

2021 Groundwater Monitoring Report Revision 1

For

Essential 1 Gas Station (Former Shoreside Texaco)

Mile 100 / 100.7 Seward Highway

Girdwood, Alaska 99587

ADEC File. No. 2105.26.001

Prepared for:

ALASKA DEPARTMENT OF ENVIRONMENTAL CONSERVATION

555 CORDOVA STREET

ANCHORAGE, ALASKA 99501

Prepared by:

ENVIRONMENTAL MANAGEMENT, INC.

206 EAST FIREWEED LANE, SUITE 201

ANCHORAGE, ALASKA 99503

April 2022

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Prepared by:

Environmental Management, Inc.
206 East Fireweed Lane, Suite 201
Anchorage, Alaska 99503

Written by:



Andy Coulson, QEP
Environmental Scientist

Reviewed by:



Glenn Hasburgh, QEP
Project Manager

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ACRONYMS AND ABBREVIATIONS

| | |
|---------|---|
| AAC | Alaska Administrative Code |
| ADEC | Alaska Department of Environmental Conservation |
| ADOT&PF | Alaska Department of Transportation and Public Facilities |
| AK | Alaska Method |
| bgs | Below ground surface |
| DRO | Diesel Range Organics |
| EMI | Environmental Management, Inc. |
| FSG | Field Sampling Guidance |
| IDW | Investigation Derived Waste |
| LCS | Laboratory Control Sample |
| LCSD | Laboratory Control Sample Duplicate |
| mg/L | Milligrams per liter |
| PAH | Polycyclic Aromatic Hydrocarbons |
| PPE | Personal Protective Equipment |
| QC | Quality Control |
| QEP | Qualified Environmental Professional |
| RPD | Relative Percent Difference |
| SGS | SGS North America |
| UST | Underground Storage Tank |
| VOC | Volatile Organic Compounds |

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This report describes the annual groundwater monitoring event conducted at the Essential 1 gas station (former Shoreside Texaco) in Girdwood, Alaska, undertaken by Environmental Management, Inc. (EMI) on September 10, 2021.

Groundwater monitoring activities were conducted under EMI's work plan that was approved by the Alaska Department of Environmental Conservation (ADEC) on August 13, 2021. This work was performed under contract with Shoreside Petroleum, Inc., the owner of the site. Previous quarterly sampling events prior to September 2020 were conducted by Tellus, Ltd.

1.0 SITE DESCRIPTION AND BACKGROUND

The site is the Essential 1 gas station (former Shoreside Texaco), which is located at 29383 Seward Highway, Girdwood, Alaska (Mile 100.7); Due to previously identified petroleum contamination the site is an ADEC listed contaminated site, ADEC File No. 2105.26.001.

According to the site chronology listed on the ADEC Contaminated Sites Database, contamination was originally identified at the site during the removal of two gasoline underground storage tanks (UST) in 1995. Additional contamination was identified in 2000 during the removal of a diesel dispenser and associated piping. A release investigation conducted in 2002 identified petroleum-based contamination extending to depths down to 62 feet below ground surface (bgs), in several locations. Between the 2002 release investigation and a 2014 borehole investigation, six monitoring wells were installed at the site. Three of the wells, MW1, MW2, and MW3, represent a deeper aquifer at the site and the total well depths extend down to approximately 80 feet below ground surface (bgs). These three wells have had sufficient water volume and have been monitored on a regular basis since at least 2016; however, the other three wells, MW4a, MW5 and MW6, which represent a shallower aquifer and extend to approximately 50 feet bgs, have had insufficient water for sampling and viable samples have never been collected from them. In 2020 free product was observed in two of the shallow aquifer wells, MW5 and MW6.

Prior to this September 2021 sampling event, sampling has been conducted on a quarterly basis. Gasoline range organics (GRO) and benzene, toluene, ethylbenzene, and xylenes (BTEX) have not been detected in any monitoring event, while diesel range organics (DRO) were only detected above their groundwater cleanup level once, in well MW2 in August 2019 (Table 2). Following EMI's September 2020 sampling event, which again showed no contamination within the deep aquifer, EMI recommended a reduction in the sampling frequency from quarterly to annually. This

sampling event, which is the subject of this report, is the first sampling event under the reduced frequency.

2.0 PROJECT PURPOSE AND OBJECTIVES

The purpose of the groundwater sampling was to collect additional data from the wells in the deeper aquifer, including MW3 which was unable to be located during the September 2020 sampling event, to show that contamination has not migrated to the groundwater in this aquifer. The objectives also included investigating the shallow aquifer wells (MW4a, MW5 and MW6) to determine if free product was still present within the wells and/or to determine if there was sufficient volume for sampling.

3.0 CONTAMINANTS OF CONCERN

The ADEC Contaminated Site's database identifies contamination associated with two gasoline USTs. Due to the age of those former USTs (installed in 1972), leaded gasoline was likely used. Previous investigations included the analyses of constituents typically found in gasoline during the time period including lead, EDB, polyaromatic hydrocarbons (PAH), and 1,2-DCB; none of these contaminants were detected. In addition, GRO and BTEX were analyzed over seven monitoring events between October 2016 and May 2020. GRO and BTEX were not detected and have thus been removed from the sampling program, however the full suite of volatile organic compounds (VOC), including MTBE, had not been analyzed. The constituents of potential concern (COPC) associated with the site as identified in previous investigations include the following petroleum-related compounds: diesel range organics (DRO), VOC (including MTBE), and PAH.

Water samples collected during the field activities compared to the cleanup levels outlined in the following regulations:

- Groundwater: Table C in 18 AAC 75.345, *Oil and Other Hazardous Substances Pollution Control* (November 18, 2021)

4.0 FIELD ACTIVITIES

Field activities were conducted on September 10, 2021. Andy Coulson, ADEC Qualified Environmental Professional (QEP), served as the field lead. Photographs of the field activities are provided in Appendix A. Field notes are included in Appendix B. A figure showing the monitoring well locations is provided as Figure 2.

4.1 Work Plan Variances

The following work plan variances are noted:

- Monitoring well MW3 could not be located using a metal detector and was therefore not sampled. This well was also unable to be located in 2020.

- Free product was observed in wells MW5 and MW6; no laboratory samples were collected from these wells. Free product was also believed present in MW2 and because of this, not sampled (see discussion related to MW2).
- There was insufficient water in well MW4a to collect a sample.

4.2 Groundwater Sampling

Due to insufficient well volume and/or the presence of free product, only one well was sampled during this event, MW1. Monitoring well MW1 was sampled and purged using a submersible pump and low flow methods as presented in the Environmental Protection Agency's 2010 *Low Stress (Low-Flow) Purging and Sampling Procedures for the Collection of Groundwater Samples from Monitoring Wells*. The pump was lowered to draw water from three feet below the water's surface in the well; the screened interval of the well was not reported in previous site documentation. Well MW1 was purged for one hour and twelve minutes prior to sampling. A YSI 556 multi-parameter meter was connected to the sampling apparatus using a flow-through cell and was used to monitor temperature, pH, conductivity, redox potential and dissolved oxygen. The well was purged until three well volumes had been removed. Purge water was collected in five-gallon buckets. The analytical water samples were collected by filling the laboratory-provided, hydrochloric acid-preserved and non-preserved sample containers from the discharge tube of the flow-through cell. Containers for VOC analysis were filled first, followed by containers for PAH analysis, then containers for DRO analysis. Sample containers were immediately placed in a cooler with frozen gel ice after being filled.

As previously stated, only MW1 was sampled during this event, which represents the deeper aquifer. From this well a primary and duplicate sample were collected; these are samples 18190-MW1-091021 and 18190-MW77-091021, respectively.

Of the other two deep aquifer wells, MW2 was not sampled due to the perceived presence of free product in the well (see discussion below regarding this) and MW3 was not sampled due to it being unable to be located, even when using a metal detector.

From the shallow aquifer wells, none were sampled. This is because wells MW5 and MW6 had free product up to 0.2 feet thick present. No sample was collected from well MW4a because there was insufficient water in the well to collect a sample.

4.3 Decontamination Procedures

All equipment was decontaminated between use in different wells and after sampling was complete. Equipment was decontaminated using a three-bucket process where it was first brushed in potable water with Alconox, then rinsed with potable water, and then rinsed with deionized water. Potable water was obtained from the faucet in the gas station building. The flow-through chamber was fully disassembled and all parts and fittings were decontaminated individually. The pump was decontaminated using a brush, and also powered up and briefly ran in water from each decontamination bucket. Long pieces of reusable equipment such as the water level meter and the electrical leader to the pump were fully submerged such that the entire length taken off the roll

during their most recent use was submerged in each stage of decontamination; in addition, this length was also pulled through a brush while in the Alconox and water bucket.

After decontaminating the interface probe after using it in well MW5, sheen was observed on water in the first (water and alconox) bucket. When this happened, the water from all three decontamination buckets was containerized in a different bucket and replaced with new water, and the product interface probe was decontaminated a second time in the new water.

5.0 INVESTIGATIVE DERIVED WASTE MANAGEMENT

Investigation derived waste (IDW) included decontamination water, purge water from well sampling, and disposable personal protective equipment (PPE) and sampling equipment.

Purge and decontamination water, which totaled approximately 15 gallons, were combined into three five-gallon buckets. A Contaminated Media Transport and Treatment or Disposal Approval form has been completed and submitted to ADEC. This request was approved by ADEC on November 11, 2021. The water was disposed of at US Ecology Alaska, LLC, located at 2020 Viking Drive, Anchorage, Alaska 99501 on December 20, 2021. The ADEC approved transport request and the bill of lading are included in Appendix D.

Disposable sampling equipment included nitrile gloves and tubing. These were disposed of as solid waste in the dumpster on site.

6.0 RESULTS

The following describes the results of the 2021 groundwater monitoring activities.

6.1 Surface Observations

The following observations were made of surface conditions at the site during groundwater monitoring on September 10, 2021:

- Well MW3 could not be located. Previously, both EMI and Shoreside personnel searched the area between the store building and AST where figures in previous reports had shown MW3, and where shoreside personnel had observed previous sampling events being conducted. During this event, a metal detector was used to try and locate the well to no avail. (Photo 1)
- In 2020, the Alaska Department of Transportation and Public Facilities (ADOT&PF) had a project ongoing on roads near the site. As part of this project, pavement around well MW4a had been removed, along with its surface monument. The road construction project has been completed and the damaged monument has been repaired. (Photo 2)
- The surface monuments of wells MW2 and MW5 (Photo 3) had filled in with sediment and water above the level of the well casing when opened. Small flecks of sheen were observed on the water inside these surface monuments. This water and sediment was removed by

hand prior to opening the plug on the well casing. In MW2, the seal between the well casing and plug appeared to be intact, but in MW5 the plug had floated loose of the casing.

6.2 Subsurface Observations

The oil/water interface probe indicated that petroleum may be present in both monitoring wells MW1 and MW2 with an approximate thickness of 0.01 feet. However, there was not visual or olfactory evidence of free product when sampling MW1.

In MW5 there was no product detected with the interface probe, however the probe did have a sheen on it when removed from the well and it also had a petroleum odor; sheen was observed on the decontamination water after decontaminating the product interface probe and tape after measuring this well, and the decontamination water was replaced with fresh potable water. In 2020, 0.25 feet of free product was measured in this well.

Within monitoring well MW6 0.21 feet of free product was detected with the interface probe. The presence of free product was confirmed by observing product on the probe once it was removed from the well along with a petroleum odor.

Table 1 Depths to product, groundwater, and well base

| Well | Well Elevation ¹ | Depth to Water ² | Depth to Well Base ² | Groundwater Elevation ¹ | Depth to Product ² | Product Thickness (2021) ² | Product Thickness (2020) ² |
|------|-----------------------------|-----------------------------|---------------------------------|------------------------------------|-------------------------------|---------------------------------------|---------------------------------------|
| MW1 | 99.87 | 69.44 | 80.77 | 30.43 | 69.43 | 0.01 | - |
| MW2 | 98.97 | 68.54 | 77.36 | 30.43 | 68.53 | 0.01 | - |
| MW4a | 98.53 | 47.79 | 50.40 | 50.74 | - | - | - |
| MW5 | 98.85 | 46.37 | 49.89 | 52.48 | - | - | 0.25 |
| MW6 | 99.85 | 47.31 | 50.38 | 52.54 | 47.10 | 0.21 | 0.16 |

Notes

- 1 Elevation recorded in feet relative to a horizontal strut on a utility pole northeast of MW2, which was arbitrarily assigned an elevation of 100 ft.
- 2 All depths are recorded in feet from the highest point of the well casing
- Product was not observed in these wells

6.3 Groundwater Analytical Results

Samples for laboratory analysis were submitted to SGS North America (SGS) in Anchorage, Alaska. In accordance with the approved work plan the samples were analyzed for diesel range organics (DRO) by method AK 102, volatile organic compounds (VOC) by EPA method 8260D, and polycyclic aromatic hydrocarbon (PAH) by EPA method 8270D.

Analysis of the two samples (primary and duplicate) collected during this effort found all analytes to be below the laboratory reporting limits. With the exception of 1,2,3-trichloropropane all reporting limits were below ADEC groundwater cleanup levels.

Table 2 presents the analytical results. The laboratory report is included in Appendix C.

