
Brandon Landfill 31 Corona Street Brandon, Vermont

VT DEC Project# RU97-0128
Solid Waste Facility ID# RU080
KAS Job# 609210052

SPRING 2023 SEMI-ANNUAL WATER QUALITY MONITORING REPORT

July 25, 2023

Prepared for:

Town of Brandon
49 Center Street
Brandon, VT 05733



589 Avenue D, Suite 10
PO Box 787
Williston, VT 05495

www.kas-consulting.com

802.383.0486 p
802.383.0490 f



Introduction

KAS, Inc. (KAS) conducted a semi-annual water quality monitoring event on May 22, 2023 at the Brandon Landfill (Site Location Map and Site Map in Appendix A). The spring 2023 groundwater monitoring was conducted in accordance with the current landfill post-closure certification. In addition, as requested by the Vermont Department of Environmental Conservation (VT DEC), the monitoring wells were tested for per- and polyfluoroalkyl substances (PFAS), an emerging group of contaminants, that have been frequently found in landfill leachate.

Background

The 5-acre facility operated as a landfill from 1940 until its closure in 1992, and currently operates as a transfer station. Post-closure groundwater monitoring has been conducted consistently since 2016, although select wells have not been sampled for various reasons (i.e., dry, inaccessible, etc.). Manganese, arsenic, and lead remain at levels above Vermont Groundwater Enforceable Standard (VGES). Other metals that have infrequently exceeded VGES in the past include cadmium and nickel. Volatile organic compounds (VOCs) have generally remained below VGES, with the exception of naphthalene, which was found to be slightly above VGES in June 2022. May 2023 was the first time groundwater was tested for PFAS.

PFAS compounds subject to regulation in Vermont include perfluorooctanoic acid (PFOA), perfluorooctanesulfonic acid (PFOS), perfluorohexanesulfonic acid (PFHxS), perfluoroheptanoic acid (PFHpA), and perfluorononanoic acid (PFNA). The VGES for PFAS is 20 nanograms per liter (ng/L) for the sum of the five regulated PFAS. There are numerous other PFAS compounds that are not regulated in Vermont, some of which are considered replacements for PFAS that have been historically phased out of production and use.

Groundwater is presumed to flow in a south-southwesterly direction, towards Otter Creek, based on the topography of the area and based on the data collected to date. The current monitoring well network consists of four (4) wells: MW-1 (upgradient), MW-3 (cross-gradient), and MW-2C and MW-5 (both of which are downgradient).

Groundwater Sampling & Results

Field measurements

At the time of sample collection, groundwater was field analyzed for temperature, pH, and specific conductance using a properly calibrated YSI® meter. The depth to water was gauged using a Geotech™ water level indicator. Field measurement data is tabulated in Appendix B.

Laboratory Results

A groundwater sample was collected from all monitoring wells (MW-1, MW-2C, MW-3, and MW-5). Low-flow groundwater purging and sampling techniques were used at MW-2C and MW-5; however, at MW-2C purging until stabilization was not conducted due to a low water column. At MW-3 and MW-5, the sample was collected via bailer due to the depth of water exceeding the capacity of the peristaltic pump (e.g., >30 feet below top of casing). The groundwater samples were analyzed for:

- Total metals (e.g., arsenic, cadmium, chromium, copper, lead, iron, manganese, mercury, nickel, sodium, and zinc) via EPA Method 6010/6020;
- Chemical oxygen demand (COD) via Method 410.4;

- Chloride via Method 300.0;
- VOCs via EPA Method 8260C; and,
- PFAS via isotope dilution.

At MW-2C, MW-3 and MW-5, manganese concentrations were above VGES (0.3 mg/L), with the highest level being at MW-5 (2.1 mg/L). At MW-3, the lead concentration of 0.0151 mg/L was slightly above VGES (0.015 mg/L). No VOCs were detected above laboratory method detection limits in the samples collected, except for diethyl ether at MW-2C, for which there is no groundwater quality standard. Total regulated PFAS exceeded VGES (20 ng/L) at MW-2C and MW-5, with the highest level being at MW-2C (305.7 ng/L). Current and historical analytical data are provided in tables and graphs in Appendix B. A copy of the laboratory reports is provided in Appendix C.

Quality Assurance/Quality Control

Quality assurance and quality control (QA/QC) samples included a duplicate sample that was analyzed for VOCs, metals, chloride, and COD. The results of the laboratory analysis of the duplicate sample were analyzed using a relative percent difference (RPD) analysis. The RPD is defined as 100 times the difference in reported concentration between sample and duplicate, divided by the mean of the two samples. A small RPD indicates good correlation between sample and duplicate. The absolute RPD values ranged between 1.9 and 15.8%, which good correlation/precision based on an EPA Region 1 guideline value of 30%.

A QA/QC sample also included a trip blank for VOC analysis. No VOC were detected in the trip blank, which indicates that potential contamination from transit, sample bottles, or laboratory conditions was not a concern.

For PFAS analysis, a QA/QC sample included an equipment rinsate blank (ERB) sample. No PFAS was detected in the ERB sample, which indicates that there was no cross contamination of PFAS from the sampler, equipment, field conditions, and/or laboratory conditions.

Trends

Additional field measurement data is needed to establish long-term trends.

Overall, metal concentrations appear to have decreased from their respective historical peaks. The exception is manganese, the levels of which continue to widely fluctuate. In general, sodium, chloride, and COD also continue to fluctuate. Trends/graphs are provided in Appendix B.

Recommendation

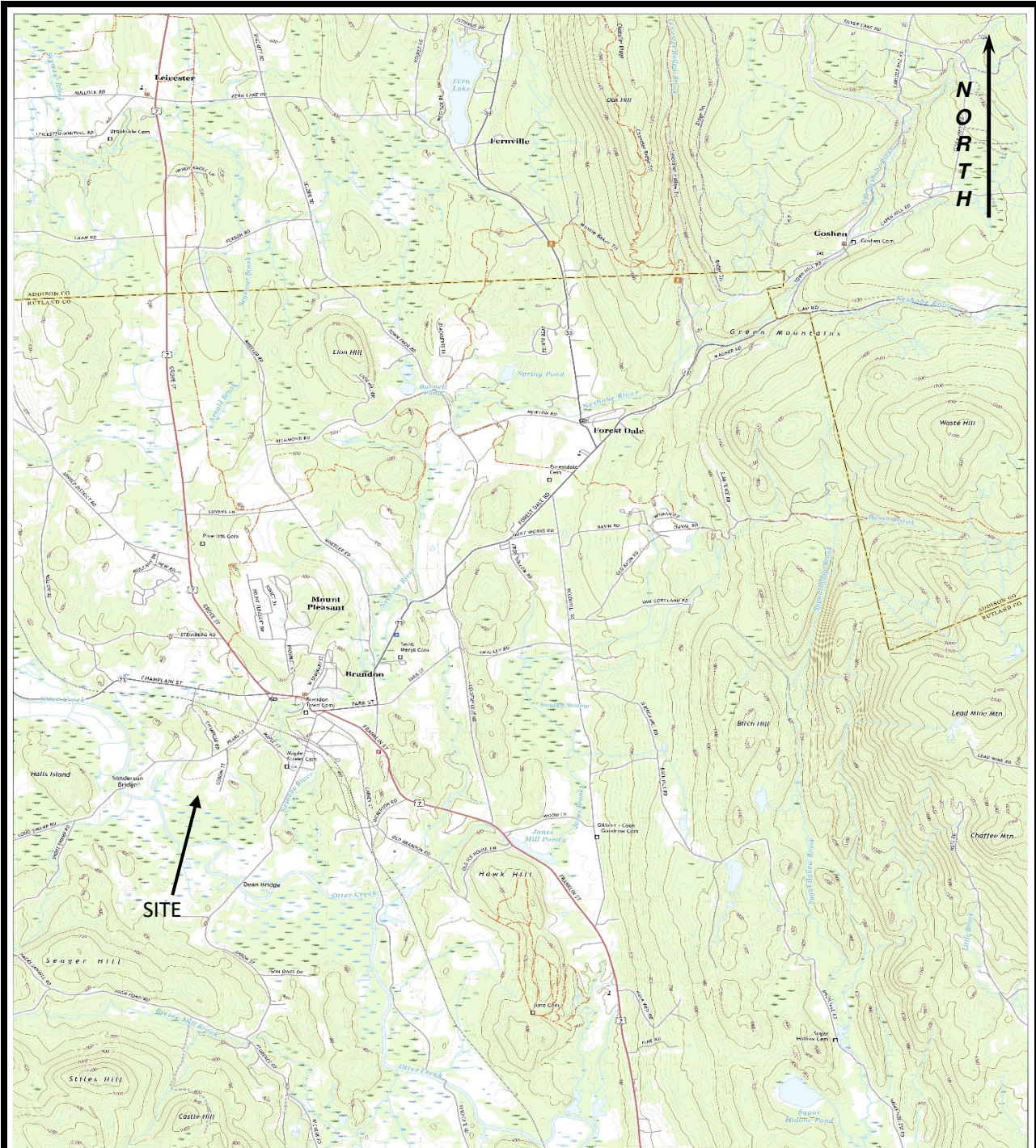
KAS recommends that groundwater monitoring continue in accordance with Brandon Solid Waste Facility Certification – Monitoring Requirements (7), with the next monitoring event to occur in October 2023. PFAS sampling and analysis should also continue at all monitoring wells.

At this time, it is unknown if PFAS has impacted the deeper aquifer. In general, the surrounding residential area is served by municipal water. As shown on the Site Map (Appendix A), there are no private water supply wells downgradient from the landfill in the nearby vicinity. The nearest cross-gradient supply well is approximately 0.16 miles to the west. Therefore, based on the distances and locations, the risk of PFAS contamination in the supply wells is considered low at this time; however, sampling and analysis would be required to fully rule out the risk.



APPENDIX A

Site Location Map and Site Map



KAS Job Number 609210052
 Source: USGS



589 Avenue D, Suite 10
 PO Box 787
 Williston, VT 05495
 www.kas-consulting.com
 802 383.0486 p
 802 383.0490 f

Brandon Town Landfill 31 Corona Street, Brandon, Vermont

Site Location Map
 USGS Mapping

| | | | |
|----------------|---------------|-----------|--------|
| Date: 04/27/22 | Drawing No. 0 | Scale NTS | By: ML |
|----------------|---------------|-----------|--------|



LEGEND

- Private Wells**
- GPS Located
 - Screen Digitized
 - E911 Address Matched
 - Welldriller/Clarion
 - Unknown Location Method
 - Incorrectly Located
- Public Water Sources**

Markups by KAS

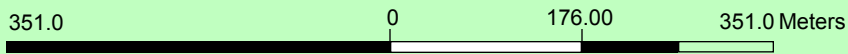
- - Approx. location of monitoring well
- - Inactive monitoring well

NOTES

Map created using ANR GIS mapping technology.

1: 6,916

September 13, 2022





APPENDIX B

Data Summaries

Brandon Closed Landfill
 Sampling Date: May 22, 2023

| PARAMETER | Monitoring Well ID: | MW-1 | MW-2C | MW-3 | MW-5 | VGES | PAL |
|--|---------------------|---------------|---------------|---------------|-------------|-------|--------|
| VOCs (ug/L) | | | | | | | |
| Diethyl Ether | | ND | 14.6 | ND | ND | - | - |
| Total Metals (mg/L) | | | | | | | |
| Arsenic | | 0.0014 | <0.0010 | 0.0075 | <0.0010 | 0.010 | 0.001 |
| Cadmium | | <0.0020 | <0.0020 | <0.0020 | <0.0020 | 0.005 | 0.001 |
| Chromium | | <0.0050 | <0.0050 | 0.0110 | <0.0050 | 0.100 | 0.050 |
| Copper | | <0.020 | <0.020 | 0.021 | <0.020 | 1.300 | 0.650 |
| Iron | | 2.4 | 2.4 | 18 | 0.52 | - | - |
| Lead | | <0.0010 | 0.0036 | 0.0151 | <0.0010 | 0.015 | 0.002 |
| Manganese | | 0.25 | 1.3 | 0.55 | 2.1 | 0.300 | 0.150 |
| Mercury | | <0.0002 | <0.0002 | <0.0002 | <0.0002 | 0.002 | 0.0005 |
| Nickel | | <0.0050 | 0.0081 | 0.0140 | <0.0050 | 0.100 | 0.050 |
| Sodium | | 41 | 28 | 26 | 34 | - | - |
| Zinc | | <0.020 | <0.020 | 0.042 | <0.020 | - | - |
| Other Analytes (mg/L) | | | | | | | |
| Chloride | | 77 | 24 | 44 | 59 | - | - |
| COD | | 74 | 77 | 79 | 35 | - | - |
| PFAS (ng/L) | | | | | | | |
| Perfluorohexanesulfonic acid (PFHxS) | | <1.8 | 36 | <1.8 | 17 | - | - |
| Perfluoroheptanoic acid (PFHpA) | | <1.8 | 18 | <1.8 | 5.1 | | |
| Perfluorooctanoic acid (PFOA) | | <1.8 | 97 | <1.8 | 30 | | |
| Perfluorooctanesulfonic acid (PFOS) | | 2.4 | 150 | 5.2 | 14 | | |
| Perfluorononanoic acid (PFNA) | | <1.8 | 4.7 | <1.8 | <1.9 | | |
| Total Regulated PFAS | | 2.4 | 305.7 | 5.2 | 66.1 | 20 | 2 |
| Total Non-Regulated PFAS | | ND | 68.7 | ND | 14.8 | - | - |
| Field Measurements (units as noted) | | | | | | | |
| pH (std units) | | 7.26 | 6.96 | 7.57 | 6.69 | - | - |
| Temperature (deg C) | | 14.5 | 11.3 | 13.0 | 10.1 | - | - |
| Conductivity (uS) | | 1,008 | 1,487 | 715 | 826 | - | - |
| Water Level (feet btoc) | | 30.67 | 8.36 | 31.50 | 4.79 | - | - |

