or usability was not affected.

## 7. Other Data Flags/Qualifiers (ACOE, AFCEE, Lab Specific, etc.)

a. Defined and appropriate?

Yes  No  N/A  Comments:

Yes.

## 8. Additional Laboratory Data Qualification

Method SW846 8260D: Continuing calibration for compound acrolein was exhibited a low bias recovery. Compound result in the associated sample IDs MW-2R-W-20220404, MW-8RR-W-20220404, MW-1R-W-20220404, MW-9-W-20220404, BD-1-W-20220404, EQB-1-W-20220404 and TRIP BLANK-W-20220404 was qualified as estimated (UJ).