7.0 QUALITY CONTROL

The ADEC Laboratory Data Review checklist was completed for the laboratory report. Both the checklist and report are included in Appendix C. At the time of receipt at the laboratory the cooler temperature was 3.8°C, which is within the acceptable range. Sample condition was documented as being acceptable with no containers broken or leaking, vials were free of air bubbles, sample containers matched the COC, and the samples were received and analyzed within hold times.

The laboratory quality control (QC) samples did have some failures with regards to recoveries in the laboratory control samples (LCS) and laboratory control sample duplicates (LCSD). The relative percent difference (RPD) for naphthalene was also outside of the acceptable range between the LCS and LCSD. Surrogates toluene-D8 and 4-bromoflorobenzene also had failed recoveries in the method blank. Additionally, in the matrix spike and matrix spike duplicate there were some failed surrogate recoveries and some RPDs that were outside of range.

The above stated issues related to the laboratory QC data does not have any significant impact on the data since analysis of the field samples found all analytes to be below the laboratory reporting limits, and with the exception of 1,2,3-trichloropropane, all reporting limits were below the ADEC cleanup levels.

The RPD could not be calculated for the field samples due to the primary and duplicate samples being non-detect for all analytes.

No other discrepancies that would affect project conclusion were noted.

8.0 CONCLUSION

During this sampling event only one well was sampled, monitoring well MW1. The oil/water interface probe indicated that free product may be present in both MW1 and MW2. Due to the perceived presence of the free product, MW2 was not sampled. Even though MW1 also had an indication of free product, the QEP decided to sample the well to get some data on the deep aquifer. At no time during the purging of the well or during the sample collection was indications of contamination observed. There was no sheen, oil droplets, odor or other indicators of contamination. The two samples collected, sample 18190-MW1-091021 and duplicate sample 18190-M77-091021 did not have any analytes detected above the laboratory's reporting limit. For all primary contaminants of concern the reporting limits were below the cleanup level. Based on this data it appears the deep aquifer still remains free of contamination. The reason for the indication of free product by the interface probe remains unknown as no visual or olfactory indicators were observed during purging and sampling. This should be further evaluated during future events if probe again indicates the potential for free product. During future events, wells MW1 and MW2 should be sampled even if product is indicated; however, sampling may be suspended if free product is visually confirmed in these wells.

During the September 2020 and September 2021 sampling events sufficient effort has been taken to try and locate MW3, to no avail. At this point MW3 should be considered to be inaccessible for sampling.

The shallow aquifer wells MW4a, MW5 and MW6 have never yielded a viable sample due to the combination of insufficient water volume and the presence of free product. These wells are still useful for monitoring the depth of the free product and they should continue to be investigated during future monitoring events.

Table 2 DRO results of samples from monitoring wells on at the Essential 1 Gas Station. All results are shown in milligrams per liter (mg/L). GRO and BTEX were analyzed for in samples from October 2016 through May 2020 and were never detected.

| Date | MW1 | MW2 | MW3 |
|--------------|----------------|----------------|----------------|
| September-21 | 0.566 u | - | - |
| September-20 | 0.652 u,1 | 0.625 u | - |
| May-20 | 0.213 j | 0.268 j | 0.329 j |
| December-19 | 0.25 u | 0.25 u | 0.25 u |
| August-19 | 0.25 | 2.2 | 0.33 |
| October-17 | 0.26 u | 0.25 u | 0.25 u |
| May-17 | 0.099 u | 0.10 u | 0.10 u |
| October-16 | 0.11 | 0.10 u | 0.10 u |

Key

- 1 Lowest non-detect reporting limit shown
- No sample was collected from this well during this sampling event
- 0.652 u DRO was not detected above the reporting limit of 0.652 mg/L
- 0.213 j** DRO was detected at an estimated concentration of 0.213 mg/L
- 0.25** DRO was detected at 0.25 mg/L
- 2.2** DRO was detected at 2.2 mg/L, and this concentration is greater than DRO's cleanup level of 1.5 mg/L



VICINITY MAP

ESSENTIAL 1 STATION,
29383 SEWARD HIGHWAY,
GIRDWOOD, ALASKA

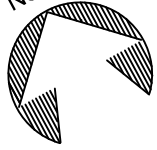


EMI JOB: 18190
DRAWN: HJD
REVIEWED: AVC
DATE: 6/25/2021

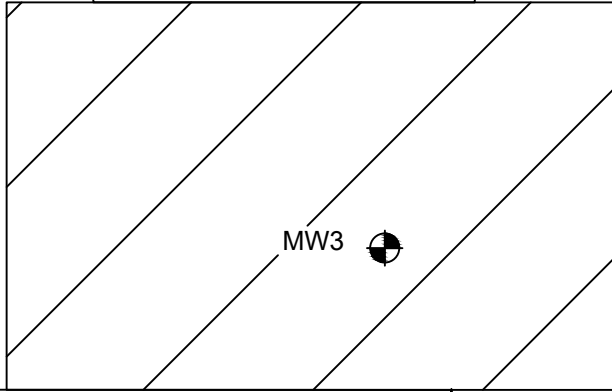
FIGURE

1

North



AST ENCLOSURE



AREA SEARCHED FOR MW3, WELL NOT FOUND

GAS STATION BUILDING



FUEL DISPENSERS

MW1



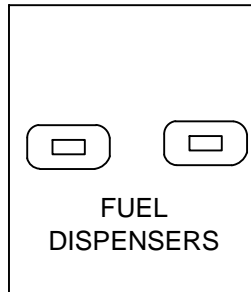
MW6



MW2



MW5



LEGEND

 MONITORING WELL LOCATION

MW4a



GRAPHIC SCALE (IN FEET)

SEWARD HIGHWAY

SITE MAP

ESSENTIAL 1 STATION,
29383 SEWARD HIGHWAY,
GIRDWOOD, ALASKA



EMI JOB: 18190
DRAWN: HJD
REVIEWED: AVC
DATE:
6/25/2021

FIGURE

2

APPENDIX A

Photo Log

Photo Log



Photo 1: Area where records indicate MW3 is located. The white van and gold truck could not be moved, but their locations were thoroughly searched in 2020 and are paved. September 10, 2021, facing west.



Photo 2: Monitoring well MW4a with new surface monument. September 10, 2021, facing south.

Photo Log



Photo 3: Silt accumulated in the surface monument of MW5. September 10, 2021.



Photo 4: Sheen observed on decontamination water after measuring MW5. September 10, 2021.

APPENDIX B

Field Notes

8th Sept 2021
 Bird Creek essential 1

Andy Carlson - EMI QEP

on site 11:10

Cloudy, varied weather, calm.
 ~ 60% RH

No luck finding MWS w/
 metal detector.

MW 4A

WL 47.73

bed 50.40

No odor noted, bottom
 of probe covered
 in mud/silt.

MW 6

product at 47.10

water at 47.31 > 7.07

base of well 50.38

Fuel odor residue on probe
 shown on log in drum imp
 after decanting fuel.

13:00 MWS - surface mound
 silted in. Plug
 appears to have floated
 off - only loosely on
 casing

WL 46.37 > 3.52
 base 49.89

~ 2 inches silt on probe
 - silt on silt on probe, odor
 - no product detection
 by interface probe

MW 2 - casing completely covered
 by silt & silt, cement
 by hand

GW Elev 30.44

product time just just 66.53

WL 66.54

well base 77.36

MW 1

product time 69.43

GW Elev 30.43

WL 69.44

well base 80.77

11.33 - 0.16 1.81 - low

5.44 - 30V
 Rate in the Rain

| TBS | OS | FS | Elev. |
|------|------|------|---------|
| TBS1 | 3.58 | | 100.00 |
| MW4 | 4.81 | 5.05 | 98.53 |
| TBS1 | | 3.34 | 100.00 |
| TBS1 | 3.77 | | 100.00 |
| MW5 | 4.79 | 4.92 | 98.85 |
| TBS1 | | 3.64 | 100.00 |
| TBS1 | 3.58 | | 100.00 |
| MW6 | 3.85 | 3.73 | 99.85 |
| TBS1 | | 3.70 | 100.00 |
| TBS1 | 3.72 | | 100.00 |
| MW2 | 4.78 | 4.74 | 98.98 |
| TBS1 | | 3.79 | 99.97 x |

Scale: 1 square =

| TBS | OS | FS | Elev. |
|------|------|--------------------|--------|
| TBS1 | 3.70 | | 100.00 |
| MW2 | 4.78 | 4.73 | 98.97 |
| TBS1 | | 3.76 | 99.99 |
| TBS1 | 3.67 | | 100.00 |
| MW1 | 3.65 | 3.80 ^{re} | 99.87 |
| TBS1 | | 3.52 | 100.00 |

MW1 - setting pump initially at 71.97

began pumping at 16:07
 pump at 76.10, WL neither not lifting

| Time | WL | °C | % O ₂ | MS/cm | pH | ORP |
|-------|-------|-----|------------------|-------|------|-------|
| 16:14 | - | 6.2 | 2.10 | 104.1 | 6.20 | 275.9 |
| 16:18 | - | 5.6 | 2.23 | 101.6 | 6.68 | 261 |
| 16:39 | 69.50 | 5.3 | 2.59 | 100 | 7.37 | 217.4 |
| 17:13 | 69.51 | 5.5 | 10.03 | 101.2 | 7.90 | 163.5 |

Scale: 1 square =

Rite in the Rain

16:20 pausing purging to
see if screen apprs.

16:22 no screen, restarted pump

17:05 restarted pump

17:26 } 3 wv purged
(> 7 gallons)
collected samples
18190 - MW1 - 091021
and its duplicate
18190 - MW77 - 091021
each

3 x 90 ml w/HCl
2 x 250 ml no pres
2 x 250 ml w/HCl

17:30 MW1 - w/v 69.49
product 69.98

Swing trees - feet

Alfa

| | N. Corner | S. Corner |
|-----------|-----------|-----------|
| N. Corner | | 36.1 |
| MW 1 | 41.8 | 21.3 |
| MW 6 | 46.8 | 29.0 |
| MW 2 | 59.8 | 61.5 |
| MW 5 | 71.2 | 65.5 |
| MW 4a | 98.6 | 70.3 |

18:16 off site.

APPENDIX C

Laboratory Reports and ADEC Laboratory Data Review Checklists



Laboratory Report of Analysis

To: Environmental Mgmt Inc (EMI)
206 E Fireweed Ln #201
Anchorage, AK 99503
(907)272-9336

Report Number: **1215973**

Client Project: **18190 Essential 1**

Dear Andy Coulson,

Enclosed are the results of the analytical services performed under the referenced project for the received samples and associated QC as applicable. The samples are certified to meet the requirements of the National Environmental Laboratory Accreditation Conference Standards. Copies of this report and supporting data will be retained in our files for a period of ten years in the event they are required for future reference. All results are intended to be used in their entirety and SGS is not responsible for use of less than the complete report. Any samples submitted to our laboratory will be retained for a maximum of fourteen (14) days from the date of this report unless other archiving requirements were included in the quote.

If there are any questions about the report or services performed during this project, please call Alexandra at (907) 562-2343. We will be happy to answer any questions or concerns which you may have.

Thank you for using SGS North America Inc. for your analytical services. We look forward to working with you again on any additional analytical needs.

Sincerely,
SGS North America Inc.

Alexandra Daniel
Project Manager
Alexandra.Daniel@sgs.com

Date

Case Narrative

SGS Client: **Environmental Mgmt Inc (EMI)**

SGS Project: **1215973**

Project Name/Site: **18190 Essential 1**

Project Contact: **Andy Coulson**

Refer to sample receipt form for information on sample condition.

LCSD for HBN 1825633 [VXX/3785 (1636753) LCSD

8260D - LCSD recoveries for several analytes do not meet QC criteria. These analytes were not reported above the LOQ in the associated samples.

8260D - LCS/LCSD RPD for naphthalene does not meet QC criteria. This analyte was not reported above the LOQ in the associated samples.

MB for HBN 1825633 [VXX/37856] (1636751) MB

8260D - Surrogate recovery for 4-bromofluorobenzene does not meet QC criteria.

MB for HBN 1825978 [VXX/37888] (1637560) MB

8260D - Surrogate recovery for toluene-d8 does not meet QC criteria, however the associated analytes were not detected above the LOQ.

1215976003MS (1636319) MS

8270D SIM - PAH MS recoveries for multiple analytes do not meet QC criteria. Refer to the LCS for accuracy requirements.

8270D SIM - PAH surrogate recoveries for 2-methylnaphthalene-d10 and fluoranthene-d10 do not meet QC criteria.

1215976003MSD (1636320) MSD

8270D SIM - PAH MS/MSD RPDs for multiple analytes do not meet QC criteria. These analytes were not detected above the LOQ in the parent sample.

*QC comments may be associated with the field samples found in this report. When applicable, comments will be applied to associated field samples.

Print Date: 09/30/2021 2:29:23PM

Laboratory Qualifiers

Enclosed are the analytical results associated with the above work order. The results apply to the samples as received. All results are intended to be used in their entirety and SGS is not responsible for use of less than the complete report. This document is issued by the Company under its General Conditions of Service accessible at <http://www.sgs.com/en/Terms-and-Conditions.aspx>. Attention is drawn to the limitation of liability, indemnification and jurisdiction issues defined therein.