Only detected or targeted VOCs are depicted

All values reported in units noted above

"-" = Not Analyzed, No Information or No Applicable Standard

ND = None Detected

<X = None Detected above Detection Limit (X)

VGES = Vermont Groundwater Enforcement Standard (GWPRS 07/06/2019)

PAL = Vermont Preventive Action Level (GWPRS 07/06/2019)

Detections are **bolded**

>VGES

Bold (italic) indicates value exceeds PAL

Brandon Closed Landfill

MW-1

| PARAMETER | Sample Date: | 2012 | 4/22/2013 | 10/1/2013 | 5/29/2014 | 6/9/2015 | 10/25/2015 | 7/24/2016 | October 2016 | 5/7/2017 | 10/28/2017 | 5/28/2018 | 10/20/2018 | 05/27/2019 | 10/26/2019 | VGES | PAL |
|-------------------------------------|--------------|-------------|-------------|-------------|-------------|-------------|-------------|-------------|--------------|--------------|--------------|---------------|--------------|---------------|--------------|-------|--------|
| VOCs (ug/L) | | | | | | | | | | | | | | | | | |
| 1,1-dichloroethane | | - | - | - | - | - | - | - | - | <1.0 | <1.0 | <1.0 | <1.0 | <1 | <1 | 70 | 35 |
| Total Metals (mg/L) | | | | | | | | | | | | | | | | | |
| Arsenic | Well | Well | Well | Well | Well | Well | Well | Well | Well | 0.001 | <0.0010 | 0.0016 | <0.0010 | <0.0010 | <0.0010 | 0.010 | 0.001 |
| Cadmium | Not Sampled | Not Sampled | Not Sampled | Not Sampled | Not Sampled | Not Sampled | Not Sampled | Not Sampled | Not Sampled | <0.0020 | <0.0020 | <0.0020 | <0.0020 | <0.0020 | <0.0020 | 0.005 | 0.001 |
| Chromium | | | | | | | | | | <0.0050 | <0.0050 | <0.0050 | <0.0050 | <0.0050 | <0.0050 | 0.100 | 0.050 |
| Copper | | | | | | | | | | <0.020 | <0.020 | <0.020 | <0.020 | <0.020 | <0.020 | 1.300 | 0.650 |
| Iron | | | | | | | | | | 0.79 | 0.51 | 3.0 | 0.62 | 0.091 | 0.26 | - | - |
| Lead | | | | | | | | | | <0.001 | <0.0010 | <0.0021 | <0.0010 | <0.0010 | <0.0010 | 0.015 | 0.002 |
| Manganese | | | | | | | | | | 0.18 | 0.14 | 0.79 | 0.21 | 0.034 | 0.028 | 0.300 | 0.150 |
| Mercury | | | | | | | | | | <0.0002 | <0.0002 | <0.0002 | <0.0002 | <0.0002 | <0.0002 | 0.002 | 0.0005 |
| Nickel | | | | | | | | | | <0.0050 | <0.0050 | <0.0050 | <0.0050 | <0.0050 | <0.0050 | 0.100 | 0.050 |
| Sodium | | | | | | | | | | 40 | 40 | 38 | 42 | 38 | 38 | - | - |
| Zinc | | | | | | | | | | 0.046 | 0.170 | 0.025 | 0.025 | <0.020 | <0.020 | - | - |
| Other Analytes (mg/L) | | | | | | | | | | | | | | | | | |
| Chloride | | | | | | | | | | 74 | 72 | 77 | 82 | 86 | 83 | - | - |
| COD | | | | | | | | | | 21 | 24 | 14 | 12 | <10 | 39 | - | - |
| Field Measurements (units as noted) | | | | | | | | | | | | | | | | | |
| pH (std units) | | | | | | | | | | 6.5 | 6.3 | 6.6 | 6.6 | 6.5 | 6.7 | - | - |
| Temperature (deg C) | | | | | | | | | | 10.7 | 11.6 | 11.8 | 11.9 | 11.2 | 12.1 | - | - |
| Conductivity (uS) | | | | | | | | | | 1050 | 1030 | 1020 | 1000 | 1100 | 1050 | - | - |
| Water Level (feet btoc) | | | | | | | | | | - | - | - | - | - | - | - | - |

| PARAMETER | Sample Date: | 10/14/2021 | 6/3/2022 | 10/27/2022 | 5/22/2023 | | | | | | | | | | | VGES | PAL |
|--------------------------------------|------------------|------------------|---------------|---------------|-----------|--|--|--|--|--|--|--|--|--|--|-------|--------|
| VOCs (ug/L) | | | | | | | | | | | | | | | | | |
| 1,1-dichloroethane | | - | - | <1.0 | <1.0 | | | | | | | | | | | 70 | 35 |
| Total Metals (mg/L) | | | | | | | | | | | | | | | | | |
| Arsenic | No Sample | No Sample | 0.0015 | 0.0014 | | | | | | | | | | | | 0.010 | 0.001 |
| Cadmium | | | <0.0020 | <0.0020 | | | | | | | | | | | | 0.005 | 0.001 |
| Chromium | | | <0.0050 | <0.0050 | | | | | | | | | | | | 0.100 | 0.050 |
| Copper | Unable To Locate | Unable To Locate | <0.020 | <0.020 | | | | | | | | | | | | 1.300 | 0.650 |
| Iron | | | 2.4 | 2.4 | | | | | | | | | | | | - | - |
| Lead | Well | Well | <0.0010 | <0.0010 | | | | | | | | | | | | 0.015 | 0.002 |
| Manganese | | | 0.20 | 0.25 | | | | | | | | | | | | 0.300 | 0.150 |
| Mercury | | | <0.0002 | <0.0002 | | | | | | | | | | | | 0.002 | 0.0005 |
| Nickel | | | <0.0050 | <0.0050 | | | | | | | | | | | | 0.100 | 0.050 |
| Sodium | | | 39 | 41 | | | | | | | | | | | | - | - |
| Zinc | | | 0.024 | <0.020 | | | | | | | | | | | | - | - |
| Other Analytes (mg/L) | | | | | | | | | | | | | | | | | |
| Chloride | | | 78 | 77 | | | | | | | | | | | | - | - |
| COD | | | 59 | 74 | | | | | | | | | | | | - | - |
| PFAS (ng/L) | | | | | | | | | | | | | | | | | |
| Perfluorohexanesulfonic acid (PFHxS) | | | - | <1.8 | | | | | | | | | | | | | |
| Perfluoroheptanoic acid (PFHpA) | | | - | <1.8 | | | | | | | | | | | | | |
| Perfluorooctanoic acid (PFOA) | | | - | <1.8 | | | | | | | | | | | | | |
| Perfluorooctanesulfonic acid (PFOS) | | | - | 2.4 | | | | | | | | | | | | | |
| Perfluorononanoic acid (PFNA) | | | - | <1.8 | | | | | | | | | | | | | |
| Total Regulated PFAS | | | - | 2.4 | | | | | | | | | | | | 20 | 2 |
| Total Non-Regulated PFAS | | | - | ND | | | | | | | | | | | | - | - |
| Field Measurements (units as noted) | | | | | | | | | | | | | | | | | |
| pH (std units) | | | | 7.21 | 7.26 | | | | | | | | | | | - | - |
| Temperature (deg C) | | | | 10.8 | 14.5 | | | | | | | | | | | - | - |
| Spec. Conductivity (uS/cm) | | | | 1,039 | 1,008 | | | | | | | | | | | - | - |
| Water Level (feet btoc) | | | | 31.17 | 30.67 | | | | | | | | | | | - | - |

Data prior to 2021 collected by others and obtained from 2019 reports prepared by EIV Technical Services

Data 2021-present collected by KAS, Inc.

Only detected or targeted VOCs are depicted

All values reported in units noted above

"-" = Not Analyzed, No Information or No Applicable Standard

ND = None Detected

<X = None Detected above Detection Limit (X)

VGES = Vermont Groundwater Enforcement Standard (GWPRS 07/06/2019)

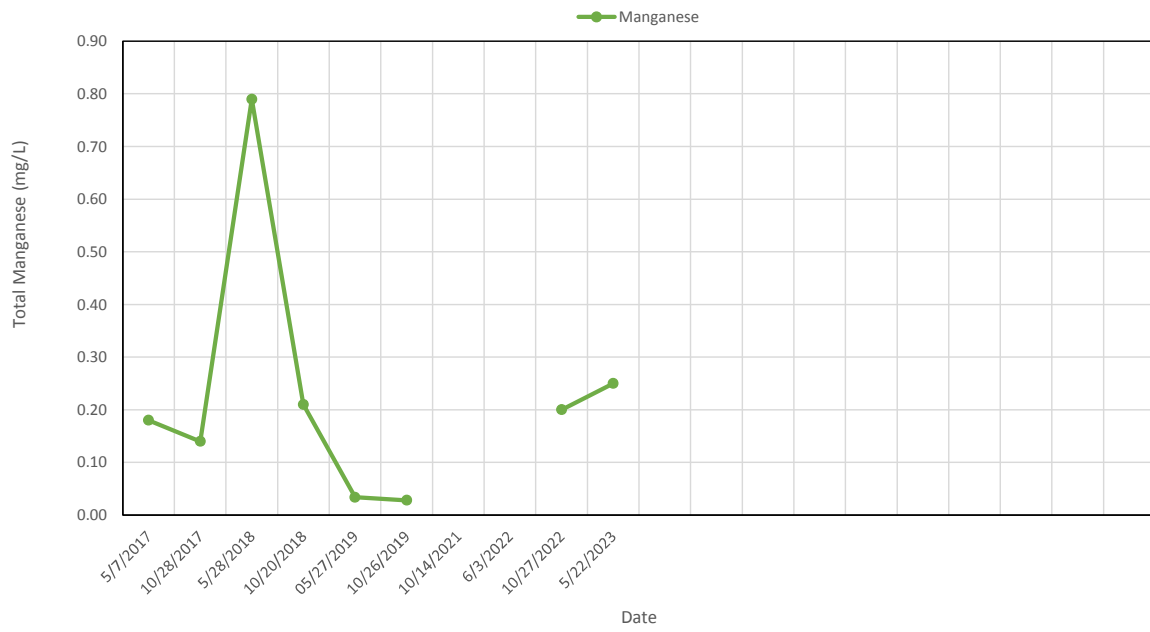
PAL = Vermont Preventive Action Level (GWPRS 07/06/2019)

Detections are **bolded**

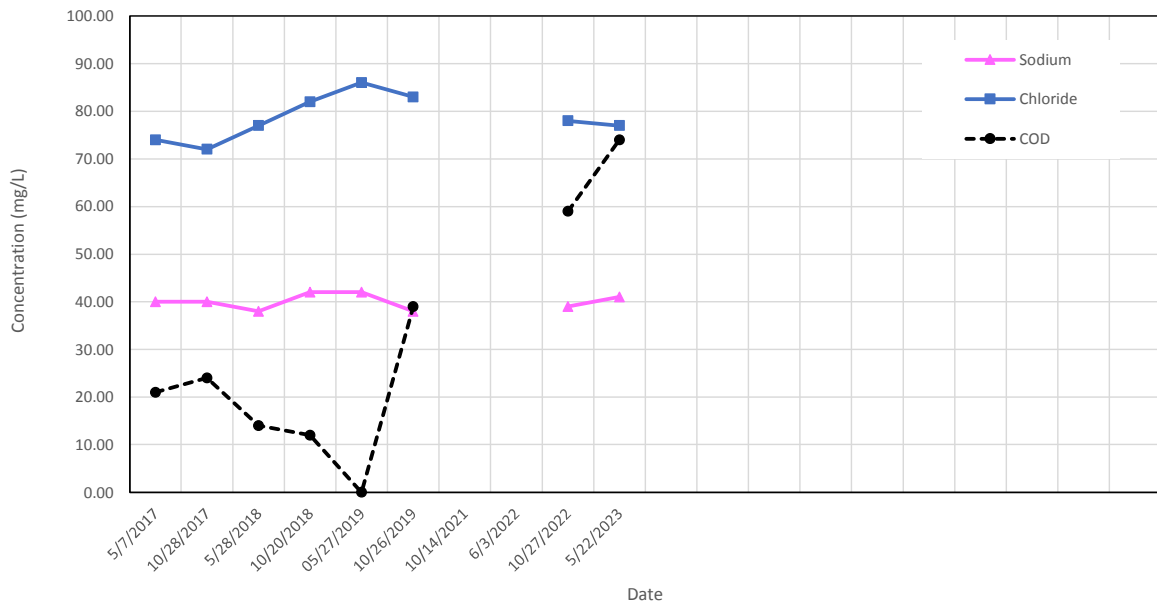
>VGES

bold (italic) indicates value exceeds PAL

Brandon Landfill
MW-1



Brandon Landfill
MW-1



Brandon Closed Landfill

MW-2C

| PARAMETER | Sample Date: | 2012 | 4/22/2013 | 10/1/2013 | 5/29/2014 | 6/9/2015 | 10/25/2015 | 7/24/2016 | October 2016 | 5/7/2017 | 10/28/2017 | 5/28/2018 | 10/20/2018 | 05/27/2019 | 10/26/2019 | VGES | PAL |
|-------------------------------------|--------------|-------------|-------------|--------------|--------------|--------------|--------------|------------|---------------|---------------|---------------|---------------|---------------|--------------|------------|--------|-------|
| VOCs (ug/L) | | | | | | | | | | | | | | | | | |
| Dichlorodifluoromethane | - | <5.0 | <5.0 | <5.0 | <5.0 | <5.0 | <5.0 | 7.2 | - | <5.0 | <5.0 | <5.0 | 5.0 | 1.8 | <1 | - | - |
| Vinyl Chloride | - | <2.0 | <2.0 | <2.0 | <2.0 | <2.0 | <2.0 | 0.7 | - | <0.5 | <0.5 | <0.5 | <0.5 | - | - | 2 | 0.5 |
| 1,4-dichlorobenzene | 2.3 | 2.4 | 2.5 | 2.3 | 2.1 | 2.0 | 2.6 | <1.0 | <1.0 | 2.4 | 2.2 | 1.7 | 2.7 | - | - | 75 | 38 |
| Acetone | <10.0 | <10.0 | <10.0 | <10.0 | <10.0 | <10.0 | <10.0 | <10.0 | <10.0 | <10.0 | <10.0 | <10.0 | <10.0 | 15.8 | <10.0 | 950 | 475 |
| Benzene | 2.0 | 2.3 | 3.3 | 2.6 | 2.3 | 1.9 | 3.4 | <0.5 | 2.9 | 2.9 | 1.3 | 3.5 | 1.2 | <0.5 | 5 | 0.5 | |
| Chlorobenzene | 5.4 | 5.3 | 4.8 | 6.0 | 5.9 | 5.0 | 7.6 | <1.0 | 8.3 | 7.2 | 4.1 | 7.4 | 4.9 | <1.0 | 100 | 50 | |
| Diethyl Ether | - | 22.5 | 36.2 | 24.0 | 23.9 | 19.5 | 26.3 | - | 19.4 | 25.2 | 17.5 | 30.5 | - | - | - | - | |
| Total Metals (mg/L) | | | | | | | | | | | | | | | | | |
| Arsenic | - | - | - | 0.074 | 0.026 | 0.025 | 0.027 | Data | 0.034 | 0.0409 | 0.071 | 0.0200 | 0.0213 | <0.0010 | 0.010 | 0.001 | |
| Cadmium | - | - | - | <0.002 | <0.002 | <0.002 | 0.021 | Not | 0.0027 | <0.0020 | <0.0020 | <0.0020 | <0.0020 | <0.0020 | <0.0020 | 0.005 | 0.001 |
| Chromium | - | - | - | 0.012 | <0.005 | <0.005 | <0.0050 | Available | <0.0050 | <0.0050 | <0.0050 | <0.0050 | <0.0050 | <0.0050 | <0.0050 | 0.100 | 0.050 |
| Copper | - | - | - | <0.020 | <0.020 | <0.020 | <0.020 | - | <0.020 | <0.020 | <0.020 | <0.020 | <0.020 | <0.020 | <0.020 | 1.300 | 0.650 |
| Iron | - | - | - | 56 | 28 | 22 | 33 | - | 32 | 33 | 65 | 33 | 28 | 0.28 | - | - | |
| Lead | - | - | - | 0.007 | <0.001 | <0.001 | <0.001 | - | <0.001 | <0.0010 | 0.0013 | <0.0010 | <0.0010 | <0.0010 | 0.015 | 0.002 | |
| Manganese | - | - | - | 0.92 | 0.54 | 0.45 | 0.67 | - | 0.54 | 0.58 | 0.57 | 0.58 | 0.53 | 0.077 | 0.300 | 0.150 | |
| Mercury | - | - | - | <0.0002 | <0.0002 | <0.0002 | <0.0002 | - | <0.0002 | <0.0002 | <0.0002 | <0.0002 | <0.0002 | <0.0002 | 0.002 | 0.0005 | |
| Nickel | - | - | - | 0.022 | 0.013 | 0.012 | <0.0050 | - | 0.0088 | 0.0074 | 0.0091 | 0.0095 | 0.0092 | <0.0050 | 0.100 | 0.050 | |
| Sodium | - | - | - | 170 | 160 | 130 | 230 | - | 270 | 230 | 270 | 290 | 240 | 18 | - | - | |
| Zinc | - | - | - | 0.043 | <0.02 | <0.02 | <0.020 | - | <0.020 | <0.020 | 0.025 | <0.020 | <0.020 | <0.020 | - | - | |
| Other Analytes (mg/L) | | | | | | | | | | | | | | | | | |
| Chloride | - | - | - | 203 | 280 | 290 | 380 | - | 480 | 440 | 450 | 500 | 420 | 32 | - | - | |
| COD | - | - | - | 57 | 100 | 63 | 62 | - | 52 | 67 | 41 | 53 | 62 | 32 | - | - | |
| Field Measurements (units as noted) | | | | | | | | | | | | | | | | | |
| pH (std units) | - | - | - | - | - | - | 6.5 | - | 6.5 | 6.4 | 6.4 | 6.3 | 6.6 | 6.5 | - | - | |
| Temperature (deg C) | - | - | - | - | - | - | 13 | - | 9.8 | 10.2 | 10.2 | 11.2 | 10.9 | 11.3 | - | - | |
| Spec. Conductivity (uS/cm) | - | - | - | - | - | - | 3,010 | - | 2,800 | 2,900 | 2,800 | 2,300 | 2,100 | 2,210 | - | - | |
| Water Level (feet btoc) | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - | |

| PARAMETER | Sample Date: | 10/14/2021 | 6/3/2022 | 10/27/2022 | 5/22/2023 | | | | | | | | | | | VGES | PAL |
|--------------------------------------|--------------|---------------|----------|---------------|-----------|---|---|---|---|---|---|---|---|---|---|-------|--------|
| VOCs (ug/L) | | | | | | | | | | | | | | | | | |
| 1,4-dichlorobenzene | - | <1 | - | <1.0 | - | - | - | - | - | - | - | - | - | - | - | 75 | 38 |
| Diethyl Ether | - | 21 | - | 14.6 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Acetone | - | 67 | - | <10.0 | - | - | - | - | - | - | - | - | - | - | - | 950 | 475 |
| Methyl-t-butyl ether (MTBE) | - | 1.4 | - | <2.0 | - | - | - | - | - | - | - | - | - | - | - | 11 | 5 |
| Tetrahydrofuran | - | 17 | - | <10.0 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Benzene | - | <1 | - | <0.5 | - | - | - | - | - | - | - | - | - | - | - | 5 | 0.5 |
| Chlorobenzene | - | <1 | - | <1.0 | - | - | - | - | - | - | - | - | - | - | - | 100 | 50 |
| Naphthalene | - | 0.56 | - | <0.5 | - | - | - | - | - | - | - | - | - | - | - | 0.5 | 0.5 |
| Total Metals (mg/L) | | | | | | | | | | | | | | | | | |
| Arsenic | No | 0.016 | No | <0.0010 | - | - | - | - | - | - | - | - | - | - | - | 0.010 | 0.001 |
| Cadmium | Sample | 0.0022 | Sample | <0.0020 | - | - | - | - | - | - | - | - | - | - | - | 0.005 | 0.001 |
| Chromium | - | 0.0093 | - | <0.0050 | - | - | - | - | - | - | - | - | - | - | - | 0.100 | 0.050 |
| Copper | Well | 0.084 | Well | <0.020 | - | - | - | - | - | - | - | - | - | - | - | 1.300 | 0.650 |
| Iron | Dry | 22 | Dry | 2.4 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Lead | - | 0.17 | - | 0.0036 | - | - | - | - | - | - | - | - | - | - | - | 0.015 | 0.002 |
| Manganese | - | 1.9 | - | 1.3 | - | - | - | - | - | - | - | - | - | - | - | 0.300 | 0.150 |
| Mercury | - | <0.0001 | - | <0.0002 | - | - | - | - | - | - | - | - | - | - | - | 0.002 | 0.0005 |
| Nickel | - | 0.033 | - | 0.0081 | - | - | - | - | - | - | - | - | - | - | - | 0.100 | 0.050 |
| Sodium | - | 38 | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Zinc | - | 0.094 | - | <0.020 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Other Analytes (mg/L) | | | | | | | | | | | | | | | | | |
| Chloride | - | 41 | - | 24 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| COD | - | 900 | - | 77 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| PFAS (ng/L) | | | | | | | | | | | | | | | | | |
| Perfluorohexanesulfonic acid (PFHxS) | - | - | - | 36 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Perfluoroheptanoic acid (PFHpA) | - | - | - | 18 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Perfluorooctanoic acid (PFOA) | - | - | - | 97 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Perfluorooctanesulfonic acid (PFOS) | - | - | - | 150 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Perfluorononanoic acid (PFNA) | - | - | - | 4.7 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Total Regulated PFAS | - | - | - | 305.7 | - | - | - | - | - | - | - | - | - | - | - | 20 | 2 |
| Total Non-Regulated PFAS | - | - | - | 68.7 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Field Measurements (units as noted) | | | | | | | | | | | | | | | | | |
| pH (std units) | - | 6.51 | - | 6.96 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Temperature (deg C) | - | 18.1 | - | 11.3 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Spec. Conductivity (uS/cm) | - | 1,643 | - | 1,487 | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Water Level (feet btoc) | - | 8.20 | - | 8.36 | - | - | - | - | - | - | - | - | - | - | - | - | - |

Data prior to 2021 collected by others and obtained from 2019 reports prepared by EIV Technical Services

Data 2021-present collected by KAS, Inc.

Only detected or targeted VOCs are depicted

All values reported in units noted above

"-" = Not Analyzed, No Information or No Applicable Standard

ND = None Detected

<X = None Detected above Detection Limit (X)

VGES = Vermont Groundwater Enforcement Standard (GWPRS 07/06/2019)

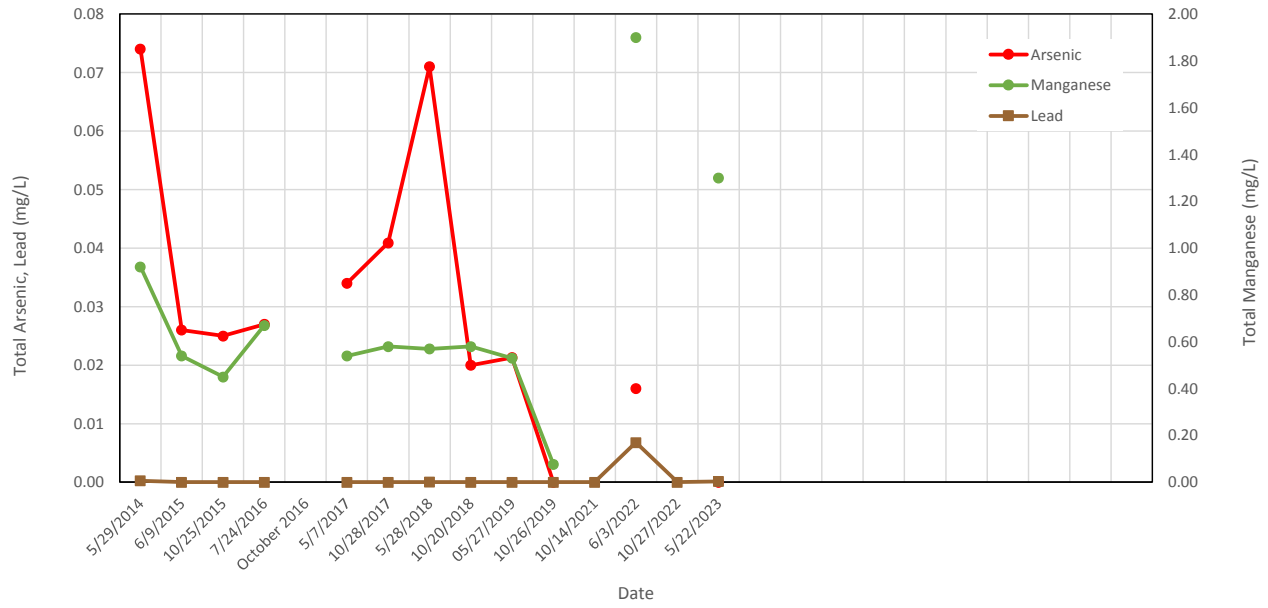
PAL = Vermont Preventive Action Level (GWPRS 07/06/2019)