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SGS maintains a formal Quality Assurance/Quality Control (QA/QC) program. A copy of our Quality Assurance Plan (QAP), which outlines this program, is available at your request. The laboratory certification numbers are AK00971 (DW Chemistry & Microbiology) & 17-021 (CS) for ADEC and 2944.01 for DOD ELAP/ISO17025 (RCRA methods: 1020B, 1311, 3010A, 3050B, 3520C, 3550C, 5030B, 5035A, 6020B, 7470A, 7471B, 8015C, 8021B, 8082A, 8260D, 8270D, 8270D-SIM, 9040C, 9045D, 9056A, 9060A, AK101 and AK102/103). SGS is only certified for the analytes listed on our Drinking Water Certification (DW methods: 200.8, 2130B, 2320B, 2510B, 300.0, 4500-CN-C,E, 4500-H-B, 4500-NO3-F, 4500-P-E and 524.2) and only those analytes will be reported to the State of Alaska for compliance. Except as specifically noted, all statements and data in this report are in conformance to the provisions set forth by the SGS QAP and, when applicable, other regulatory authorities.

The following descriptors or qualifiers may be found in your report:

| | |
|--------------------|---|
| * | The analyte has exceeded allowable regulatory or control limits. |
| ! | Surrogate out of control limits. |
| B | Indicates the analyte is found in a blank associated with the sample. |
| CCV/CVA/CVB | Continuing Calibration Verification |
| CCCV/CVC/CVCA/CVCB | Closing Continuing Calibration Verification |
| CL | Control Limit |
| DF | Analytical Dilution Factor |
| DL | Detection Limit (i.e., maximum method detection limit) |
| E | The analyte result is above the calibrated range. |
| GT | Greater Than |
| IB | Instrument Blank |
| ICV | Initial Calibration Verification |
| J | The quantitation is an estimation. |
| LCS(D) | Laboratory Control Spike (Duplicate) |
| LLQC/LLIQC | Low Level Quantitation Check |
| LOD | Limit of Detection (i.e., 1/2 of the LOQ) |
| LOQ | Limit of Quantitation (i.e., reporting or practical quantitation limit) |
| LT | Less Than |
| MB | Method Blank |
| MS(D) | Matrix Spike (Duplicate) |
| ND | Indicates the analyte is not detected. |
| RPD | Relative Percent Difference |
| TNTC | Too Numerous To Count |
| U | Indicates the analyte was analyzed for but not detected. |

Note: Sample summaries which include a result for "Total Solids" have already been adjusted for moisture content. All DRO/RRO analyses are integrated per SOP.

Sample Summary

| <u>Client Sample ID</u> | <u>Lab Sample ID</u> | <u>Collected</u> | <u>Received</u> | <u>Matrix</u> |
|-------------------------|----------------------|------------------|-----------------|-------------------------------|
| 18190-MW1-091021 | 1215973001 | 09/10/2021 | 09/13/2021 | Water (Surface, Eff., Ground) |
| 18190-MW77-091021 | 1215973002 | 09/10/2021 | 09/13/2021 | Water (Surface, Eff., Ground) |
| Trip Blank | 1215973003 | 09/10/2021 | 09/13/2021 | Water (Surface, Eff., Ground) |

| <u>Method</u> | <u>Method Description</u> |
|--------------------|-------------------------------------|
| 8270D SIM LV (PAH) | 8270 PAH SIM GC/MS LV |
| AK102 | DRO Low Volume (W) |
| SW8260D | Volatile Organic Compounds (W) FULL |

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Results of 18190-MW1-091021

Client Sample ID: 18190-MW1-091021
Client Project ID: 18190 Essential 1
Lab Sample ID: 1215973001
Lab Project ID: 1215973

Collection Date: 09/10/21 17:26
Received Date: 09/13/21 14:00
Matrix: Water (Surface, Eff., Ground)
Solids (%):
Location:

Results by Polynuclear Aromatics GC/MS

Table with 8 columns: Parameter, Result Qual, LOQ/CL, DL, Units, DF, Allowable Limits, Date Analyzed. Lists various polynuclear aromatic hydrocarbons and their surrogate values.

Batch Information

Analytical Batch: XMS12899
Analytical Method: 8270D SIM LV (PAH)
Analyst: LAW
Analytical Date/Time: 09/20/21 23:27
Container ID: 1215973001-F

Prep Batch: XXX45561
Prep Method: SW3535A
Prep Date/Time: 09/14/21 12:30
Prep Initial Wt./Vol.: 275 mL
Prep Extract Vol: 1 mL



Results of **18190-MW1-091021**

Client Sample ID: **18190-MW1-091021**
Client Project ID: **18190 Essential 1**
Lab Sample ID: 1215973001
Lab Project ID: 1215973

Collection Date: 09/10/21 17:26
Received Date: 09/13/21 14:00
Matrix: Water (Surface, Eff., Ground)
Solids (%):
Location:

Results by **Semivolatile Organic Fuels**

| <u>Parameter</u> | <u>Result Qual</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> | <u>DF</u> | <u>Allowable Limits</u> | <u>Date Analyzed</u> |
|-----------------------|--------------------|---------------|-----------|--------------|-----------|-------------------------|----------------------|
| Diesel Range Organics | 0.566 U | 0.566 | 0.189 | mg/L | 1 | | 09/22/21 20:18 |
| Surrogates | | | | | | | |
| 5a Androstane (surr) | 79.6 | 50-150 | | % | 1 | | 09/22/21 20:18 |

Batch Information

Analytical Batch: XFC16086
Analytical Method: AK102
Analyst: JMG
Analytical Date/Time: 09/22/21 20:18
Container ID: 1215973001-D

Prep Batch: XXX45584
Prep Method: SW3520C
Prep Date/Time: 09/18/21 15:55
Prep Initial Wt./Vol.: 265 mL
Prep Extract Vol: 1 mL

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Results of 18190-MW1-091021

Client Sample ID: 18190-MW1-091021
Client Project ID: 18190 Essential 1
Lab Sample ID: 1215973001
Lab Project ID: 1215973

Collection Date: 09/10/21 17:26
Received Date: 09/13/21 14:00
Matrix: Water (Surface, Eff., Ground)
Solids (%):
Location:

Results by Volatile GC/MS

| <u>Parameter</u> | <u>Result Qual</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> | <u>DF</u> | <u>Allowable Limits</u> | <u>Date Analyzed</u> |
|-----------------------------|--------------------|---------------|-----------|--------------|-----------|-------------------------|----------------------|
| 1,1,1,2-Tetrachloroethane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:06 |
| 1,1,1-Trichloroethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 1,1,2,2-Tetrachloroethane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:06 |
| 1,1,2-Trichloroethane | 0.400 U | 0.400 | 0.120 | ug/L | 1 | | 09/15/21 15:06 |
| 1,1-Dichloroethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 1,1-Dichloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 1,1-Dichloropropene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 1,2,3-Trichlorobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 1,2,3-Trichloropropane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 1,2,4-Trichlorobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 1,2,4-Trimethylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 1,2-Dibromo-3-chloropropane | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:06 |
| 1,2-Dibromoethane | 0.0750 U | 0.0750 | 0.0180 | ug/L | 1 | | 09/15/21 15:06 |
| 1,2-Dichlorobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 1,2-Dichloroethane | 0.500 U | 0.500 | 0.200 | ug/L | 1 | | 09/15/21 15:06 |
| 1,2-Dichloropropane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 1,3,5-Trimethylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 1,3-Dichlorobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 1,3-Dichloropropane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:06 |
| 1,4-Dichlorobenzene | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:06 |
| 2,2-Dichloropropane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 2-Butanone (MEK) | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:06 |
| 2-Chlorotoluene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 2-Hexanone | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:06 |
| 4-Chlorotoluene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 4-Isopropyltoluene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| 4-Methyl-2-pentanone (MIBK) | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:06 |
| Benzene | 0.400 U | 0.400 | 0.120 | ug/L | 1 | | 09/15/21 15:06 |
| Bromobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Bromochloromethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Bromodichloromethane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:06 |
| Bromoform | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Bromomethane | 6.00 U | 6.00 | 3.00 | ug/L | 1 | | 09/15/21 15:06 |
| Carbon disulfide | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:06 |
| Carbon tetrachloride | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Chlorobenzene | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:06 |
| Chloroethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |

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Results of **18190-MW1-091021**

Client Sample ID: **18190-MW1-091021**
 Client Project ID: **18190 Essential 1**
 Lab Sample ID: 1215973001
 Lab Project ID: 1215973

Collection Date: 09/10/21 17:26
 Received Date: 09/13/21 14:00
 Matrix: Water (Surface, Eff., Ground)
 Solids (%):
 Location:

Results by **Volatile GC/MS**

| <u>Parameter</u> | <u>Result Qual</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> | <u>DF</u> | <u>Allowable Limits</u> | <u>Date Analyzed</u> |
|------------------------------|--------------------|---------------|-----------|--------------|-----------|-------------------------|----------------------|
| Chloroform | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Chloromethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| cis-1,2-Dichloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| cis-1,3-Dichloropropene | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:06 |
| Dibromochloromethane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:06 |
| Dibromomethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Dichlorodifluoromethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Ethylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Freon-113 | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:06 |
| Hexachlorobutadiene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Isopropylbenzene (Cumene) | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Methylene chloride | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:06 |
| Methyl-t-butyl ether | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:06 |
| Naphthalene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| n-Butylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| n-Propylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| o-Xylene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| P & M -Xylene | 2.00 U | 2.00 | 0.620 | ug/L | 1 | | 09/15/21 15:06 |
| sec-Butylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Styrene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| tert-Butylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Tetrachloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Toluene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| trans-1,2-Dichloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| trans-1,3-Dichloropropene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Trichloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Trichlorofluoromethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:06 |
| Vinyl acetate | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:06 |
| Vinyl chloride | 0.150 U | 0.150 | 0.0500 | ug/L | 1 | | 09/15/21 15:06 |
| Xylenes (total) | 3.00 U | 3.00 | 1.00 | ug/L | 1 | | 09/15/21 15:06 |
| Surrogates | | | | | | | |
| 1,2-Dichloroethane-D4 (surr) | 118 | 81-118 | | % | 1 | | 09/15/21 15:06 |
| 4-Bromofluorobenzene (surr) | 96.4 | 85-114 | | % | 1 | | 09/15/21 15:06 |
| Toluene-d8 (surr) | 99.5 | 89-112 | | % | 1 | | 09/15/21 15:06 |

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Results of 18190-MW1-091021

Client Sample ID: **18190-MW1-091021**
Client Project ID: **18190 Essential 1**
Lab Sample ID: 1215973001
Lab Project ID: 1215973

Collection Date: 09/10/21 17:26
Received Date: 09/13/21 14:00
Matrix: Water (Surface, Eff., Ground)
Solids (%):
Location:

Results by Volatile GC/MS

Batch Information

Analytical Batch: VMS21184
Analytical Method: SW8260D
Analyst: NRB
Analytical Date/Time: 09/15/21 15:06
Container ID: 1215973001-A

Prep Batch: VXX37856
Prep Method: SW5030B
Prep Date/Time: 09/15/21 09:00
Prep Initial Wt./Vol.: 5 mL
Prep Extract Vol: 5 mL



Results of 18190-MW77-091021

Client Sample ID: 18190-MW77-091021
Client Project ID: 18190 Essential 1
Lab Sample ID: 1215973002
Lab Project ID: 1215973

Collection Date: 09/10/21 17:36
Received Date: 09/13/21 14:00
Matrix: Water (Surface, Eff., Ground)
Solids (%):
Location:

Results by Polynuclear Aromatics GC/MS

Table with 8 columns: Parameter, Result Qual, LOQ/CL, DL, Units, DF, Allowable Limits, Date Analyzed. Lists various polynuclear aromatic hydrocarbons and their surrogate compounds with associated quality and detection data.