Detections are bolded

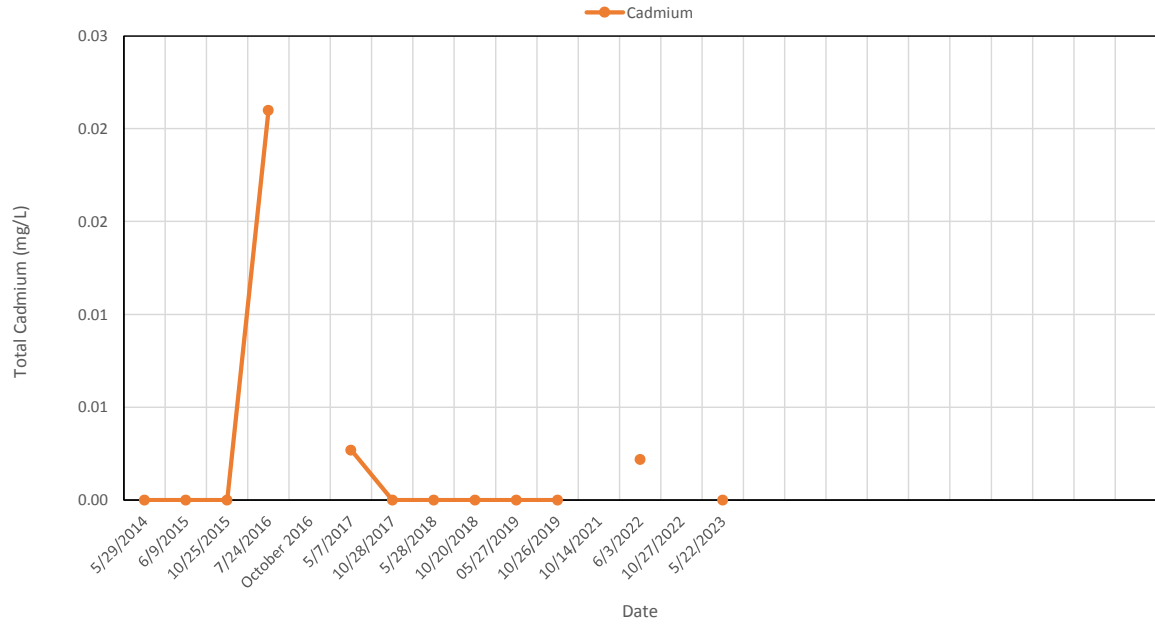
>VGES

Bold (italic) indicates value exceeds PAL

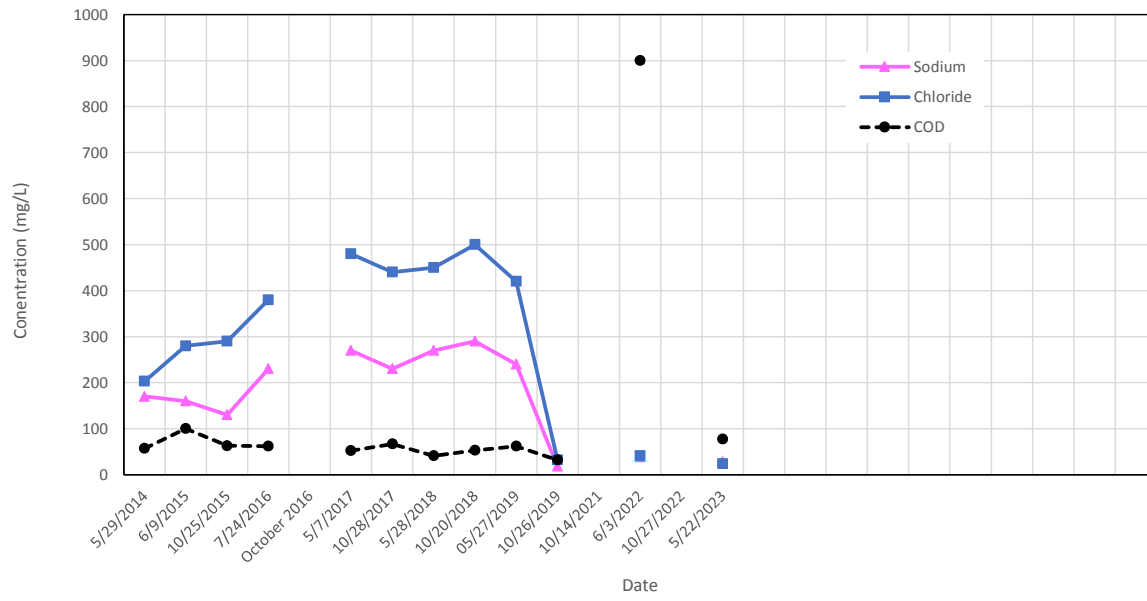
Brandon Landfill
MW-2C



Brandon Landfill
MW-2C



Brandon Landfill
MW-2C



Brandon Closed Landfill

MW-3

| PARAMETER | Sample Date: | 2012 | 4/22/2013 | 10/1/2013 | 5/29/2014 | 6/9/2015 | 10/25/2015 | 7/24/2016 | October 2016 | 5/7/2017 | 10/28/2017 | 5/28/2018 | 10/20/2018 | 05/27/2019 | 10/26/2019 | VGES | PAL |
|-------------------------------------|--------------|------|-----------|-----------|--------------|--------------|--------------|-----------|--------------|----------|------------|-----------|------------|--------------|--------------|-------|--------|
| VOCs (ug/L) | | | | | | | | | | | | | | | | | |
| 1,1-dichloroethane | | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 | - | - | - | - | - | - | - | - | 70 | 35 |
| Total Metals (mg/L) | | | | | | | | | | | | | | | | | |
| Arsenic | | - | - | - | <0.001 | 0.002 | 0.009 | Well | Well | Well | Well | Well | Well | No | No | 0.010 | 0.001 |
| Cadmium | | - | - | - | <0.002 | <0.002 | <0.002 | not | not | not | not | not | not | Sample | Sample | 0.005 | 0.001 |
| Chromium | | - | - | - | <0.005 | <0.005 | <0.0052 | sampled | sampled | sampled | sampled | sampled | sampled | - | - | 0.100 | 0.050 |
| Copper | | - | - | - | <0.020 | <0.020 | <0.022 | - | - | - | - | - | - | Insufficient | Insufficient | 1.300 | 0.650 |
| Iron | | - | - | - | 0.41 | 9.1 | 29 | - | - | - | - | - | - | Amount | Amount | - | - |
| Lead | | - | - | - | <0.001 | <0.001 | 0.008 | - | - | - | - | - | - | of Water | of Water | 0.015 | 0.002 |
| Manganese | | - | - | - | <0.020 | 1.1 | 1.2 | - | - | - | - | - | - | in Well | in Well | 0.300 | 0.150 |
| Mercury | | - | - | - | <0.0002 | <0.0002 | <0.0002 | - | - | - | - | - | - | Column | Column | 0.002 | 0.0005 |
| Nickel | | - | - | - | 0.005 | <0.005 | 0.016 | - | - | - | - | - | - | - | - | 0.100 | 0.050 |
| Sodium | | - | - | - | 23 | 20 | 15 | - | - | - | - | - | - | - | - | - | - |
| Zinc | | - | - | - | 0.020 | <0.020 | 0.024 | - | - | - | - | - | - | - | - | - | - |
| Other Analytes (mg/L) | | | | | | | | | | | | | | | | | |
| Chloride | | - | - | - | 34 | 31 | 30 | - | - | - | - | - | - | - | - | - | - |
| COD | | - | - | - | 11 | 34 | 34 | - | - | - | - | - | - | - | - | - | - |
| Field Measurements (units as noted) | | | | | | | | | | | | | | | | | |
| pH (std units) | | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Temperature (deg C) | | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Conductivity (uS) | | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - |
| Water Level (feet btoc) | | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - |

| PARAMETER | Sample Date: | 10/14/2021 | 6/3/2022 | 10/27/2022 | 5/22/2023 | | | | | | | | | | | VGES | PAL |
|--------------------------------------|--------------|-------------|--------------|---------------|---------------|--|--|--|--|--|--|--|--|--|--|-------|--------|
| VOCs (ug/L) | | | | | | | | | | | | | | | | | |
| 1,1-dichloroethane | | <0.5 | <0.5 | <1.0 | <1.0 | | | | | | | | | | | 70 | 35 |
| Total Metals (mg/L) | | | | | | | | | | | | | | | | | |
| Arsenic | | 0.17 | 0.015 | 0.030 | 0.0075 | | | | | | | | | | | 0.010 | 0.001 |
| Cadmium | | <0.005 | <0.001 | <0.0020 | <0.0020 | | | | | | | | | | | 0.005 | 0.001 |
| Chromium | | 0.32 | 0.022 | <0.050 | 0.0110 | | | | | | | | | | | 0.100 | 0.050 |
| Copper | | 0.82 | 0.062 | <0.20 | 0.021 | | | | | | | | | | | 1.300 | 0.650 |
| Iron | | 370 | 47 | 57 | 18 | | | | | | | | | | | | |
| Lead | | 0.51 | 0.036 | 0.0505 | 0.0151 | | | | | | | | | | | 0.015 | 0.002 |
| Manganese | | 25 | 1.5 | 2.5 | 0.55 | | | | | | | | | | | 0.300 | 0.150 |
| Mercury | | <0.001 | <0.0001 | <0.0002 | <0.0002 | | | | | | | | | | | 0.002 | 0.0005 |
| Nickel | | 0.49 | 0.036 | 0.0538 | 0.0140 | | | | | | | | | | | 0.100 | 0.050 |
| Sodium | | 26 | 28 | 23 | 26 | | | | | | | | | | | - | - |
| Zinc | | 1.4 | 0.11 | <0.20 | 0.042 | | | | | | | | | | | - | - |
| Other Analytes (mg/L) | | | | | | | | | | | | | | | | | |
| Chloride | | 34 | 56 | 98 | 44 | | | | | | | | | | | - | - |
| COD | | <10 | <10 | 220 | 79 | | | | | | | | | | | - | - |
| PFAS (ng/L) | | | | | | | | | | | | | | | | | |
| Perfluorohexanesulfonic acid (PFHxS) | | - | - | - | <1.8 | | | | | | | | | | | | |
| Perfluoroheptanoic acid (PFHpA) | | - | - | - | <1.8 | | | | | | | | | | | | |
| Perfluorooctanoic acid (PFOA) | | - | - | - | <1.8 | | | | | | | | | | | | |
| Perfluorooctanesulfonic acid (PFOS) | | - | - | - | 5.2 | | | | | | | | | | | | |
| Perfluorononanoic acid (PFNA) | | - | - | - | <1.8 | | | | | | | | | | | | |
| Total Regulated PFAS | | - | - | - | 5.2 | | | | | | | | | | | 20 | 2 |
| Total Non-Regulated PFAS | | - | - | - | ND | | | | | | | | | | | - | - |
| Field Measurements (units as noted) | | | | | | | | | | | | | | | | | |
| pH (std units) | | 6.8 | 7.14 | 7.27 | 7.57 | | | | | | | | | | | - | - |
| Temperature (deg C) | | 11.9 | 12.3 | 10.7 | 13.0 | | | | | | | | | | | - | - |
| Spec. Conductivity (uS/cm) | | 740 | 773 | 378.6 | 715 | | | | | | | | | | | - | - |
| Water Level (feet btoc) | | 33.69 | 30.78 | 33.60 | 31.50 | | | | | | | | | | | - | - |

Data prior to 2021 collected by others and obtained from 2019 reports prepared by EIV Technical Services

Data 2021-present collected by KAS, Inc.

Only detected or targeted VOCs are depicted

All values reported in units noted above

"-" = Not Analyzed, No Information or No Applicable Standard

ND = None Detected

<X = None Detected above Detection Limit (X)

VGES = Vermont Groundwater Enforcement Standard (GWPRS 07/06/2019)

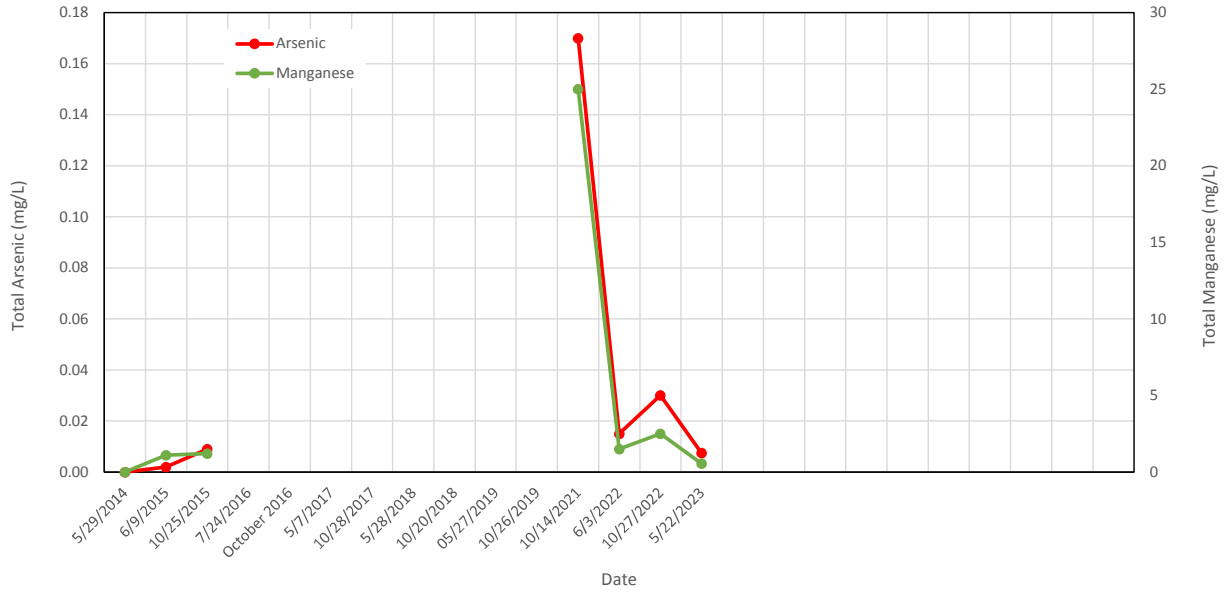
PAL = Vermont Preventive Action Level (GWPRS 07/06/2019)

Detections are **bolded**

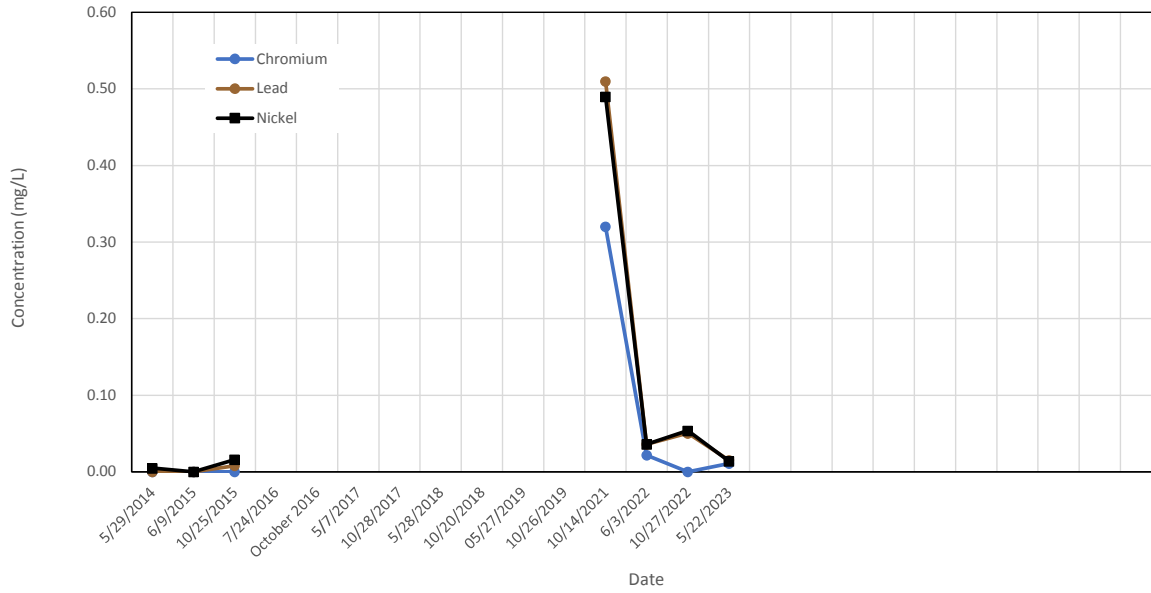
>VGES

bold (italic) indicates value exceeds PAL

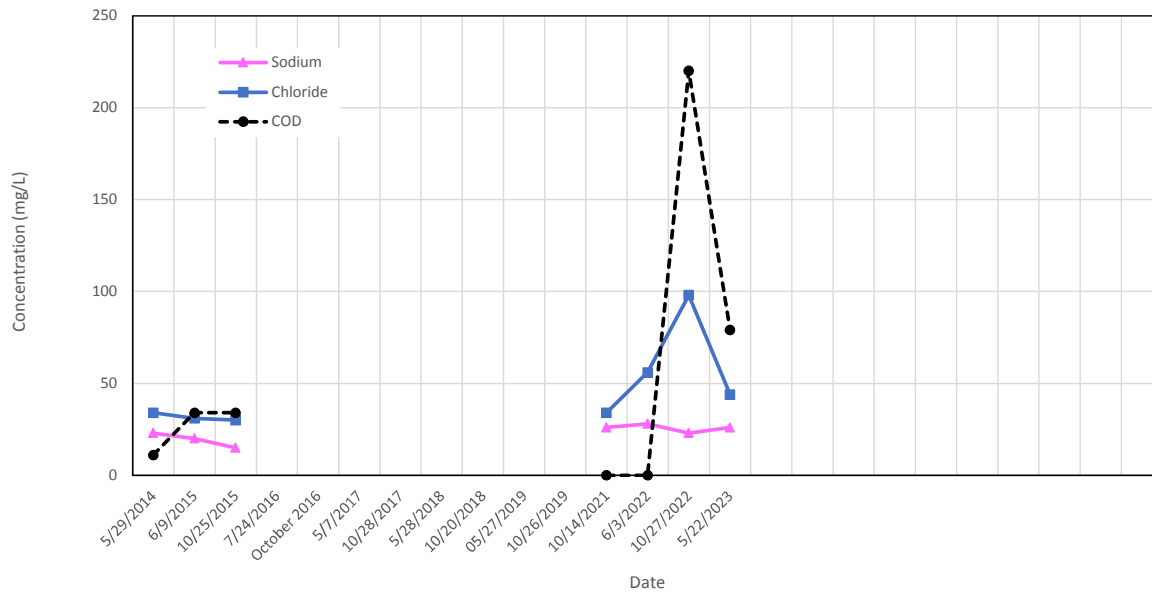
Brandon Landfill MW-3



Brandon Landfill MW-3



Brandon Landfill
MW-3



**Brandon Closed Landfill
MW-5**

| PARAMETER | Sample Date: | 2012 | 4/22/2013 | 10/1/2013 | 5/29/2014 | 6/9/2015 | 10/25/2015 | 7/24/2016 | October 2016 | 5/7/2017 | 10/28/2017 | 5/28/2018 | 10/20/2018 | 05/27/2019 | 10/26/2019 | VGES | PAL |
|--|--------------|------|-----------|------------|--------------|---------------|------------|--------------|--------------|---------------|---------------|---------------|---------------|-------------|---------------|-------|--------|
| VOCs (ug/L) | | | | | | | | | | | | | | | | | |
| 1,1-dichloroethane | | <1.0 | - | 1.1 | <1.0 | <1.0 | - | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 | <1 | <1 | 70 | 35 |
| Diethyl Ether | | - | - | 6.0 | <5.0 | <5.0 | - | <5.0 | - | <5.0 | - | - | - | - | - | - | - |
| Total Metals (mg/L) | | | | | | | | | | | | | | | | | |
| Arsenic | | - | Well | - | 0.004 | 0.003 | Well | 0.032 | Data | 0.006 | 0.006 | 0.0040 | 0.0011 | <0.0010 | 0.121 | 0.010 | 0.001 |
| Cadmium | | - | Not | - | <0.002 | <0.002 | not | 0.010 | not | 0.0061 | 0.0061 | 0.0083 | 0.0027 | <0.0020 | <0.0020 | 0.005 | 0.001 |
| Chromium | | - | Sampled | - | <0.005 | <0.005 | sampled | 0.020 | available | 0.0056 | 0.0056 | <0.0050 | <0.0050 | <0.0050 | <0.0050 | 0.100 | 0.050 |
| Copper | | - | - | - | <0.020 | <0.020 | - | 0.076 | - | <0.020 | <0.020 | <0.020 | <0.020 | <0.020 | <0.020 | 1.300 | 0.650 |
| Iron | | - | - | - | 31 | 16 | - | 120 | - | 21 | 21 | 13 | 4 | 0.22 | 48 | - | - |
| Lead | | - | - | - | 0.003 | 0.003 | - | 0.044 | - | 0.0080 | 0.0080 | 0.0080 | 0.0022 | <0.0010 | <0.0010 | 0.015 | 0.002 |
| Manganese | | - | - | - | 1.4 | 1.3 | - | 2.6 | - | 0.78 | 0.78 | 1.2 | 0.38 | 1.0 | 0.89 | 0.300 | 0.150 |
| Mercury | | - | - | - | <0.0002 | <0.0002 | - | <0.0002 | - | <0.0002 | <0.0002 | <0.0002 | <0.0002 | <0.0002 | <0.0002 | 0.002 | 0.0005 |
| Nickel | | - | - | - | 0.007 | 0.0066 | - | 0.025 | - | 0.0084 | 0.0084 | 0.0082 | <0.0050 | <0.0050 | 0.0077 | 0.100 | 0.050 |
| Sodium | | - | - | - | 26 | 21 | - | 18 | - | 24 | 24 | 25 | 31 | 33 | 150 | - | - |
| Zinc | | - | - | - | 0.020 | 0.020 | - | 0.11 | - | 0.026 | 0.026 | <0.020 | <0.020 | <0.020 | <0.020 | - | - |
| Other Analytes (mg/L) | | | | | | | | | | | | | | | | | |
| Chloride | | - | - | - | 38 | 32 | - | 33 | - | 42 | 43 | 40 | 64 | 64 | 260 | - | - |
| COD | | - | - | - | 30 | 25 | - | 54 | - | 31 | 13 | 30 | <10 | <10 | 45 | - | - |
| Field Measurements (units as noted) | | | | | | | | | | | | | | | | | |
| pH (std units) | | - | - | - | - | - | - | 6.4 | - | 6.0 | 6.1 | 6.2 | 6.2 | 6.3 | 6.4 | - | - |
| Temperature (deg C) | | - | - | - | - | - | - | 15 | - | 10.7 | 10.6 | 10.8 | 12.0 | 11.8 | 11.9 | - | - |
| Conductivity (uS) | | - | - | - | - | - | - | - | - | 1,160 | 1,090 | 1,080 | 1,120 | 1,080 | 1,100 | - | - |
| Water Level (feet btoc) | | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - | - |

| PARAMETER | Sample Date: | 10/14/2021 | 6/3/2022 | 10/27/2022 | 5/22/2023 | | | | | | | | | | | VGES | PAL |
|--|--------------|--------------|---------------|---------------|-------------|--|--|--|--|--|--|--|--|--|--|-------|--------|
| VOCs (ug/L) | | | | | | | | | | | | | | | | | |
| 1,1-dichloroethane | | - | <0.5 | <1.0 | <1.0 | | | | | | | | | | | 70 | 35 |
| Total Metals (mg/L) | | | | | | | | | | | | | | | | | |
| Arsenic | | No | 0.0042 | 0.0170 | <0.0010 | | | | | | | | | | | 0.010 | 0.001 |
| Cadmium | | Sample | 0.0016 | <0.0020 | <0.0020 | | | | | | | | | | | 0.005 | 0.001 |
| Chromium | | | 0.0013 | 0.0052 | <0.0050 | | | | | | | | | | | 0.100 | 0.050 |
| Copper | | Well | 0.0053 | 0.024 | <0.020 | | | | | | | | | | | 1.300 | 0.650 |
| Iron | | Inaccessible | 8.0 | 30 | 0.52 | | | | | | | | | | | - | - |
| Lead | | Due to | 0.0029 | 0.0175 | <0.0010 | | | | | | | | | | | 0.015 | 0.002 |
| Manganese | | Lock | 1.9 | 1.3 | 2.1 | | | | | | | | | | | 0.300 | 0.150 |
| Mercury | | | <0.0001 | <0.0002 | <0.0002 | | | | | | | | | | | 0.002 | 0.0005 |
| Nickel | | | 0.0074 | 0.0197 | <0.0050 | | | | | | | | | | | 0.100 | 0.050 |
| Sodium | | | 30 | 34 | 34 | | | | | | | | | | | - | - |
| Zinc | | | 0.014 | 0.047 | <0.020 | | | | | | | | | | | - | - |
| Other Analytes (mg/L) | | | | | | | | | | | | | | | | | |
| Chloride | | - | 52 | 59 | 59 | | | | | | | | | | | - | - |
| COD | | - | <10 | 58 | 35 | | | | | | | | | | | - | - |
| PFAS (ng/L) | | | | | | | | | | | | | | | | | |
| Perfluorohexanesulfonic acid (PFHxS) | | - | - | - | 17 | | | | | | | | | | | | |
| Perfluoroheptanoic acid (PFHpA) | | - | - | - | 5.1 | | | | | | | | | | | | |
| Perfluorooctanoic acid (PFOA) | | - | - | - | 30 | | | | | | | | | | | | |
| Perfluorooctanesulfonic acid (PFOS) | | - | - | - | 14 | | | | | | | | | | | | |
| Perfluorononanoic acid (PFNA) | | - | - | - | <1.9 | | | | | | | | | | | | |
| Total Regulated PFAS | | - | - | - | 66.1 | | | | | | | | | | | 20 | 2 |
| Total Non-Regulated PFAS | | - | - | - | 14.8 | | | | | | | | | | | - | - |
| Field Measurements (units as noted) | | | | | | | | | | | | | | | | | |
| pH (std units) | | - | 6.58 | 7.18 | 6.69 | | | | | | | | | | | - | - |
| Temperature (deg C) | | - | 13.2 | 11.7 | 10.1 | | | | | | | | | | | - | - |
| Spec. Conductivity (uS/cm) | | - | 1,109 | - | 826 | | | | | | | | | | | - | - |
| Water Level (feet btoc) | | - | 4.79 | 4.97 | 4.79 | | | | | | | | | | | - | - |

Data prior to 2021 collected by others and obtained from 2019 reports prepared by EIV Technical Services

Data 2021-present collected by KAS, Inc.