Batch Information

Analytical Batch: XMS12899
Analytical Method: 8270D SIM LV (PAH)
Analyst: LAW
Analytical Date/Time: 09/20/21 23:48
Container ID: 1215973002-F

Prep Batch: XXX45561
Prep Method: SW3535A
Prep Date/Time: 09/14/21 12:30
Prep Initial Wt./Vol.: 245 mL
Prep Extract Vol: 1 mL



Results of **18190-MW77-091021**

Client Sample ID: **18190-MW77-091021**
Client Project ID: **18190 Essential 1**
Lab Sample ID: 1215973002
Lab Project ID: 1215973

Collection Date: 09/10/21 17:36
Received Date: 09/13/21 14:00
Matrix: Water (Surface, Eff., Ground)
Solids (%):
Location:

Results by **Semivolatile Organic Fuels**

| <u>Parameter</u> | <u>Result Qual</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> | <u>DF</u> | <u>Allowable Limits</u> | <u>Date Analyzed</u> |
|-----------------------|--------------------|---------------|-----------|--------------|-----------|-------------------------|----------------------|
| Diesel Range Organics | 0.588 U | 0.588 | 0.196 | mg/L | 1 | | 09/22/21 20:28 |
| Surrogates | | | | | | | |
| 5a Androstane (surr) | 80.1 | 50-150 | | % | 1 | | 09/22/21 20:28 |

Batch Information

Analytical Batch: XFC16086
Analytical Method: AK102
Analyst: JMG
Analytical Date/Time: 09/22/21 20:28
Container ID: 1215973002-D

Prep Batch: XXX45584
Prep Method: SW3520C
Prep Date/Time: 09/18/21 15:55
Prep Initial Wt./Vol.: 255 mL
Prep Extract Vol: 1 mL

Print Date: 09/30/2021 2:29:29PM



Results of 18190-MW77-091021

Client Sample ID: 18190-MW77-091021
 Client Project ID: 18190 Essential 1
 Lab Sample ID: 1215973002
 Lab Project ID: 1215973

Collection Date: 09/10/21 17:36
 Received Date: 09/13/21 14:00
 Matrix: Water (Surface, Eff., Ground)
 Solids (%):
 Location:

Results by Volatile GC/MS

| <u>Parameter</u> | <u>Result Qual</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> | <u>DF</u> | <u>Allowable Limits</u> | <u>Date Analyzed</u> |
|-----------------------------|--------------------|---------------|-----------|--------------|-----------|-------------------------|----------------------|
| 1,1,1,2-Tetrachloroethane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:21 |
| 1,1,1-Trichloroethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| 1,1,2,2-Tetrachloroethane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/21/21 17:57 |
| 1,1,2-Trichloroethane | 0.400 U | 0.400 | 0.120 | ug/L | 1 | | 09/15/21 15:21 |
| 1,1-Dichloroethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| 1,1-Dichloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| 1,1-Dichloropropene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| 1,2,3-Trichlorobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| 1,2,3-Trichloropropane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| 1,2,4-Trichlorobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| 1,2,4-Trimethylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| 1,2-Dibromo-3-chloropropane | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/21/21 17:57 |
| 1,2-Dibromoethane | 0.0750 U | 0.0750 | 0.0180 | ug/L | 1 | | 09/15/21 15:21 |
| 1,2-Dichlorobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| 1,2-Dichloroethane | 0.500 U | 0.500 | 0.200 | ug/L | 1 | | 09/15/21 15:21 |
| 1,2-Dichloropropane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| 1,3,5-Trimethylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| 1,3-Dichlorobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| 1,3-Dichloropropane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:21 |
| 1,4-Dichlorobenzene | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/21/21 17:57 |
| 2,2-Dichloropropane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| 2-Butanone (MEK) | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:21 |
| 2-Chlorotoluene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| 2-Hexanone | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:21 |
| 4-Chlorotoluene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| 4-Isopropyltoluene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| 4-Methyl-2-pentanone (MIBK) | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:21 |
| Benzene | 0.400 U | 0.400 | 0.120 | ug/L | 1 | | 09/15/21 15:21 |
| Bromobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| Bromochloromethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| Bromodichloromethane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:21 |
| Bromoform | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| Bromomethane | 6.00 U | 6.00 | 3.00 | ug/L | 1 | | 09/15/21 15:21 |
| Carbon disulfide | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:21 |
| Carbon tetrachloride | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| Chlorobenzene | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:21 |
| Chloroethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |

Print Date: 09/30/2021 2:29:29PM



Results of 18190-MW77-091021

Client Sample ID: 18190-MW77-091021
 Client Project ID: 18190 Essential 1
 Lab Sample ID: 1215973002
 Lab Project ID: 1215973

Collection Date: 09/10/21 17:36
 Received Date: 09/13/21 14:00
 Matrix: Water (Surface, Eff., Ground)
 Solids (%):
 Location:

Results by Volatile GC/MS

| <u>Parameter</u> | <u>Result Qual</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> | <u>DF</u> | <u>Allowable Limits</u> | <u>Date Analyzed</u> |
|------------------------------|--------------------|---------------|-----------|--------------|-----------|-------------------------|----------------------|
| Chloroform | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| Chloromethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| cis-1,2-Dichloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| cis-1,3-Dichloropropene | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:21 |
| Dibromochloromethane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 15:21 |
| Dibromomethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| Dichlorodifluoromethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| Ethylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| Freon-113 | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:21 |
| Hexachlorobutadiene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| Isopropylbenzene (Cumene) | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| Methylene chloride | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:21 |
| Methyl-t-butyl ether | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:21 |
| Naphthalene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| n-Butylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| n-Propylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| o-Xylene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| P & M -Xylene | 2.00 U | 2.00 | 0.620 | ug/L | 1 | | 09/15/21 15:21 |
| sec-Butylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| Styrene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| tert-Butylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/21/21 17:57 |
| Tetrachloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| Toluene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| trans-1,2-Dichloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| trans-1,3-Dichloropropene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| Trichloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| Trichlorofluoromethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 15:21 |
| Vinyl acetate | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 15:21 |
| Vinyl chloride | 0.150 U | 0.150 | 0.0500 | ug/L | 1 | | 09/15/21 15:21 |
| Xylenes (total) | 3.00 U | 3.00 | 1.00 | ug/L | 1 | | 09/15/21 15:21 |
| Surrogates | | | | | | | |
| 1,2-Dichloroethane-D4 (surr) | 105 | 81-118 | | % | 1 | | 09/15/21 15:21 |
| 4-Bromofluorobenzene (surr) | 112 | 85-114 | | % | 1 | | 09/21/21 17:57 |
| Toluene-d8 (surr) | 95.5 | 89-112 | | % | 1 | | 09/15/21 15:21 |

Print Date: 09/30/2021 2:29:29PM

Results of 18190-MW77-091021

Client Sample ID: **18190-MW77-091021**
Client Project ID: **18190 Essential 1**
Lab Sample ID: 1215973002
Lab Project ID: 1215973

Collection Date: 09/10/21 17:36
Received Date: 09/13/21 14:00
Matrix: Water (Surface, Eff., Ground)
Solids (%):
Location:

Results by Volatile GC/MS

Batch Information

Analytical Batch: VMS21184
Analytical Method: SW8260D
Analyst: NRB
Analytical Date/Time: 09/15/21 15:21
Container ID: 1215973002-A

Prep Batch: VXX37856
Prep Method: SW5030B
Prep Date/Time: 09/15/21 09:00
Prep Initial Wt./Vol.: 5 mL
Prep Extract Vol: 5 mL

Analytical Batch: VMS21201
Analytical Method: SW8260D
Analyst: MDT
Analytical Date/Time: 09/21/21 17:57
Container ID: 1215973002-B

Prep Batch: VXX37888
Prep Method: SW5030B
Prep Date/Time: 09/21/21 06:00
Prep Initial Wt./Vol.: 5 mL
Prep Extract Vol: 5 mL

Print Date: 09/30/2021 2:29:29PM



Results of Trip Blank

Client Sample ID: **Trip Blank**
 Client Project ID: **18190 Essential 1**
 Lab Sample ID: 1215973003
 Lab Project ID: 1215973

Collection Date: 09/10/21 17:26
 Received Date: 09/13/21 14:00
 Matrix: Water (Surface, Eff., Ground)
 Solids (%):
 Location:

Results by Volatile GC/MS

| <u>Parameter</u> | <u>Result Qual</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> | <u>DF</u> | <u>Allowable Limits</u> | <u>Date Analyzed</u> |
|-----------------------------|--------------------|---------------|-----------|--------------|-----------|-------------------------|----------------------|
| 1,1,1,2-Tetrachloroethane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 13:07 |
| 1,1,1-Trichloroethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 1,1,2,2-Tetrachloroethane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 13:07 |
| 1,1,2-Trichloroethane | 0.400 U | 0.400 | 0.120 | ug/L | 1 | | 09/15/21 13:07 |
| 1,1-Dichloroethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 1,1-Dichloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 1,1-Dichloropropene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 1,2,3-Trichlorobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 1,2,3-Trichloropropane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 1,2,4-Trichlorobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 1,2,4-Trimethylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 1,2-Dibromo-3-chloropropane | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 13:07 |
| 1,2-Dibromoethane | 0.0750 U | 0.0750 | 0.0180 | ug/L | 1 | | 09/15/21 13:07 |
| 1,2-Dichlorobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 1,2-Dichloroethane | 0.500 U | 0.500 | 0.200 | ug/L | 1 | | 09/15/21 13:07 |
| 1,2-Dichloropropane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 1,3,5-Trimethylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 1,3-Dichlorobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 1,3-Dichloropropane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 13:07 |
| 1,4-Dichlorobenzene | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 13:07 |
| 2,2-Dichloropropane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 2-Butanone (MEK) | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 13:07 |
| 2-Chlorotoluene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 2-Hexanone | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 13:07 |
| 4-Chlorotoluene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 4-Isopropyltoluene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| 4-Methyl-2-pentanone (MIBK) | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 13:07 |
| Benzene | 0.400 U | 0.400 | 0.120 | ug/L | 1 | | 09/15/21 13:07 |
| Bromobenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Bromochloromethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Bromodichloromethane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 13:07 |
| Bromoform | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Bromomethane | 6.00 U | 6.00 | 3.00 | ug/L | 1 | | 09/15/21 13:07 |
| Carbon disulfide | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 13:07 |
| Carbon tetrachloride | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Chlorobenzene | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 13:07 |
| Chloroethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |

Print Date: 09/30/2021 2:29:29PM



Results of Trip Blank

Client Sample ID: **Trip Blank**
 Client Project ID: **18190 Essential 1**
 Lab Sample ID: 1215973003
 Lab Project ID: 1215973

Collection Date: 09/10/21 17:26
 Received Date: 09/13/21 14:00
 Matrix: Water (Surface, Eff., Ground)
 Solids (%):
 Location:

Results by Volatile GC/MS

| <u>Parameter</u> | <u>Result Qual</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> | <u>DF</u> | <u>Allowable Limits</u> | <u>Date Analyzed</u> |
|------------------------------|--------------------|---------------|-----------|--------------|-----------|-------------------------|----------------------|
| Chloroform | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Chloromethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| cis-1,2-Dichloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| cis-1,3-Dichloropropene | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 13:07 |
| Dibromochloromethane | 0.500 U | 0.500 | 0.150 | ug/L | 1 | | 09/15/21 13:07 |
| Dibromomethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Dichlorodifluoromethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Ethylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Freon-113 | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 13:07 |
| Hexachlorobutadiene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Isopropylbenzene (Cumene) | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Methylene chloride | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 13:07 |
| Methyl-t-butyl ether | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 13:07 |
| Naphthalene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| n-Butylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| n-Propylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| o-Xylene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| P & M -Xylene | 2.00 U | 2.00 | 0.620 | ug/L | 1 | | 09/15/21 13:07 |
| sec-Butylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Styrene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| tert-Butylbenzene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Tetrachloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Toluene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| trans-1,2-Dichloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| trans-1,3-Dichloropropene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Trichloroethene | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Trichlorofluoromethane | 1.00 U | 1.00 | 0.310 | ug/L | 1 | | 09/15/21 13:07 |
| Vinyl acetate | 10.0 U | 10.0 | 3.10 | ug/L | 1 | | 09/15/21 13:07 |
| Vinyl chloride | 0.150 U | 0.150 | 0.0500 | ug/L | 1 | | 09/15/21 13:07 |
| Xylenes (total) | 3.00 U | 3.00 | 1.00 | ug/L | 1 | | 09/15/21 13:07 |
| Surrogates | | | | | | | |
| 1,2-Dichloroethane-D4 (surr) | 107 | 81-118 | | % | 1 | | 09/15/21 13:07 |
| 4-Bromofluorobenzene (surr) | 103 | 85-114 | | % | 1 | | 09/15/21 13:07 |
| Toluene-d8 (surr) | 99.1 | 89-112 | | % | 1 | | 09/15/21 13:07 |

Print Date: 09/30/2021 2:29:29PM

Results of Trip Blank

Client Sample ID: **Trip Blank**
Client Project ID: **18190 Essential 1**
Lab Sample ID: 1215973003
Lab Project ID: 1215973

Collection Date: 09/10/21 17:26
Received Date: 09/13/21 14:00
Matrix: Water (Surface, Eff., Ground)
Solids (%):
Location:

Results by Volatile GC/MS

Batch Information

Analytical Batch: VMS21184
Analytical Method: SW8260D
Analyst: NRB
Analytical Date/Time: 09/15/21 13:07
Container ID: 1215973003-A

Prep Batch: VXX37856
Prep Method: SW5030B
Prep Date/Time: 09/15/21 09:00
Prep Initial Wt./Vol.: 5 mL
Prep Extract Vol: 5 mL

Print Date: 09/30/2021 2:29:29PM



Method Blank

Blank ID: MB for HBN 1825633 [VXX/37856]

Blank Lab ID: 1636751

QC for Samples:

1215973001, 1215973002, 1215973003

Matrix: Water (Surface, Eff., Ground)