Only detected or targeted VOCs are depicted

All values reported in units noted above

"-" = Not Analyzed, No Information or No Applicable Standard

ND = None Detected

<X = None Detected above Detection Limit (X)

VGES = Vermont Groundwater Enforcement Standard (GWPRS 07/06/2019)

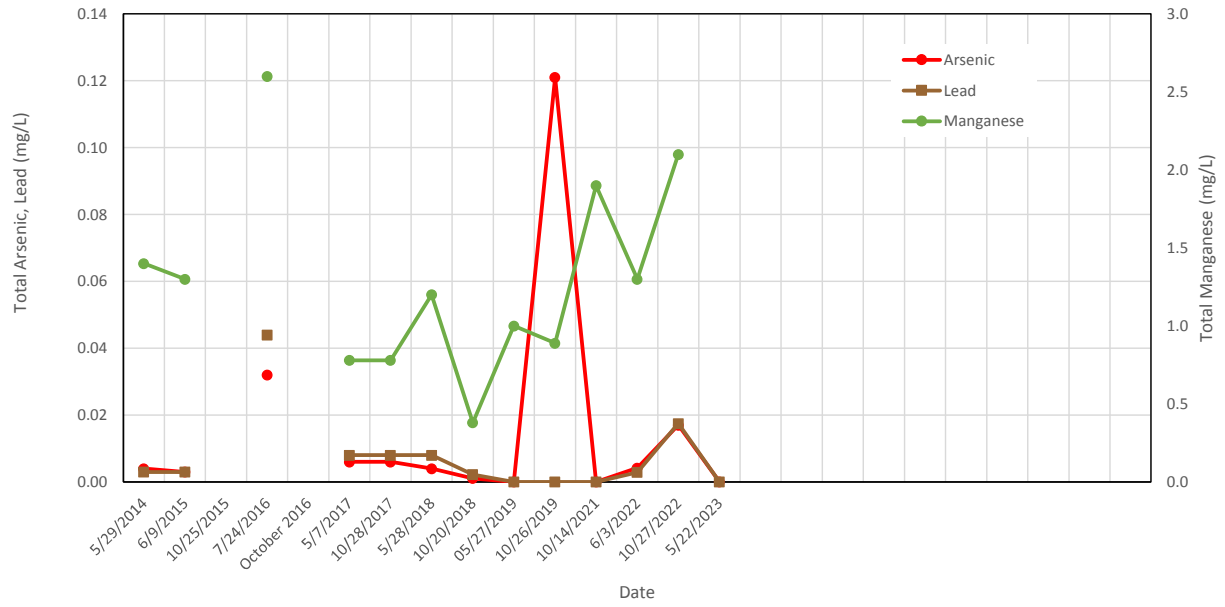
PAL = Vermont Preventive Action Level (GWPRS 07/06/2019)

Detections are **bolded**

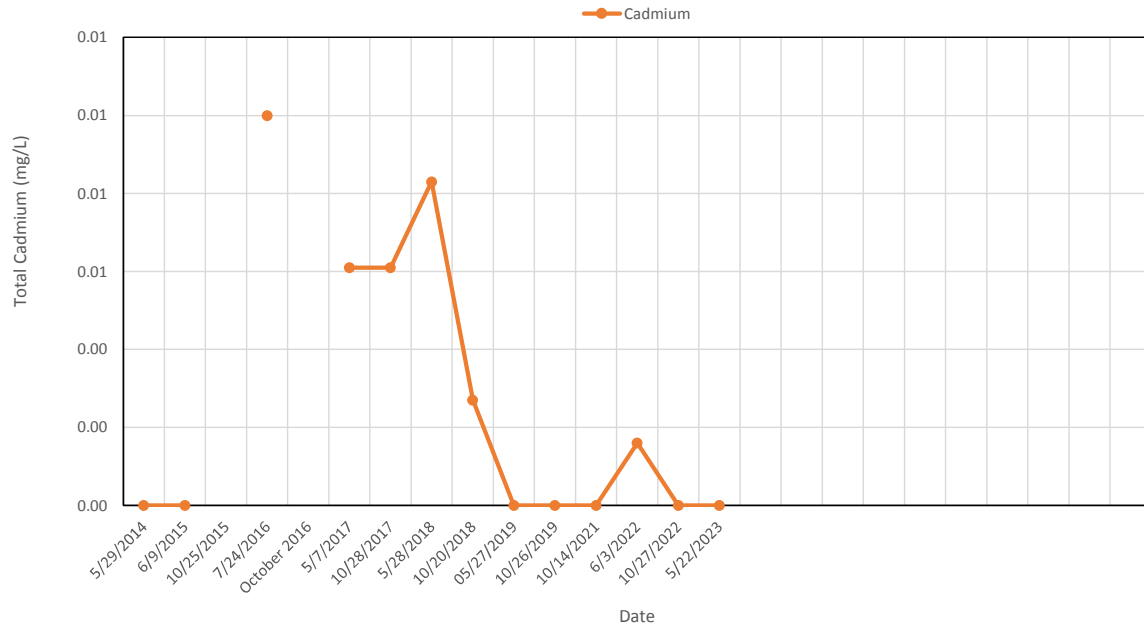
>VGES

Italic (italic) indicates value exceeds PAL

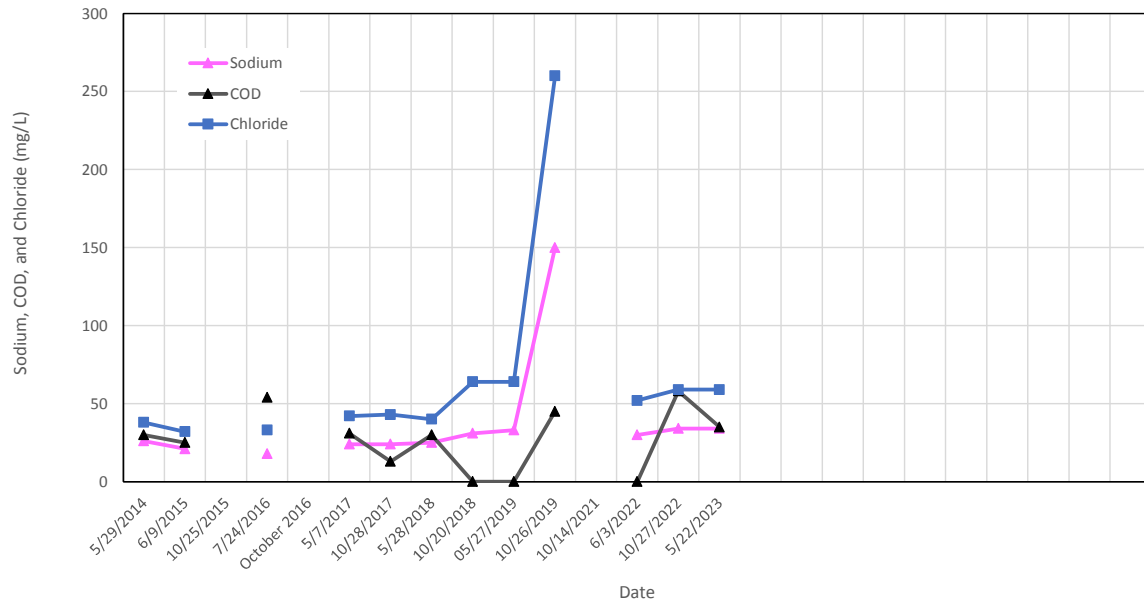
Brandon Landfill MW-5



Brandon Landfill MW-5



Brandon Landfill
MW-5



Brandon Closed Landfill
Quality Assurance/Quality Control Samples

| <i>Sample ID:</i> <i>Sample Date:</i> | Trip Blank 5/22/2023 | Duplicate 5/22/2023 | MW-5 5/22/2023 | RPD (%) |
|--|-------------------------|------------------------|-------------------|---------|
| PARAMETER | | | | |
| VOCs (ug/L) | | | | |
| Total VOCs | ND | ND | ND | - |
| Total Metals (mg/L) | | | | |
| Arsenic | - | <0.0010 | <0.0010 | - |
| Cadmium | - | <0.0020 | <0.0020 | - |
| Chromium | - | <0.0050 | <0.0050 | - |
| Copper | - | <0.020 | <0.020 | - |
| Iron | - | 0.51 | 0.52 | 1.9 |
| Lead | - | <0.0010 | <0.0010 | - |
| Manganese | - | 2.1 | 2.1 | 0.0 |
| Mercury | - | <0.0002 | <0.0002 | - |
| Nickel | - | <0.0050 | <0.0050 | - |
| Sodium | - | 35 | 34 | 2.9 |
| Zinc | - | <0.020 | <0.020 | - |
| Other Analytes (mg/L) | | | | |
| Chloride | - | 58 | 59 | 1.7 |
| COD | - | 41 | 35 | 15.8 |

Only detected or targeted VOCs are depicted

All values reported in units noted above

"-" = Not Analyzed, RPD could not be calculated due to non-detects or No Applicable Standard

ND = None Detected

<X = None Detected above Detection Limit (X)

RPD = The results of the laboratory analysis of the duplicate sample were analyzed using a relative percent difference (RPD) analysis. The RPD is defined as 100 times the difference in reported concentration between sample and duplicate, divided by the mean of the two samples. A small RPD indicates good correlation between sample and duplicate.



APPENDIX C

Laboratory Reports



Laboratory Report

| | |
|---------------------|--------|
| KAS, Inc | 100306 |
| PO Box 787 | |
| Williston, VT 05495 | |
| Atten: Clare Santos | |

PROJECT: Brandon Landfill
 WORK ORDER: **2305-14321**
 DATE RECEIVED: May 26, 2023
 DATE REPORTED: June 26, 2023
 SAMPLER: HG/KG

Enclosed please find the results of the analyses performed for the samples referenced on the attached chain of custody. All required method quality control elements including instrument calibration were performed in accordance with method requirements and determined to be acceptable unless otherwise noted.

The column labeled Lab/Tech in the accompanying report denotes the laboratory facility where the testing was performed and the technician who conducted the assay. A "W" designates the Williston, VT lab under NELAC certification ELAP 11263; "R" designates the Lebanon, NH facility under certification NH 2037 and "N" the Plattsburgh, NY lab under certification ELAP 11892. "Sub" indicates the testing was performed by a subcontracted laboratory. The accreditation status of the subcontracted lab is referenced in the corresponding NELAC and Qual fields.

The NELAC column also denotes the accreditation status of each laboratory for each reported parameter. "A" indicates the referenced laboratory is NELAC accredited for the parameter reported. "N" indicates the laboratory is not accredited. "U" indicates that NELAC does not offer accreditation for that parameter in that specific matrix. Test results denoted with an "A" meet all National Environmental Laboratory Accreditation Program requirements except where denoted by pertinent data qualifiers. Test results are representative of the samples as they were received at the laboratory

Endyne, Inc. warrants, to the best of its knowledge and belief, the accuracy of the analytical test results contained in this report, but makes no other warranty, expressed or implied, especially no warranties of merchantability or fitness for a particular purpose.