Results by SW8260D

| <u>Parameter</u> | <u>Results</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> |
|-----------------------------|----------------|---------------|-----------|--------------|
| 1,1,1,2-Tetrachloroethane | 0.250U | 0.500 | 0.150 | ug/L |
| 1,1,1-Trichloroethane | 0.500U | 1.00 | 0.310 | ug/L |
| 1,1,2,2-Tetrachloroethane | 0.250U | 0.500 | 0.150 | ug/L |
| 1,1,2-Trichloroethane | 0.200U | 0.400 | 0.120 | ug/L |
| 1,1-Dichloroethane | 0.500U | 1.00 | 0.310 | ug/L |
| 1,1-Dichloroethene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,1-Dichloropropene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,2,3-Trichlorobenzene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,2,3-Trichloropropane | 0.500U | 1.00 | 0.310 | ug/L |
| 1,2,4-Trichlorobenzene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,2,4-Trimethylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,2-Dibromo-3-chloropropane | 5.00U | 10.0 | 3.10 | ug/L |
| 1,2-Dibromoethane | 0.0375U | 0.0750 | 0.0180 | ug/L |
| 1,2-Dichlorobenzene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,2-Dichloroethane | 0.250U | 0.500 | 0.200 | ug/L |
| 1,2-Dichloropropane | 0.500U | 1.00 | 0.310 | ug/L |
| 1,3,5-Trimethylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,3-Dichlorobenzene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,3-Dichloropropane | 0.250U | 0.500 | 0.150 | ug/L |
| 1,4-Dichlorobenzene | 0.250U | 0.500 | 0.150 | ug/L |
| 2,2-Dichloropropane | 0.500U | 1.00 | 0.310 | ug/L |
| 2-Butanone (MEK) | 5.00U | 10.0 | 3.10 | ug/L |
| 2-Chlorotoluene | 0.500U | 1.00 | 0.310 | ug/L |
| 2-Hexanone | 5.00U | 10.0 | 3.10 | ug/L |
| 4-Chlorotoluene | 0.500U | 1.00 | 0.310 | ug/L |
| 4-Isopropyltoluene | 0.500U | 1.00 | 0.310 | ug/L |
| 4-Methyl-2-pentanone (MIBK) | 5.00U | 10.0 | 3.10 | ug/L |
| Benzene | 0.200U | 0.400 | 0.120 | ug/L |
| Bromobenzene | 0.500U | 1.00 | 0.310 | ug/L |
| Bromochloromethane | 0.500U | 1.00 | 0.310 | ug/L |
| Bromodichloromethane | 0.250U | 0.500 | 0.150 | ug/L |
| Bromoform | 0.500U | 1.00 | 0.310 | ug/L |
| Bromomethane | 3.00U | 6.00 | 3.00 | ug/L |
| Carbon disulfide | 5.00U | 10.0 | 3.10 | ug/L |
| Carbon tetrachloride | 0.500U | 1.00 | 0.310 | ug/L |
| Chlorobenzene | 0.250U | 0.500 | 0.150 | ug/L |
| Chloroethane | 0.500U | 1.00 | 0.310 | ug/L |
| Chloroform | 0.500U | 1.00 | 0.310 | ug/L |

Print Date: 09/30/2021 2:29:32PM



Method Blank

Blank ID: MB for HBN 1825633 [VXX/37856]

Blank Lab ID: 1636751

QC for Samples:

1215973001, 1215973002, 1215973003

Matrix: Water (Surface, Eff., Ground)

Results by SW8260D

| <u>Parameter</u> | <u>Results</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> |
|------------------------------|----------------|---------------|-----------|--------------|
| Chloromethane | 0.500U | 1.00 | 0.310 | ug/L |
| cis-1,2-Dichloroethene | 0.500U | 1.00 | 0.310 | ug/L |
| cis-1,3-Dichloropropene | 0.250U | 0.500 | 0.150 | ug/L |
| Dibromochloromethane | 0.250U | 0.500 | 0.150 | ug/L |
| Dibromomethane | 0.500U | 1.00 | 0.310 | ug/L |
| Dichlorodifluoromethane | 0.500U | 1.00 | 0.310 | ug/L |
| Ethylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| Freon-113 | 5.00U | 10.0 | 3.10 | ug/L |
| Hexachlorobutadiene | 0.500U | 1.00 | 0.310 | ug/L |
| Isopropylbenzene (Cumene) | 0.500U | 1.00 | 0.310 | ug/L |
| Methylene chloride | 5.00U | 10.0 | 3.10 | ug/L |
| Methyl-t-butyl ether | 5.00U | 10.0 | 3.10 | ug/L |
| Naphthalene | 0.500U | 1.00 | 0.310 | ug/L |
| n-Butylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| n-Propylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| o-Xylene | 0.500U | 1.00 | 0.310 | ug/L |
| P & M -Xylene | 1.00U | 2.00 | 0.620 | ug/L |
| sec-Butylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| Styrene | 0.500U | 1.00 | 0.310 | ug/L |
| tert-Butylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| Tetrachloroethene | 0.500U | 1.00 | 0.310 | ug/L |
| Toluene | 0.500U | 1.00 | 0.310 | ug/L |
| trans-1,2-Dichloroethene | 0.500U | 1.00 | 0.310 | ug/L |
| trans-1,3-Dichloropropene | 0.500U | 1.00 | 0.310 | ug/L |
| Trichloroethene | 0.500U | 1.00 | 0.310 | ug/L |
| Trichlorofluoromethane | 0.500U | 1.00 | 0.310 | ug/L |
| Vinyl acetate | 5.00U | 10.0 | 3.10 | ug/L |
| Vinyl chloride | 0.0750U | 0.150 | 0.0500 | ug/L |
| Xylenes (total) | 1.50U | 3.00 | 1.00 | ug/L |
| Surrogates | | | | |
| 1,2-Dichloroethane-D4 (surr) | 106 | 81-118 | | % |
| 4-Bromofluorobenzene (surr) | 115* | 85-114 | | % |
| Toluene-d8 (surr) | 100 | 89-112 | | % |

Print Date: 09/30/2021 2:29:32PM



Method Blank

Blank ID: MB for HBN 1825633 [VXX/37856]
Blank Lab ID: 1636751

Matrix: Water (Surface, Eff., Ground)

QC for Samples:
1215973001, 1215973002, 1215973003

Results by SW8260D

| <u>Parameter</u> | <u>Results</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> |
|------------------|----------------|---------------|-----------|--------------|
|------------------|----------------|---------------|-----------|--------------|

Batch Information

Analytical Batch: VMS21184
Analytical Method: SW8260D
Instrument: Agilent 7890-75MS
Analyst: NRB
Analytical Date/Time: 9/15/2021 9:30:00AM

Prep Batch: VXX37856
Prep Method: SW5030B
Prep Date/Time: 9/15/2021 9:00:00AM
Prep Initial Wt./Vol.: 5 mL
Prep Extract Vol: 5 mL

Print Date: 09/30/2021 2:29:32PM



Blank Spike Summary

Blank Spike ID: LCS for HBN 1215973 [VXX37856]
 Blank Spike Lab ID: 1636752
 Date Analyzed: 09/15/2021 09:45

Spike Duplicate ID: LCSD for HBN 1215973
 [VXX37856]
 Spike Duplicate Lab ID: 1636753
 Matrix: Water (Surface, Eff., Ground)

QC for Samples: 1215973001, 1215973002, 1215973003

Results by SW8260D

| Parameter | Blank Spike (ug/L) | | | Spike Duplicate (ug/L) | | | CL | RPD (%) | RPD CL |
|-----------------------------|--------------------|--------|---------|------------------------|--------|---------|--------------|---------|-----------|
| | Spike | Result | Rec (%) | Spike | Result | Rec (%) | | | |
| 1,1,1,2-Tetrachloroethane | 30 | 26.9 | 90 | 30 | 32.0 | 107 | (78-124) | 17.10 | (< 20) |
| 1,1,1-Trichloroethane | 30 | 31.8 | 106 | 30 | 29.4 | 98 | (74-131) | 7.70 | (< 20) |
| 1,1,2,2-Tetrachloroethane | 30 | 31.8 | 106 | 30 | 32.1 | 107 | (71-121) | 0.90 | (< 20) |
| 1,1,2-Trichloroethane | 30 | 30.7 | 102 | 30 | 32.2 | 107 | (80-119) | 5.10 | (< 20) |
| 1,1-Dichloroethane | 30 | 31.0 | 103 | 30 | 29.7 | 99 | (77-125) | 4.30 | (< 20) |
| 1,1-Dichloroethene | 30 | 29.9 | 100 | 30 | 28.5 | 95 | (71-131) | 4.90 | (< 20) |
| 1,1-Dichloropropene | 30 | 29.5 | 98 | 30 | 30.1 | 100 | (79-125) | 2.00 | (< 20) |
| 1,2,3-Trichlorobenzene | 30 | 30.7 | 102 | 30 | 37.6 | 125 | (69-129) | 20.10 | * (< 20) |
| 1,2,3-Trichloropropane | 30 | 32.8 | 109 | 30 | 35.5 | 118 | (73-122) | 7.90 | (< 20) |
| 1,2,4-Trichlorobenzene | 30 | 29.0 | 97 | 30 | 34.0 | 113 | (69-130) | 16.00 | (< 20) |
| 1,2,4-Trimethylbenzene | 30 | 32.1 | 107 | 30 | 31.7 | 106 | (79-124) | 1.50 | (< 20) |
| 1,2-Dibromo-3-chloropropane | 30 | 31.7 | 106 | 30 | 30.8 | 103 | (62-128) | 2.80 | (< 20) |
| 1,2-Dibromoethane | 30 | 27.8 | 93 | 30 | 29.2 | 97 | (77-121) | 4.70 | (< 20) |
| 1,2-Dichlorobenzene | 30 | 30.2 | 101 | 30 | 30.8 | 103 | (80-119) | 2.00 | (< 20) |
| 1,2-Dichloroethane | 30 | 30.6 | 102 | 30 | 30.6 | 102 | (73-128) | 0.06 | (< 20) |
| 1,2-Dichloropropane | 30 | 30.4 | 101 | 30 | 25.7 | 86 | (78-122) | 16.60 | (< 20) |
| 1,3,5-Trimethylbenzene | 30 | 32.3 | 108 | 30 | 35.3 | 118 | (75-124) | 8.70 | (< 20) |
| 1,3-Dichlorobenzene | 30 | 29.1 | 97 | 30 | 30.4 | 101 | (80-119) | 4.60 | (< 20) |
| 1,3-Dichloropropane | 30 | 27.9 | 93 | 30 | 29.4 | 98 | (80-119) | 5.30 | (< 20) |
| 1,4-Dichlorobenzene | 30 | 28.8 | 96 | 30 | 30.1 | 100 | (79-118) | 4.40 | (< 20) |
| 2,2-Dichloropropane | 30 | 31.0 | 103 | 30 | 29.8 | 99 | (60-139) | 3.80 | (< 20) |
| 2-Butanone (MEK) | 90 | 82.1 | 91 | 90 | 77.0 | 86 | (56-143) | 6.40 | (< 20) |
| 2-Chlorotoluene | 30 | 33.9 | 113 | 30 | 38.3 | 128 | * (79-122) | 12.10 | (< 20) |
| 2-Hexanone | 90 | 83.8 | 93 | 90 | 82.9 | 92 | (57-139) | 1.00 | (< 20) |
| 4-Chlorotoluene | 30 | 34.9 | 116 | 30 | 37.7 | 126 | * (78-122) | 7.60 | (< 20) |
| 4-Isopropyltoluene | 30 | 28.2 | 94 | 30 | 29.5 | 99 | (77-127) | 4.80 | (< 20) |
| 4-Methyl-2-pentanone (MIBK) | 90 | 85.1 | 95 | 90 | 75.5 | 84 | (67-130) | 11.90 | (< 20) |
| Benzene | 30 | 30.3 | 101 | 30 | 29.1 | 97 | (79-120) | 4.10 | (< 20) |
| Bromobenzene | 30 | 31.1 | 104 | 30 | 35.1 | 117 | (80-120) | 12.00 | (< 20) |
| Bromochloromethane | 30 | 27.1 | 90 | 30 | 27.2 | 91 | (78-123) | 0.37 | (< 20) |
| Bromodichloromethane | 30 | 31.0 | 103 | 30 | 27.5 | 92 | (79-125) | 11.80 | (< 20) |
| Bromoform | 30 | 26.0 | 87 | 30 | 30.6 | 102 | (66-130) | 16.20 | (< 20) |
| Bromomethane | 30 | 33.5 | 112 | 30 | 34.5 | 115 | (53-141) | 2.80 | (< 20) |
| Carbon disulfide | 45 | 46.1 | 102 | 45 | 43.4 | 97 | (64-133) | 5.90 | (< 20) |

Print Date: 09/30/2021 2:29:34PM

Blank Spike Summary

Blank Spike ID: LCS for HBN 1215973 [VXX37856]
 Blank Spike Lab ID: 1636752
 Date Analyzed: 09/15/2021 09:45

Spike Duplicate ID: LCSD for HBN 1215973 [VXX37856]
 Spike Duplicate Lab ID: 1636753
 Matrix: Water (Surface, Eff., Ground)