Reviewed by:

Harry B. Locker, Ph.D.
 Laboratory Director

www.endynelabs.com



160 James Brown Dr., Williston, VT 05495
 Ph 802-879-4333 Fax 802-879-7103

56 Etna Road, Lebanon, NH 03755
 Ph 603-678-4891 Fax 603-678-4893



Laboratory Report

REPORT DATE: 6/26/2023

CLIENT: KAS, Inc
PROJECT: Brandon Landfill

WORK ORDER: 2305-14321
DATE RECEIVED: 05/26/2023

001 Site: MW-1 Date Sampled: 5/22/23 Time: 12:25

| Parameter | Result | Units | Method | Analysis Date | Lab/Tech | NELAC | Qual. |
|------------------|----------|-------|-----------|---------------|----------|-------|-------|
| Chloride | 77 | mg/L | EPA 300.0 | 5/30/23 15:52 | W ECM | A | |
| COD | 74 | mg/L | EPA 410.4 | 6/6/23 | N WEP | A | |
| Metals Digestion | Digested | | EPA 3015A | 6/22/23 | W MGT | A | |
| Arsenic, Total | 0.0014 | mg/L | EPA 6020B | 6/24/23 0:04 | W MGT | A | |
| Cadmium, Total | < 0.0020 | mg/L | EPA 6020B | 6/24/23 0:04 | W MGT | A | |
| Chromium, Total | < 0.0050 | mg/L | EPA 6020B | 6/24/23 0:04 | W MGT | A | |
| Copper, Total | < 0.020 | mg/L | EPA 6020B | 6/24/23 0:04 | W MGT | A | |
| Iron, Total | 2.4 | mg/L | EPA 6010C | 6/23/23 | W MGT | A | |
| Lead, Total | < 0.0010 | mg/L | EPA 6020B | 6/24/23 0:04 | W MGT | A | |
| Manganese, Total | 0.25 | mg/L | EPA 6020B | 6/24/23 0:04 | W MGT | A | |
| Mercury, Total | < 0.0002 | mg/L | EPA 6020B | 6/24/23 0:04 | W MGT | N | |
| Nickel, Total | < 0.0050 | mg/L | EPA 6020B | 6/24/23 0:04 | W MGT | A | |
| Sodium, Total | 41 | mg/L | EPA 6010C | 6/23/23 | W MGT | A | |
| Zinc, Total | < 0.020 | mg/L | EPA 6020B | 6/24/23 0:04 | W MGT | A | |

002 Site: MW-2C Date Sampled: 5/22/23 Time: 9:30

| Parameter | Result | Units | Method | Analysis Date | Lab/Tech | NELAC | Qual. |
|------------------|----------|-------|-----------|---------------|----------|-------|-------|
| Chloride | 24 | mg/L | EPA 300.0 | 5/30/23 16:12 | W ECM | A | |
| COD | 77 | mg/L | EPA 410.4 | 6/6/23 | N WEP | A | |
| Metals Digestion | Digested | | EPA 3015A | 6/22/23 | W MGT | A | |
| Arsenic, Total | < 0.0010 | mg/L | EPA 6020B | 6/24/23 0:09 | W MGT | A | |
| Cadmium, Total | < 0.0020 | mg/L | EPA 6020B | 6/24/23 0:09 | W MGT | A | |
| Chromium, Total | < 0.0050 | mg/L | EPA 6020B | 6/24/23 0:09 | W MGT | A | |
| Copper, Total | < 0.020 | mg/L | EPA 6020B | 6/24/23 0:09 | W MGT | A | |
| Iron, Total | 2.4 | mg/L | EPA 6010C | 6/23/23 | W MGT | A | |
| Lead, Total | 0.0036 | mg/L | EPA 6020B | 6/24/23 0:09 | W MGT | A | |
| Manganese, Total | 1.3 | mg/L | EPA 6020B | 6/24/23 0:09 | W MGT | A | |
| Mercury, Total | < 0.0002 | mg/L | EPA 6020B | 6/24/23 0:09 | W MGT | N | |
| Nickel, Total | 0.0081 | mg/L | EPA 6020B | 6/24/23 0:09 | W MGT | A | |
| Sodium, Total | 28 | mg/L | EPA 6010C | 6/23/23 | W MGT | A | |
| Zinc, Total | < 0.020 | mg/L | EPA 6020B | 6/24/23 0:09 | W MGT | A | |

003 Site: MW-3 Date Sampled: 5/22/23 Time: 11:46

| Parameter | Result | Units | Method | Analysis Date | Lab/Tech | NELAC | Qual. |
|------------------|----------|-------|-----------|---------------|----------|-------|-------|
| Chloride | 44 | mg/L | EPA 300.0 | 5/30/23 17:32 | W ECM | A | |
| COD | 79 | mg/L | EPA 410.4 | 6/6/23 | N WEP | A | |
| Metals Digestion | Digested | | EPA 3015A | 6/22/23 | W MGT | A | |
| Arsenic, Total | 0.0075 | mg/L | EPA 6020B | 6/24/23 0:14 | W MGT | A | |
| Cadmium, Total | < 0.0020 | mg/L | EPA 6020B | 6/24/23 0:14 | W MGT | A | |
| Chromium, Total | 0.0110 | mg/L | EPA 6020B | 6/24/23 0:14 | W MGT | A | |
| Copper, Total | 0.021 | mg/L | EPA 6020B | 6/24/23 0:14 | W MGT | A | |
| Iron, Total | 18 | mg/L | EPA 6010C | 6/23/23 | W MGT | A | |
| Lead, Total | 0.0151 | mg/L | EPA 6020B | 6/24/23 0:14 | W MGT | A | |
| Manganese, Total | 0.55 | mg/L | EPA 6020B | 6/24/23 0:14 | W MGT | A | |

Laboratory Report

REPORT DATE: 6/26/2023

CLIENT: KAS, Inc
PROJECT: Brandon Landfill

WORK ORDER: 2305-14321
DATE RECEIVED: 05/26/2023

003 Site: MW-3 Date Sampled: 5/22/23 Time: 11:46

| Parameter | Result | Units | Method | Analysis Date | Lab/Tech | NELAC | Qual. |
|----------------|----------|-------|-----------|---------------|----------|-------|-------|
| Mercury, Total | < 0.0002 | mg/L | EPA 6020B | 6/24/23 0:14 | W MGT | N | |
| Nickel, Total | 0.0140 | mg/L | EPA 6020B | 6/24/23 0:14 | W MGT | A | |
| Sodium, Total | 26 | mg/L | EPA 6010C | 6/23/23 | W MGT | A | |
| Zinc, Total | 0.042 | mg/L | EPA 6020B | 6/24/23 0:14 | W MGT | A | |

004 Site: MW-5 Date Sampled: 5/22/23 Time: 10:35

| Parameter | Result | Units | Method | Analysis Date | Lab/Tech | NELAC | Qual. |
|------------------|----------|-------|-----------|---------------|----------|-------|-------|
| Chloride | 59 | mg/L | EPA 300.0 | 5/30/23 17:52 | W ECM | A | |
| COD | 35 | mg/L | EPA 410.4 | 6/6/23 | N WEP | A | |
| Metals Digestion | Digested | | EPA 3015A | 6/22/23 | W MGT | A | |
| Arsenic, Total | < 0.0010 | mg/L | EPA 6020B | 6/24/23 0:19 | W MGT | A | |
| Cadmium, Total | < 0.0020 | mg/L | EPA 6020B | 6/24/23 0:19 | W MGT | A | |
| Chromium, Total | < 0.0050 | mg/L | EPA 6020B | 6/24/23 0:19 | W MGT | A | |
| Copper, Total | < 0.020 | mg/L | EPA 6020B | 6/24/23 0:19 | W MGT | A | |
| Iron, Total | 0.52 | mg/L | EPA 6010C | 6/23/23 | W MGT | A | B |
| Lead, Total | < 0.0010 | mg/L | EPA 6020B | 6/24/23 0:19 | W MGT | A | |
| Manganese, Total | 2.1 | mg/L | EPA 6020B | 6/24/23 0:19 | W MGT | A | |
| Mercury, Total | < 0.0002 | mg/L | EPA 6020B | 6/24/23 0:19 | W MGT | N | |
| Nickel, Total | < 0.0050 | mg/L | EPA 6020B | 6/24/23 0:19 | W MGT | A | |
| Sodium, Total | 34 | mg/L | EPA 6010C | 6/23/23 | W MGT | A | |
| Zinc, Total | < 0.020 | mg/L | EPA 6020B | 6/24/23 0:19 | W MGT | A | |

005 Site: Duplicate Date Sampled: 5/22/23 Time: 10:35

| Parameter | Result | Units | Method | Analysis Date | Lab/Tech | NELAC | Qual. |
|------------------|----------|-------|-----------|---------------|----------|-------|-------|
| Chloride | 58 | mg/L | EPA 300.0 | 5/30/23 18:12 | W ECM | A | |
| COD | 41 | mg/L | EPA 410.4 | 6/6/23 | N WEP | A | |
| Metals Digestion | Digested | | EPA 3015A | 6/22/23 | W MGT | A | |
| Arsenic, Total | < 0.0010 | mg/L | EPA 6020B | 6/24/23 0:24 | W MGT | A | |
| Cadmium, Total | < 0.0020 | mg/L | EPA 6020B | 6/24/23 0:24 | W MGT | A | |
| Chromium, Total | < 0.0050 | mg/L | EPA 6020B | 6/24/23 0:24 | W MGT | A | |
| Copper, Total | < 0.020 | mg/L | EPA 6020B | 6/24/23 0:24 | W MGT | A | |
| Iron, Total | 0.51 | mg/L | EPA 6010C | 6/23/23 | W MGT | A | B |
| Lead, Total | < 0.0010 | mg/L | EPA 6020B | 6/24/23 0:24 | W MGT | A | |
| Manganese, Total | 2.1 | mg/L | EPA 6020B | 6/24/23 0:24 | W MGT | A | |
| Mercury, Total | < 0.0002 | mg/L | EPA 6020B | 6/24/23 0:24 | W MGT | N | |
| Nickel, Total | < 0.0050 | mg/L | EPA 6020B | 6/24/23 0:24 | W MGT | A | |
| Sodium, Total | 35 | mg/L | EPA 6010C | 6/23/23 | W MGT | A | |
| Zinc, Total | < 0.020 | mg/L | EPA 6020B | 6/24/23 0:24 | W MGT | A | |