QC for Samples: 1215973001, 1215973002, 1215973003

Results by SW8260D

| Parameter | Blank Spike (ug/L) | | | Spike Duplicate (ug/L) | | | CL | RPD (%) | RPD CL |
|---------------------------|--------------------|--------|---------|------------------------|--------|---------|--------------|---------|-----------|
| | Spike | Result | Rec (%) | Spike | Result | Rec (%) | | | |
| Carbon tetrachloride | 30 | 31.2 | 104 | 30 | 30.7 | 102 | (72-136) | 1.60 | (< 20) |
| Chlorobenzene | 30 | 27.0 | 90 | 30 | 31.2 | 104 | (82-118) | 14.50 | (< 20) |
| Chloroethane | 30 | 35.0 | 117 | 30 | 34.3 | 114 | (60-138) | 2.20 | (< 20) |
| Chloroform | 30 | 30.0 | 100 | 30 | 28.1 | 94 | (79-124) | 6.30 | (< 20) |
| Chloromethane | 30 | 25.1 | 84 | 30 | 24.2 | 81 | (50-139) | 3.50 | (< 20) |
| cis-1,2-Dichloroethene | 30 | 30.0 | 100 | 30 | 29.6 | 99 | (78-123) | 1.10 | (< 20) |
| cis-1,3-Dichloropropene | 30 | 30.3 | 101 | 30 | 29.0 | 97 | (75-124) | 4.40 | (< 20) |
| Dibromochloromethane | 30 | 29.9 | 100 | 30 | 31.7 | 106 | (74-126) | 5.90 | (< 20) |
| Dibromomethane | 30 | 28.7 | 96 | 30 | 29.6 | 99 | (79-123) | 3.20 | (< 20) |
| Dichlorodifluoromethane | 30 | 31.5 | 105 | 30 | 30.0 | 100 | (32-152) | 5.00 | (< 20) |
| Ethylbenzene | 30 | 27.9 | 93 | 30 | 32.7 | 109 | (79-121) | 15.70 | (< 20) |
| Freon-113 | 45 | 46.6 | 104 | 45 | 44.4 | 99 | (70-136) | 5.00 | (< 20) |
| Hexachlorobutadiene | 30 | 28.1 | 94 | 30 | 32.1 | 107 | (66-134) | 13.30 | (< 20) |
| Isopropylbenzene (Cumene) | 30 | 29.6 | 99 | 30 | 32.5 | 108 | (72-131) | 9.60 | (< 20) |
| Methylene chloride | 30 | 28.6 | 96 | 30 | 28.1 | 94 | (74-124) | 2.10 | (< 20) |
| Methyl-t-butyl ether | 45 | 47.7 | 106 | 45 | 46.9 | 104 | (71-124) | 1.70 | (< 20) |
| Naphthalene | 30 | 25.9 | 86 | 30 | 33.4 | 111 | (61-128) | 25.40 | * (< 20) |
| n-Butylbenzene | 30 | 29.5 | 98 | 30 | 30.5 | 102 | (75-128) | 3.60 | (< 20) |
| n-Propylbenzene | 30 | 35.8 | 119 | 30 | 39.1 | 130 | * (76-126) | 8.70 | (< 20) |
| o-Xylene | 30 | 28.3 | 94 | 30 | 33.6 | 112 | (78-122) | 17.20 | (< 20) |
| P & M -Xylene | 60 | 58.7 | 98 | 60 | 69.6 | 116 | (80-121) | 17.00 | (< 20) |
| sec-Butylbenzene | 30 | 34.5 | 115 | 30 | 31.8 | 106 | (77-126) | 8.20 | (< 20) |
| Styrene | 30 | 27.1 | 90 | 30 | 32.3 | 108 | (78-123) | 17.70 | (< 20) |
| tert-Butylbenzene | 30 | 34.5 | 115 | 30 | 36.3 | 121 | (78-124) | 5.00 | (< 20) |
| Tetrachloroethene | 30 | 27.9 | 93 | 30 | 29.3 | 98 | (74-129) | 5.00 | (< 20) |
| Toluene | 30 | 26.6 | 89 | 30 | 32.4 | 108 | (80-121) | 19.80 | (< 20) |
| trans-1,2-Dichloroethene | 30 | 30.4 | 101 | 30 | 29.3 | 98 | (75-124) | 3.80 | (< 20) |
| trans-1,3-Dichloropropene | 30 | 31.7 | 106 | 30 | 32.6 | 109 | (73-127) | 2.70 | (< 20) |
| Trichloroethene | 30 | 27.6 | 92 | 30 | 29.3 | 98 | (79-123) | 6.10 | (< 20) |
| Trichlorofluoromethane | 30 | 34.1 | 114 | 30 | 32.9 | 110 | (65-141) | 3.40 | (< 20) |
| Vinyl acetate | 30 | 32.0 | 107 | 30 | 32.0 | 107 | (54-146) | 0.08 | (< 20) |
| Vinyl chloride | 30 | 31.1 | 104 | 30 | 29.5 | 99 | (58-137) | 5.10 | (< 20) |
| Xylenes (total) | 90 | 87.0 | 97 | 90 | 103 | 115 | (79-121) | 17.00 | (< 20) |

Blank Spike Summary

Blank Spike ID: LCS for HBN 1215973 [VXX37856]
 Blank Spike Lab ID: 1636752
 Date Analyzed: 09/15/2021 09:45

Spike Duplicate ID: LCSD for HBN 1215973 [VXX37856]
 Spike Duplicate Lab ID: 1636753
 Matrix: Water (Surface, Eff., Ground)

QC for Samples: 1215973001, 1215973002, 1215973003

Results by SW8260D

| Parameter | Blank Spike (%) | | | Spike Duplicate (%) | | | CL | RPD (%) | RPD CL |
|------------------------------|-----------------|--------|---------|---------------------|--------|---------|------------|---------|--------|
| | Spike | Result | Rec (%) | Spike | Result | Rec (%) | | | |
| Surrogates | | | | | | | | | |
| 1,2-Dichloroethane-D4 (surr) | 30 | | 104 | 30 | | 106 | (81-118) | 1.90 | |
| 4-Bromofluorobenzene (surr) | 30 | | 107 | 30 | | 111 | (85-114) | 3.50 | |
| Toluene-d8 (surr) | 30 | | 99 | 30 | | 107 | (89-112) | 8.40 | |

Batch Information

Analytical Batch: **VMS21184**
 Analytical Method: **SW8260D**
 Instrument: **Agilent 7890-75MS**
 Analyst: **NRB**

Prep Batch: **VXX37856**
 Prep Method: **SW5030B**
 Prep Date/Time: **09/15/2021 09:00**
 Spike Init Wt./Vol.: 30 ug/L Extract Vol: 5 mL
 Dupe Init Wt./Vol.: 30 ug/L Extract Vol: 5 mL

Method Blank

Blank ID: MB for HBN 1825978 [VXX/37888]
 Blank Lab ID: 1637560

Matrix: Water (Surface, Eff., Ground)

QC for Samples:
 1215973002

Results by SW8260D

| <u>Parameter</u> | <u>Results</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> |
|------------------------------|----------------|---------------|-----------|--------------|
| 1,1,2,2-Tetrachloroethane | 0.250U | 0.500 | 0.150 | ug/L |
| 1,2,3-Trichlorobenzene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,2,3-Trichloropropane | 0.500U | 1.00 | 0.310 | ug/L |
| 1,2,4-Trichlorobenzene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,2,4-Trimethylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,2-Dibromo-3-chloropropane | 5.00U | 10.0 | 3.10 | ug/L |
| 1,2-Dichlorobenzene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,3,5-Trimethylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,3-Dichlorobenzene | 0.500U | 1.00 | 0.310 | ug/L |
| 1,4-Dichlorobenzene | 0.250U | 0.500 | 0.150 | ug/L |
| 2-Chlorotoluene | 0.500U | 1.00 | 0.310 | ug/L |
| 4-Chlorotoluene | 0.500U | 1.00 | 0.310 | ug/L |
| 4-Isopropyltoluene | 0.500U | 1.00 | 0.310 | ug/L |
| Bromobenzene | 0.500U | 1.00 | 0.310 | ug/L |
| Hexachlorobutadiene | 0.500U | 1.00 | 0.310 | ug/L |
| Naphthalene | 0.500U | 1.00 | 0.310 | ug/L |
| n-Butylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| n-Propylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| sec-Butylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| tert-Butylbenzene | 0.500U | 1.00 | 0.310 | ug/L |
| Surrogates | | | | |
| 1,2-Dichloroethane-D4 (surr) | 106 | 81-118 | | % |
| 4-Bromofluorobenzene (surr) | 106 | 85-114 | | % |
| Toluene-d8 (surr) | 114* | 89-112 | | % |

Batch Information

Analytical Batch: VMS21201
 Analytical Method: SW8260D
 Instrument: Agilent 7890-75MS
 Analyst: MDT
 Analytical Date/Time: 9/21/2021 11:33:00AM

Prep Batch: VXX37888
 Prep Method: SW5030B
 Prep Date/Time: 9/21/2021 6:00:00AM
 Prep Initial Wt./Vol.: 5 mL
 Prep Extract Vol: 5 mL

Blank Spike Summary

Blank Spike ID: LCS for HBN 1215973 [VXX37888]
 Blank Spike Lab ID: 1637561
 Date Analyzed: 09/21/2021 11:48

Spike Duplicate ID: LCSD for HBN 1215973 [VXX37888]
 Spike Duplicate Lab ID: 1637562
 Matrix: Water (Surface, Eff., Ground)

QC for Samples: 1215973002

Results by SW8260D

| Parameter | Blank Spike (ug/L) | | | Spike Duplicate (ug/L) | | | CL | RPD (%) | RPD CL |
|-----------------------------|--------------------|--------|---------|------------------------|--------|---------|------------|---------|---------|
| | Spike | Result | Rec (%) | Spike | Result | Rec (%) | | | |
| 1,1,2,2-Tetrachloroethane | 30 | 29.9 | 100 | 30 | 30.3 | 101 | (71-121) | 1.40 | (< 20) |
| 1,2,3-Trichlorobenzene | 30 | 29.2 | 97 | 30 | 30.4 | 101 | (69-129) | 4.20 | (< 20) |
| 1,2,3-Trichloropropane | 30 | 31.2 | 104 | 30 | 31.7 | 106 | (73-122) | 1.60 | (< 20) |
| 1,2,4-Trichlorobenzene | 30 | 28.6 | 95 | 30 | 29.6 | 99 | (69-130) | 3.40 | (< 20) |
| 1,2,4-Trimethylbenzene | 30 | 32.3 | 108 | 30 | 32.5 | 108 | (79-124) | 0.80 | (< 20) |
| 1,2-Dibromo-3-chloropropane | 30 | 30.0 | 100 | 30 | 30.4 | 101 | (62-128) | 1.30 | (< 20) |
| 1,2-Dichlorobenzene | 30 | 30.0 | 100 | 30 | 30.5 | 102 | (80-119) | 1.80 | (< 20) |
| 1,3,5-Trimethylbenzene | 30 | 30.9 | 103 | 30 | 31.4 | 105 | (75-124) | 1.50 | (< 20) |
| 1,3-Dichlorobenzene | 30 | 30.0 | 100 | 30 | 30.5 | 102 | (80-119) | 1.80 | (< 20) |
| 1,4-Dichlorobenzene | 30 | 29.8 | 99 | 30 | 30.3 | 101 | (79-118) | 1.70 | (< 20) |
| 2-Chlorotoluene | 30 | 32.5 | 108 | 30 | 32.2 | 107 | (79-122) | 1.10 | (< 20) |
| 4-Chlorotoluene | 30 | 32.1 | 107 | 30 | 32.5 | 108 | (78-122) | 1.10 | (< 20) |
| 4-Isopropyltoluene | 30 | 29.7 | 99 | 30 | 30.3 | 101 | (77-127) | 2.00 | (< 20) |
| Bromobenzene | 30 | 29.6 | 99 | 30 | 30.1 | 100 | (80-120) | 1.80 | (< 20) |
| Hexachlorobutadiene | 30 | 27.4 | 92 | 30 | 27.8 | 93 | (66-134) | 1.20 | (< 20) |
| Naphthalene | 30 | 27.7 | 92 | 30 | 28.9 | 96 | (61-128) | 4.10 | (< 20) |
| n-Butylbenzene | 30 | 29.6 | 99 | 30 | 30.1 | 100 | (75-128) | 1.40 | (< 20) |
| n-Propylbenzene | 30 | 34.1 | 114 | 30 | 34.1 | 114 | (76-126) | 0.06 | (< 20) |
| sec-Butylbenzene | 30 | 31.8 | 106 | 30 | 32.2 | 107 | (77-126) | 1.30 | (< 20) |
| tert-Butylbenzene | 30 | 32.1 | 107 | 30 | 32.6 | 109 | (78-124) | 1.70 | (< 20) |

Surrogates

| | | | | | | | | | |
|------------------------------|----|--|-----|----|--|-----|------------|------|--|
| 1,2-Dichloroethane-D4 (surr) | 30 | | 111 | 30 | | 112 | (81-118) | 0.21 | |
| 4-Bromofluorobenzene (surr) | 30 | | 104 | 30 | | 105 | (85-114) | 0.67 | |
| Toluene-d8 (surr) | 30 | | 100 | 30 | | 100 | (89-112) | 0.33 | |

Batch Information

Analytical Batch: **VMS21201**
 Analytical Method: **SW8260D**
 Instrument: **Agilent 7890-75MS**
 Analyst: **MDT**

Prep Batch: **VXX37888**
 Prep Method: **SW5030B**
 Prep Date/Time: **09/21/2021 06:00**
 Spike Init Wt./Vol.: 30 ug/L Extract Vol: 5 mL
 Dupe Init Wt./Vol.: 30 ug/L Extract Vol: 5 mL

Method Blank

Blank ID: MB for HBN 1825521 [XXX/45561]

Blank Lab ID: 1636317

QC for Samples:

1215973001, 1215973002

Matrix: Water (Surface, Eff., Ground)

Results by 8270D SIM LV (PAH)

| <u>Parameter</u> | <u>Results</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> |
|--------------------------------|----------------|---------------|-----------|--------------|
| 1-Methylnaphthalene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| 2-Methylnaphthalene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Acenaphthene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Acenaphthylene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Anthracene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Benzo(a)Anthracene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Benzo[a]pyrene | 0.0100U | 0.0200 | 0.00620 | ug/L |
| Benzo[b]Fluoranthene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Benzo[g,h,i]perylene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Benzo[k]fluoranthene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Chrysene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Dibenzo[a,h]anthracene | 0.0100U | 0.0200 | 0.00620 | ug/L |
| Fluoranthene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Fluorene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Indeno[1,2,3-c,d] pyrene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Naphthalene | 0.0500U | 0.100 | 0.0310 | ug/L |
| Phenanthrene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Pyrene | 0.0250U | 0.0500 | 0.0150 | ug/L |
| Surrogates | | | | |
| 2-Methylnaphthalene-d10 (surr) | 53.8 | 42-86 | | % |
| Fluoranthene-d10 (surr) | 66.3 | 50-97 | | % |

Batch Information

Analytical Batch: XMS12899
 Analytical Method: 8270D SIM LV (PAH)
 Instrument: Agilent GC 7890B/5977A SWA
 Analyst: LAW
 Analytical Date/Time: 9/20/2021 9:03:00PM

Prep Batch: XXX45561
 Prep Method: SW3535A
 Prep Date/Time: 9/14/2021 12:30:03PM
 Prep Initial Wt./Vol.: 250 mL
 Prep Extract Vol: 1 mL

Blank Spike Summary

Blank Spike ID: LCS for HBN 1215973 [XXX45561]

Blank Spike Lab ID: 1636318

Date Analyzed: 09/20/2021 21:24

Matrix: Water (Surface, Eff., Ground)

QC for Samples: 1215973001, 1215973002

Results by 8270D SIM LV (PAH)

Blank Spike (ug/L)

| Parameter | Spike | Result | Rec (%) | CL |
|--------------------------|-------|--------|---------|------------|
| 1-Methylnaphthalene | 2 | 1.32 | 66 | (41-115) |
| 2-Methylnaphthalene | 2 | 1.24 | 62 | (39-114) |
| Acenaphthene | 2 | 1.47 | 74 | (48-114) |
| Acenaphthylene | 2 | 1.55 | 77 | (35-121) |
| Anthracene | 2 | 1.51 | 75 | (53-119) |
| Benzo(a)Anthracene | 2 | 1.46 | 73 | (59-120) |
| Benzo[a]pyrene | 2 | 1.48 | 74 | (53-120) |
| Benzo[b]Fluoranthene | 2 | 1.36 | 68 | (53-126) |
| Benzo[g,h,i]perylene | 2 | 1.68 | 84 | (44-128) |
| Benzo[k]fluoranthene | 2 | 1.61 | 81 | (54-125) |
| Chrysene | 2 | 1.50 | 75 | (57-120) |
| Dibenzo[a,h]anthracene | 2 | 1.68 | 84 | (44-131) |
| Fluoranthene | 2 | 1.45 | 72 | (58-120) |
| Fluorene | 2 | 1.53 | 77 | (50-118) |
| Indeno[1,2,3-c,d] pyrene | 2 | 1.63 | 81 | (48-130) |
| Naphthalene | 2 | 1.41 | 71 | (43-114) |
| Phenanthrene | 2 | 1.54 | 77 | (53-115) |
| Pyrene | 2 | 1.47 | 74 | (53-121) |

Surrogates

| | | | | |
|--------------------------------|---|--|----|-----------|
| 2-Methylnaphthalene-d10 (surr) | 2 | | 55 | (42-86) |
| Fluoranthene-d10 (surr) | 2 | | 64 | (50-97) |

Batch Information

Analytical Batch: XMS12899

Analytical Method: 8270D SIM LV (PAH)

Instrument: Agilent GC 7890B/5977A SWA

Analyst: LAW

Prep Batch: XXX45561

Prep Method: SW3535A

Prep Date/Time: 09/14/2021 12:30

Spike Init Wt./Vol.: 2 ug/L Extract Vol: 1 mL

Dupe Init Wt./Vol.: Extract Vol:

Matrix Spike Summary

Original Sample ID: 1215976003
 MS Sample ID: 1636319 MS
 MSD Sample ID: 1636320 MSD

Analysis Date: 09/20/2021 21:44
 Analysis Date: 09/20/2021 22:05
 Analysis Date: 09/20/2021 22:26
 Matrix: Water (Surface, Eff., Ground)

QC for Samples: 1215973001, 1215973002

Results by 8270D SIM LV (PAH)

| Parameter | Sample | Matrix Spike (ug/L) | | | Spike Duplicate (ug/L) | | | CL | RPD (%) | RPD CL |
|--------------------------------|----------|---------------------|--------|---------|------------------------|--------|---------|--------|---------|----------|
| | | Spike | Result | Rec (%) | Spike | Result | Rec (%) | | | |
| 1-Methylnaphthalene | 0.0232U | 1.82 | .874 | 48 | 1.82 | 0.937 | 52 | 41-115 | 7.00 | (< 20) |
| 2-Methylnaphthalene | 0.0232U | 1.82 | .827 | 46 | 1.82 | 0.888 | 49 | 39-114 | 7.00 | (< 20) |
| Acenaphthene | 0.0232U | 1.82 | .855 | 47 * | 1.82 | 0.958 | 53 | 48-114 | 11.30 | (< 20) |
| Acenaphthylene | 0.0232U | 1.82 | 1.05 | 58 | 1.82 | 1.09 | 60 | 35-121 | 3.20 | (< 20) |
| Anthracene | 0.0232U | 1.82 | .774 | 43 * | 1.82 | 1.03 | 57 | 53-119 | 28.10 | * (< 20) |
| Benzo(a)Anthracene | 0.0232U | 1.82 | .761 | 42 * | 1.82 | 1.22 | 67 | 59-120 | 46.10 | * (< 20) |
| Benzo(a)pyrene | 0.00925U | 1.82 | .749 | 41 * | 1.82 | 1.12 | 61 | 53-120 | 39.40 | * (< 20) |
| Benzo(b)Fluoranthene | 0.0232U | 1.82 | .693 | 38 * | 1.82 | 1.13 | 62 | 53-126 | 47.90 | * (< 20) |
| Benzo(g,h,i)perylene | 0.0232U | 1.82 | .796 | 44 * | 1.82 | 1.09 | 60 | 44-128 | 30.90 | * (< 20) |
| Benzo(k)fluoranthene | 0.0232U | 1.82 | .818 | 45 * | 1.82 | 1.19 | 65 | 54-125 | 36.70 | * (< 20) |
| Chrysene | 0.0232U | 1.82 | .801 | 44 * | 1.82 | 1.19 | 66 | 57-120 | 39.30 | * (< 20) |
| Dibenzo(a,h)anthracene | 0.00925U | 1.82 | .81 | 45 | 1.82 | 1.10 | 60 | 44-131 | 30.10 | * (< 20) |
| Fluoranthene | 0.0232U | 1.82 | .745 | 41 * | 1.82 | 1.13 | 62 | 58-120 | 40.90 | * (< 20) |
| Fluorene | 0.0232U | 1.82 | .869 | 48 * | 1.82 | 1.04 | 57 | 50-118 | 17.70 | (< 20) |
| Indeno[1,2,3-c,d] pyrene | 0.0232U | 1.82 | .767 | 42 * | 1.82 | 1.09 | 60 | 48-130 | 35.00 | * (< 20) |
| Naphthalene | 0.0463U | 1.82 | 1.12 | 62 | 1.82 | 1.14 | 63 | 43-114 | 1.70 | (< 20) |
| Phenanthrene | 0.0268J | 1.82 | .817 | 44 * | 1.82 | 1.05 | 56 | 53-115 | 25.00 | * (< 20) |
| Pyrene | 0.0167J | 1.82 | .76 | 41 * | 1.82 | 1.16 | 63 | 53-121 | 41.50 | * (< 20) |
| Surrogates | | | | | | | | | | |
| 2-Methylnaphthalene-d10 (surr) | | 1.82 | .747 | 41 * | 1.82 | 0.831 | 46 | 42-86 | 10.60 | |
| Fluoranthene-d10 (surr) | | 1.82 | .678 | 37 * | 1.82 | 1.03 | 57 | 50-97 | 41.10 | |

Batch Information

Analytical Batch: XMS12899
 Analytical Method: 8270D SIM LV (PAH)
 Instrument: Agilent GC 7890B/5977A SWA
 Analyst: LAW
 Analytical Date/Time: 9/20/2021 10:05:00PM

Prep Batch: XXX45561
 Prep Method: 3535 Solid Phase Ext for 8270 PAH SIM LV
 Prep Date/Time: 9/14/2021 12:30:00PM
 Prep Initial Wt./Vol.: 275.00mL
 Prep Extract Vol: 1.00mL

Method Blank

Blank ID: MB for HBN 1825731 [XXX/45584]

Blank Lab ID: 1636904

QC for Samples:

1215973001, 1215973002

Matrix: Water (Surface, Eff., Ground)

Results by AK102

| <u>Parameter</u> | <u>Results</u> | <u>LOQ/CL</u> | <u>DL</u> | <u>Units</u> |
|-----------------------|----------------|---------------|-----------|--------------|
| Diesel Range Organics | 0.300U | 0.600 | 0.200 | mg/L |
| Surrogates | | | | |
| 5a Androstane (surr) | 84.8 | 60-120 | | % |

Batch Information

Analytical Batch: XFC16088

Analytical Method: AK102

Instrument: Agilent 7890B F

Analyst: JMG

Analytical Date/Time: 9/24/2021 1:04:00PM

Prep Batch: XXX45584

Prep Method: SW3520C

Prep Date/Time: 9/18/2021 3:55:00PM

Prep Initial Wt./Vol.: 250 mL

Prep Extract Vol: 1 mL

Blank Spike Summary

Blank Spike ID: LCS for HBN 1215973 [XXX45584]
 Blank Spike Lab ID: 1636905
 Date Analyzed: 09/22/2021 18:10

Spike Duplicate ID: LCSD for HBN 1215973
 [XXX45584]
 Spike Duplicate Lab ID: 1636906
 Matrix: Water (Surface, Eff., Ground)

QC for Samples: 1215973001, 1215973002

Results by AK102

| Parameter | Blank Spike (mg/L) | | | Spike Duplicate (mg/L) | | | CL | RPD (%) | RPD CL |
|-----------------------|--------------------|--------|---------|------------------------|--------|---------|------------|---------|---------|
| | Spike | Result | Rec (%) | Spike | Result | Rec (%) | | | |
| Diesel Range Organics | 20 | 18.4 | 92 | 20 | 16.4 | 82 | (75-125) | 11.60 | (< 20) |

Surrogates

| | | | | | | |
|----------------------|-----|-----|-----|----|------------|------|
| 5a Androstane (surr) | 0.4 | 104 | 0.4 | 97 | (60-120) | 7.20 |
|----------------------|-----|-----|-----|----|------------|------|

Batch Information

Analytical Batch: **XFC16086**
 Analytical Method: **AK102**
 Instrument: **Agilent 7890B F**
 Analyst: **JMG**

Prep Batch: **XXX45584**
 Prep Method: **SW3520C**
 Prep Date/Time: **09/18/2021 15:55**
 Spike Init Wt./Vol.: 20 mg/L Extract Vol: 1 mL
 Dupe Init Wt./Vol.: 20 mg/L Extract Vol: 1 mL



1215973



SGS NORTH AMERICA INC. CHAIN OF CUSTODY RECORD

SGS Environmental Services
200 West Potter Road
Anchorage, AK 99518
(907) 562-2343
www.sgs.com/alaska

358307 of

Form with sections: SECTION 1 (Client: EMI, Contact: Andy Coulson), SECTION 2 (Sample Identification table with 3 rows), SECTION 3 (Preservative table), SECTION 4 (Relinquished/Received by), SECTION 5 (Temperature/Seal information). Includes handwritten signatures and dates.

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e-Sample Receipt Form

SGS Workorder #:

1215973



1 2 1 5 9 7 3

| Review Criteria | Condition (Yes, No, N/A) | Exceptions Noted below |
|--|---|---|
| Chain of Custody / Temperature Requirements | | <input checked="" type="checkbox"/> Yes Exemption permitted if sampler hand carries/delivers. |
| Were Custody Seals intact? Note # & location | <input type="checkbox"/> N/A | Absent, HD |
| COC accompanied samples? | <input checked="" type="checkbox"/> Yes | |
| DOD: Were samples received in COC corresponding coolers? | <input type="checkbox"/> N/A | |
| <input type="checkbox"/> N/A **Exemption permitted if chilled & collected <8 hours ago, or for samples where chilling is not required | | |
| Temperature blank compliant* (i.e., 0-6 °C after CF)? | <input checked="" type="checkbox"/> Yes | Cooler ID: 1 @ 3.8 °C Therm. ID: D23 |
| If samples received without a temperature blank, the "cooler temperature" will be documented instead & "COOLER TEMP" will be noted to the right. "ambient" or "chilled" will be noted if neither is available. | <input type="checkbox"/> | Cooler ID: @ °C Therm. ID: |
| | <input type="checkbox"/> | Cooler ID: @ °C Therm. ID: |
| | <input type="checkbox"/> | Cooler ID: @ °C Therm. ID: |
| | <input type="checkbox"/> | Cooler ID: @ °C Therm. ID: |
| *If >6°C, were samples collected <8 hours ago? | <input type="checkbox"/> N/A | |
| If <0°C, were sample containers ice free? | <input type="checkbox"/> N/A | |
| Note: Identify containers received at non-compliant temperature . Use form FS-0029 if more space is needed. | | |
| Holding Time / Documentation / Sample Condition Requirements | | Note: Refer to form F-083 "Sample Guide" for specific holding times. |
| Were samples received within holding time? | <input checked="" type="checkbox"/> Yes | |
| Do samples match COC** (i.e., sample IDs, dates/times collected)? | <input checked="" type="checkbox"/> Yes | |
| **Note: If times differ <1hr, record details & login per COC. ***Note: If sample information on containers differs from COC, SGS will default to COC information | | |
| Were analytical requests clear? (i.e., method is specified for analyses with multiple option for analysis (Ex: BTEX, Metals) | <input checked="" type="checkbox"/> Yes | |
| Were proper containers (type/mass/volume/preservative***) used? | <input checked="" type="checkbox"/> Yes | <input type="checkbox"/> N/A ***Exemption permitted for metals (e.g.200.8/6020B). |
| Volatile / LL-Hg Requirements | | |
| Were Trip Blanks (i.e., VOAs, LL-Hg) in cooler with samples? | <input checked="" type="checkbox"/> Yes | |
| Were all water VOA vials free of headspace (i.e., bubbles ≤ 6mm)? | <input checked="" type="checkbox"/> Yes | |
| Were all soil VOAs field extracted with MeOH+BFB? | <input type="checkbox"/> N/A | |
| Note to Client: Any "No", answer above indicates non-compliance with standard procedures and may impact data quality. | | |
| Additional notes (if applicable): | | |



Sample Containers and Preservatives

| <u>Container Id</u> | <u>Preservative</u> | <u>Container Condition</u> | <u>Container Id</u> | <u>Preservative</u> | <u>Container Condition</u> |
|---------------------|--------------------------|----------------------------|---------------------|---------------------|----------------------------|
| 1215973001-A | HCL to pH < 2 | OK | | | |
| 1215973001-B | HCL to pH < 2 | OK | | | |
| 1215973001-C | HCL to pH < 2 | OK | | | |
| 1215973001-D | HCL to pH < 2 | OK | | | |
| 1215973001-E | HCL to pH < 2 | OK | | | |
| 1215973001-F | No Preservative Required | OK | | | |
| 1215973001-G | No Preservative Required | OK | | | |
| 1215973002-A | HCL to pH < 2 | OK | | | |
| 1215973002-B | HCL to pH < 2 | OK | | | |
| 1215973002-C | HCL to pH < 2 | OK | | | |
| 1215973002-D | HCL to pH < 2 | OK | | | |
| 1215973002-E | HCL to pH < 2 | OK | | | |
| 1215973002-F | No Preservative Required | OK | | | |
| 1215973002-G | No Preservative Required | OK | | | |
| 1215973003-A | HCL to pH < 2 | OK | | | |
| 1215973003-B | HCL to pH < 2 | OK | | | |
| 1215973003-C | HCL to pH < 2 | OK | | | |

Container Condition Glossary

Containers for bacteriological, low level mercury and VOA vials are not opened prior to analysis and will be assigned condition code OK unless evidence indicates than an inappropriate container was submitted.

OK - The container was received at an acceptable pH for the analysis requested.

BU - The container was received with headspace greater than 6mm.

DM - The container was received damaged.

FR - The container was received frozen and not usable for Bacteria or BOD analyses.

IC - The container provided for microbiology analysis was not a laboratory-supplied, pre-sterilized container and therefore was not suitable for analysis.

NC- The container provided was not preserved or was under-preserved. The method does not allow for additional preservative added after collection.

PA - The container was received outside of the acceptable pH for the analysis requested. Preservative was added upon receipt and the container is now at the correct pH. See the Sample Receipt Form for details on the amount and lot # of the preservative added.

PH - The container was received outside of the acceptable pH for the analysis requested. Preservative was added upon receipt, but was insufficient to bring the container to the correct pH for the analysis requested. See the Sample Receipt Form for details on the amount and lot # of the preservative added.

QN - Insufficient sample quantity provided.

Laboratory Data Review Checklist

Completed By:

Glenn Hasburgh

Title:

Environmental Scientist / QEP

Date:

12/13/2021

CS Report Name:

Report Date:

9/30/2021

Consultant Firm:

Environmental Management, Inc.

Laboratory Name:

SGS North America – Anchorage, Alaska

Laboratory Report Number:

1215973

ADEC File Number:

2105.26.001

Hazard Identification Number:

23166

1. Laboratory

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

 Yes No

Comments:

- 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

Comments:

2. Chain of Custody (CoC)

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

 Yes No

Comments:

- b. Correct Analyses requested?

 Yes No

Comments:

3. Laboratory Sample Receipt Documentation

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

 Yes No

Comments:

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

 Yes No

Comments:

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

 Yes No

Comments:

- 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

Comments:

There were no discrepancies.

- e. Data quality or usability affected?

Comments:

No, there is nothing to indicate data quality or usability has been affected.

4. Case Narrative

- a. Present and understandable?

Yes No

Comments:

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

Yes No

Comments:

- c. Were all corrective actions documented?

Yes No

Comments:

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

Comments:

There were multiple failed recoveries of surrogates in the lab QC samples. Many were not detected above the LOQ in the primary samples. In others, the narrative does not state specific effect.

5. Samples Results

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

Yes No

Comments:

- b. All applicable holding times met?

Yes No

Comments:

c. All soils reported on a dry weight basis?

Yes No

Comments:

NA, soils were not sampled.

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

Yes No

Comments:

The LOQ for 1,2,3-Trichloropropane is above the cleanup level. However, this is not a primary contaminant of concern.

e. Data quality or usability affected?

Yes No

Comments:

There slightly elevated LOQ does not impact the usability of the data.

6. QC Samples

a. Method Blank

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

Yes No

Comments:

ii. All method blank results less than limit of quantitation (LOQ)?

Yes No

Comments:

iii. If above LOQ, what samples are affected?

Comments:

NA, all were below the LOQ.

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

Yes No

Comments:

NA, there were no affected samples.

v. Data quality or usability affected?

Comments:

No, there is nothing to indicate data quality or usability has been 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

Comments:

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

Yes No

Comments:

NA, metals nor inorganics were analyzed.

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

Yes No

Comments:

No, there were elevated recoveries for multiple analytes, however, these analytes were not detected above the LOQ in the field samples.

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

Yes No

Comments:

RPD for naphthalene was high in LCS/LCSD. This analyte was not detected in the field samples.

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

Comments:

None of the samples are affected because the analytes with failures were not detected in the field samples.

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

Yes No

Comments:

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

Comments:

This has no effect on usability since the analytes were not detected in the field samples.

c. Surrogates – Organics Only

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

 Yes No

Comments:

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

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

 Yes No

Comments:

| |
|--|
| There were failed recoveries in multiple QC samples. All recoveries for field samples were acceptable. |
|--|

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

 Yes No

Comments:

| |
|-----|
| Yes |
|-----|

iv. Data quality or usability affected?

Comments:

| |
|--|
| Data quality nor usability are affected because the analytes with the failed recoveries were not detected above the LOQ. |
|--|

d. Trip blank – Volatile analyses only (GRO, BTEX, Volatile Chlorinated Solvents, etc.): Water and Soil

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

 Yes No

Comments:

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

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

Comments:

| |
|---|
| Only one cooler was used which was hand delivered to the lab. |
|---|

iii. All results less than LOQ?

 Yes No

Comments:

| |
|-----|
| Yes |
|-----|

iv. If above LOQ, what samples are affected?

Comments:

None, all results were below the LOQ.

v. Data quality or usability affected?

Comments:

No, there is nothing to indicate data quality or usability has been affected.

e. Field Duplicate

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

Yes No

Comments:

ii. Submitted blind to lab?

Yes No

Comments:

iii. Precision – All relative percent differences (RPD) less than specified DQOs?

(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

Comments:

All results were below LOQ, therefore, RPD could not be calculated.

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

Comments:

No, there is nothing to indicate data quality or usability has been affected.

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

Yes No Not Applicable

i. All results less than LOQ?

Yes No

Comments:

NA, an equipment blank was not collected.

ii. If above LOQ, what samples are affected?

Comments:

NA, an equipment blank was not collected.

iii. Data quality or usability affected?

Comments:

Data quality or usability is not affected because all analytes in the field samples were below the LOQ.

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

a. Defined and appropriate?

Yes No

Comments:

APPENDIX D

Disposal Documentation



CERTIFICATE OF DISPOSAL/RECYCLE

GENERATOR: SHORESIDE PETROLEUM
6401 LAKE OTIS PKWY
ANCHORAGE, AK 99507

DISPOSAL FACILITY: US ECOLOGY ALASKA LLC
2020 VIKING DRIVE
ANCHORAGE, AK 99501

EPA ID NUMBER: EXEMPT
MANIFEST/DOCUMENT #: BOL80565
DATE OF DISPOSAL/RECYCLE: DEC-20-2021

| <u>LINE</u> | <u>WASTE DESCRIPTION</u> | <u>CONTAINERS</u> | <u>TYPE</u> | <u>QUANTITY</u> | <u>UOM</u> |
|-------------|--------------------------|-------------------|-------------|-----------------|------------|
| 1 | OILY WATER (DM05) | 3 | DM | 15 | G |

I certify, on behalf of the above listed treatment facility, that to the best of my knowledge, the above described waste was managed in compliance with all applicable laws, regulations, permits and licenses on the date listed above.

PREPARED BY: Michael Cooney

SIGNATURE: [Signature]

DATE: DEC 20 2021



**ALASKA DEPARTMENT OF ENVIRONMENTAL CONSERVATION
DIVISION OF SPILL PREVENTION AND RESPONSE
Contaminated Sites and Prevention Preparedness and Response Programs**

Contaminated Media Transport and Treatment or Disposal Approval Form

| | | | |
|---|--|--|--|
| DEC HAZARD/SPILL ID # | NAME OF CONTAMINATED SITE OR SPILL | | |
| 23166 | Essential 1 (former Shoreside Texaco) Bird Creek | | |
| CONTAMINATED SITE OR SPILL LOCATION – ADDRESS OR OTHER APPROPRIATE DESCRIPTION | | | |
| Mile 101 / 100.7 Seward Highway, Girdwood, Alaska 99587 | | | |
| CURRENT PHYSICAL LOCATION OF MEDIA | | SOURCE OF THE CONTAMINATION (DAY TANK, WASH BAY, FIRE TRAINING PIT, LUST, ETC.) | |
| Same, in 3 closed 5 gallon buckets | | leaking underground storage tank | |
| CONTAMINANTS OF CONCERN | ESTIMATED VOLUME | DATE(S) GENERATED | |
| Petroleum / Diesel | 15 gallons | September 10, 2021 | |
| POST TREATMENT ANALYSIS REQUIRED (such as GRO, DRO, RRO, VOCs, metals, PFAS, and/or Chlorinated Solvents) | | | |
| None | | | |
| COMMENTS OR OTHER IMPORTANT INFORMATION | | | |
| Decontamination and purge water from groundwater monitoring event; 1 well sampled. No DRO, VOC, or PAH analytes were detected in the sampled well. This water will be transported to Shoreside Petroleum's facility at 6401 Lake Otis Pkwy, Anchorage, AK, where it will be placed into their used oil tank which is regularly pumped out and treated by NRC. | | | |

| | |
|--|--|
| TREATMENT FACILITY, LANDFILL, AND/OR FINAL DESTINATION OF MEDIA | PHYSICAL ADDRESS/PHONE NUMBER |
| US Ecology | 2020 Viking Drive, Anchorage, Alaska 99501 - (907) 258-1558 |
| RESPONSIBLE PARTY | ADDRESS/PHONE NUMBER |
| Shoreside Petroleum | 6401 Lake Otis Parkway, Anchorage, Alaska 99507 - (907) 344-4571 |
| WASTE MANAGEMENT CO. / ORGANIZER | ADDRESS/PHONE NUMBER |
| US Ecology | 619 East Ship Creek Avenue, Suite 309, Alaska 99501 - (907) 258-1558 |

*Note, disposal of polluted soil in a landfill requires prior approval from the landfill operator and ADEC Solid Waste Program.

Andy Coulson

Name of the Person Requesting Approval (printed)

Andy Coulson

Signature

Jessica Hall

Environmental Scientist / EMI

Title/Association

11/10/2021 **(907) 272-9336**

Date

Phone Number

-----DEC USE ONLY-----

Based on the information provided, ADEC approves transport of the above mentioned material. The Responsible Party or their consultant must submit to the DEC Project Manager a copy of weight receipts of the loads transported and a post treatment analytical report, if disposed of at an approved treatment facility. The contaminated soil shall be transported as a covered load in compliance with 18 AAC 60.015.

DEC Project Manager Name (printed)

Project Manager Title

Signature

Date

Phone Number