Laboratory Report

REPORT DATE: 6/26/2023

CLIENT: KAS, Inc
PROJECT: Brandon Landfill

WORK ORDER: 2305-14321
DATE RECEIVED: 05/26/2023

TEST METHOD: EPA 8260C

| 001 | Site: MW-1 | Sampled: 5/22/23 12:25 | | Test Date: 6/2/23 | | W | TRP | | |
|--------------------------------|------------|------------------------|-------|-------------------|---------------------------|--------|------|-------|------|
| Parameter | Result | Unit | Nelac | Qual | Parameter | Result | Unit | Nelac | Qual |
| Dichlorodifluoromethane | < 5.0 | ug/L | A | | Chloromethane | < 3.0 | ug/L | A | |
| Vinyl chloride | < 0.5 | ug/L | A | | Bromomethane | < 0.5 | ug/L | A | |
| Chloroethane | < 5.0 | ug/L | A | | Trichlorofluoromethane | < 2.0 | ug/L | A | |
| Diethyl ether | < 5.0 | ug/L | N | | 1,1-Dichloroethene | < 0.7 | ug/L | A | |
| Acetone | < 10.0 | ug/L | A | | Carbon disulfide | < 5.0 | ug/L | A | |
| Methylene chloride | < 5.0 | ug/L | A | | t-Butanol | < 20.0 | ug/L | N | QA- |
| Methyl-t-butyl ether (MTBE) | < 2.0 | ug/L | A | | trans-1,2-Dichloroethene | < 1.0 | ug/L | A | |
| Di-isopropyl ether (DIPE) | < 2.0 | ug/L | N | | 1,1-Dichloroethane | < 1.0 | ug/L | A | |
| Ethyl-t-butyl ether (ETBE) | < 2.0 | ug/L | N | | 2-Butanone | < 10.0 | ug/L | A | |
| 2,2-Dichloropropane | < 1.0 | ug/L | N | | cis-1,2-Dichloroethene | < 1.0 | ug/L | A | |
| Bromochloromethane | < 0.8 | ug/L | A | | Chloroform | < 1.0 | ug/L | A | |
| Tetrahydrofuran | < 10.0 | ug/L | N | | 1,1,1-Trichloroethane | < 1.0 | ug/L | A | |
| Carbon tetrachloride | < 0.5 | ug/L | A | | 1,1-Dichloropropene | < 1.0 | ug/L | N | |
| Benzene | < 0.5 | ug/L | A | | t-Amylmethyl ether (TAME) | < 2.0 | ug/L | N | |
| 1,2-Dichloroethane | < 0.5 | ug/L | A | | Trichloroethene | < 0.5 | ug/L | A | |
| 1,2-Dichloropropane | < 0.5 | ug/L | A | | Dibromomethane | < 2.0 | ug/L | A | |
| Bromodichloromethane | < 0.5 | ug/L | A | | cis-1,3-Dichloropropene | < 1.0 | ug/L | A | |
| 4-Methyl-2-pentanone (MIBK) | < 10.0 | ug/L | A | | Toluene | < 1.0 | ug/L | A | |
| trans-1,3-Dichloropropene | < 1.0 | ug/L | A | | 1,1,2-Trichloroethane | < 1.0 | ug/L | A | |
| Tetrachloroethene | < 0.5 | ug/L | A | | 1,3-Dichloropropane | < 1.0 | ug/L | N | |
| 2-Hexanone | < 10.0 | ug/L | A | | Dibromochloromethane | < 1.0 | ug/L | A | |
| 1,2-Dibromoethane | < 2.0 | ug/L | A | | Chlorobenzene | < 1.0 | ug/L | A | |
| Ethylbenzene | < 1.0 | ug/L | A | | 1,1,1,2-Tetrachloroethane | < 2.0 | ug/L | A | |
| Xylenes, Total | < 2.0 | ug/L | A | | Styrene | < 1.0 | ug/L | A | |
| Bromoform | < 2.0 | ug/L | A | | Isopropylbenzene | < 1.0 | ug/L | A | |
| 1,1,2,2-Tetrachloroethane | < 2.0 | ug/L | A | | Bromobenzene | < 1.0 | ug/L | A | |
| n-Propylbenzene | < 1.0 | ug/L | A | | 1,2,3-Trichloropropane | < 2.0 | ug/L | A | |
| 2-Chlorotoluene | < 1.0 | ug/L | A | | 1,3,5-Trimethylbenzene | < 1.0 | ug/L | A | |
| 4-Chlorotoluene | < 1.0 | ug/L | A | | t-Butylbenzene | < 1.0 | ug/L | A | |
| 1,2,4-Trimethylbenzene | < 1.0 | ug/L | A | | s-Butylbenzene | < 1.0 | ug/L | A | |
| 4-Isopropyltoluene | < 1.0 | ug/L | A | | 1,3-Dichlorobenzene | < 1.0 | ug/L | A | |
| 1,4-Dichlorobenzene | < 1.0 | ug/L | A | | 1,2,3-Trimethylbenzene | < 1.0 | ug/L | U | |
| n-Butylbenzene | < 1.0 | ug/L | A | | 1,2-Dichlorobenzene | < 1.0 | ug/L | A | |
| 1,2-Dibromo-3-Chloropropane | < 2.0 | ug/L | A | | 1,2,4-Trichlorobenzene | < 2.0 | ug/L | A | |
| 1,3,5-Trichlorobenzene | < 2.0 | ug/L | N | | Hexachlorobutadiene | < 0.5 | ug/L | A | |
| Naphthalene | < 0.5 | ug/L | A | | 1,2,3-Trichlorobenzene | < 0.5 | ug/L | A | |
| Surr. 1 (Dibromofluoromethane) | 97 | % | A | | Surr. 2 (Toluene d8) | 99 | % | A | |
| Surr. 3 (4-Bromofluorobenzene) | 97 | % | A | | Unidentified Peaks | 0 | | U | |

CLIENT: KAS, Inc
PROJECT: Brandon Landfill

WORK ORDER: 2305-14321
DATE RECEIVED: 05/26/2023

TEST METHOD: EPA 8260C

| 002 | Site: MW-2C | Sampled: 5/22/23 9:30 | | Test Date: 6/2/23 | | W | TRP | | |
|--------------------------------|-------------|-----------------------|-------|-------------------|---------------------------|--------|------|-------|------|
| Parameter | Result | Unit | Nelac | Qual | Parameter | Result | Unit | Nelac | Qual |
| Dichlorodifluoromethane | < 5.0 | ug/L | A | | Chloromethane | < 3.0 | ug/L | A | |
| Vinyl chloride | < 0.5 | ug/L | A | | Bromomethane | < 0.5 | ug/L | A | |
| Chloroethane | < 5.0 | ug/L | A | | Trichlorofluoromethane | < 2.0 | ug/L | A | |
| Diethyl ether | 14.6 | ug/L | N | | 1,1-Dichloroethene | < 0.7 | ug/L | A | |
| Acetone | < 10.0 | ug/L | A | | Carbon disulfide | < 5.0 | ug/L | A | |
| Methylene chloride | < 5.0 | ug/L | A | | t-Butanol | < 20.0 | ug/L | N | QA- |
| Methyl-t-butyl ether (MTBE) | < 2.0 | ug/L | A | | trans-1,2-Dichloroethene | < 1.0 | ug/L | A | |
| Di-isopropyl ether (DIPE) | < 2.0 | ug/L | N | | 1,1-Dichloroethane | < 1.0 | ug/L | A | |
| Ethyl-t-butyl ether (ETBE) | < 2.0 | ug/L | N | | 2-Butanone | < 10.0 | ug/L | A | |
| 2,2-Dichloropropane | < 1.0 | ug/L | N | | cis-1,2-Dichloroethene | < 1.0 | ug/L | A | |
| Bromochloromethane | < 0.8 | ug/L | A | | Chloroform | < 1.0 | ug/L | A | |
| Tetrahydrofuran | < 10.0 | ug/L | N | | 1,1,1-Trichloroethane | < 1.0 | ug/L | A | |
| Carbon tetrachloride | < 0.5 | ug/L | A | | 1,1-Dichloropropene | < 1.0 | ug/L | N | |
| Benzene | < 0.5 | ug/L | A | | t-Amylmethyl ether (TAME) | < 2.0 | ug/L | N | |
| 1,2-Dichloroethane | < 0.5 | ug/L | A | | Trichloroethene | < 0.5 | ug/L | A | |
| 1,2-Dichloropropane | < 0.5 | ug/L | A | | Dibromomethane | < 2.0 | ug/L | A | |
| Bromodichloromethane | < 0.5 | ug/L | A | | cis-1,3-Dichloropropene | < 1.0 | ug/L | A | |
| 4-Methyl-2-pentanone (MIBK) | < 10.0 | ug/L | A | | Toluene | < 1.0 | ug/L | A | |
| trans-1,3-Dichloropropene | < 1.0 | ug/L | A | | 1,1,2-Trichloroethane | < 1.0 | ug/L | A | |
| Tetrachloroethene | < 0.5 | ug/L | A | | 1,3-Dichloropropane | < 1.0 | ug/L | N | |
| 2-Hexanone | < 10.0 | ug/L | A | | Dibromochloromethane | < 1.0 | ug/L | A | |
| 1,2-Dibromoethane | < 2.0 | ug/L | A | | Chlorobenzene | < 1.0 | ug/L | A | |
| Ethylbenzene | < 1.0 | ug/L | A | | 1,1,1,2-Tetrachloroethane | < 2.0 | ug/L | A | |
| Xylenes, Total | < 2.0 | ug/L | A | | Styrene | < 1.0 | ug/L | A | |
| Bromoform | < 2.0 | ug/L | A | | Isopropylbenzene | < 1.0 | ug/L | A | |
| 1,1,2,2-Tetrachloroethane | < 2.0 | ug/L | A | | Bromobenzene | < 1.0 | ug/L | A | |
| n-Propylbenzene | < 1.0 | ug/L | A | | 1,2,3-Trichloropropane | < 2.0 | ug/L | A | |
| 2-Chlorotoluene | < 1.0 | ug/L | A | | 1,3,5-Trimethylbenzene | < 1.0 | ug/L | A | |
| 4-Chlorotoluene | < 1.0 | ug/L | A | | t-Butylbenzene | < 1.0 | ug/L | A | |
| 1,2,4-Trimethylbenzene | < 1.0 | ug/L | A | | s-Butylbenzene | < 1.0 | ug/L | A | |
| 4-Isopropyltoluene | < 1.0 | ug/L | A | | 1,3-Dichlorobenzene | < 1.0 | ug/L | A | |
| 1,4-Dichlorobenzene | < 1.0 | ug/L | A | | 1,2,3-Trimethylbenzene | < 1.0 | ug/L | U | |
| n-Butylbenzene | < 1.0 | ug/L | A | | 1,2-Dichlorobenzene | < 1.0 | ug/L | A | |
| 1,2-Dibromo-3-Chloropropane | < 2.0 | ug/L | A | | 1,2,4-Trichlorobenzene | < 2.0 | ug/L | A | |
| 1,3,5-Trichlorobenzene | < 2.0 | ug/L | N | | Hexachlorobutadiene | < 0.5 | ug/L | A | |
| Naphthalene | < 0.5 | ug/L | A | | 1,2,3-Trichlorobenzene | < 0.5 | ug/L | A | |
| Surr. 1 (Dibromofluoromethane) | 98 | % | A | | Surr. 2 (Toluene d8) | 99 | % | A | |
| Surr. 3 (4-Bromofluorobenzene) | 98 | % | A | | Unidentified Peaks | 0 | | U | |

CLIENT: KAS, Inc
PROJECT: Brandon Landfill

WORK ORDER: 2305-14321
DATE RECEIVED: 05/26/2023

TEST METHOD: EPA 8260C

| 003 Site: MW-3 | | Sampled: 5/22/23 11:46 | | Test Date: 6/2/23 | | W TRP | | | |
|--------------------------------|--------|------------------------|-------|-------------------|---------------------------|--------|------|-------|------|
| Parameter | Result | Unit | Nelac | Qual | Parameter | Result | Unit | Nelac | Qual |
| Dichlorodifluoromethane | < 5.0 | ug/L | A | | Chloromethane | < 3.0 | ug/L | A | |
| Vinyl chloride | < 0.5 | ug/L | A | | Bromomethane | < 0.5 | ug/L | A | |
| Chloroethane | < 5.0 | ug/L | A | | Trichlorofluoromethane | < 2.0 | ug/L | A | |
| Diethyl ether | < 5.0 | ug/L | N | | 1,1-Dichloroethene | < 0.7 | ug/L | A | |
| Acetone | < 10.0 | ug/L | A | | Carbon disulfide | < 5.0 | ug/L | A | |
| Methylene chloride | < 5.0 | ug/L | A | | t-Butanol | < 20.0 | ug/L | N | QA- |
| Methyl-t-butyl ether (MTBE) | < 2.0 | ug/L | A | | trans-1,2-Dichloroethene | < 1.0 | ug/L | A | |
| Di-isopropyl ether (DIPE) | < 2.0 | ug/L | N | | 1,1-Dichloroethane | < 1.0 | ug/L | A | |
| Ethyl-t-butyl ether (ETBE) | < 2.0 | ug/L | N | | 2-Butanone | < 10.0 | ug/L | A | |
| 2,2-Dichloropropane | < 1.0 | ug/L | N | | cis-1,2-Dichloroethene | < 1.0 | ug/L | A | |
| Bromochloromethane | < 0.8 | ug/L | A | | Chloroform | < 1.0 | ug/L | A | |
| Tetrahydrofuran | < 10.0 | ug/L | N | | 1,1,1-Trichloroethane | < 1.0 | ug/L | A | |
| Carbon tetrachloride | < 0.5 | ug/L | A | | 1,1-Dichloropropene | < 1.0 | ug/L | N | |
| Benzene | < 0.5 | ug/L | A | | t-Amylmethyl ether (TAME) | < 2.0 | ug/L | N | |
| 1,2-Dichloroethane | < 0.5 | ug/L | A | | Trichloroethene | < 0.5 | ug/L | A | |
| 1,2-Dichloropropane | < 0.5 | ug/L | A | | Dibromomethane | < 2.0 | ug/L | A | |
| Bromodichloromethane | < 0.5 | ug/L | A | | cis-1,3-Dichloropropene | < 1.0 | ug/L | A | |
| 4-Methyl-2-pentanone (MIBK) | < 10.0 | ug/L | A | | Toluene | < 1.0 | ug/L | A | |
| trans-1,3-Dichloropropene | < 1.0 | ug/L | A | | 1,1,2-Trichloroethane | < 1.0 | ug/L | A | |
| Tetrachloroethene | < 0.5 | ug/L | A | | 1,3-Dichloropropane | < 1.0 | ug/L | N | |
| 2-Hexanone | < 10.0 | ug/L | A | | Dibromochloromethane | < 1.0 | ug/L | A | |
| 1,2-Dibromoethane | < 2.0 | ug/L | A | | Chlorobenzene | < 1.0 | ug/L | A | |
| Ethylbenzene | < 1.0 | ug/L | A | | 1,1,1,2-Tetrachloroethane | < 2.0 | ug/L | A | |
| Xylenes, Total | < 2.0 | ug/L | A | | Styrene | < 1.0 | ug/L | A | |
| Bromoform | < 2.0 | ug/L | A | | Isopropylbenzene | < 1.0 | ug/L | A | |
| 1,1,2,2-Tetrachloroethane | < 2.0 | ug/L | A | | Bromobenzene | < 1.0 | ug/L | A | |
| n-Propylbenzene | < 1.0 | ug/L | A | | 1,2,3-Trichloropropane | < 2.0 | ug/L | A | |
| 2-Chlorotoluene | < 1.0 | ug/L | A | | 1,3,5-Trimethylbenzene | < 1.0 | ug/L | A | |
| 4-Chlorotoluene | < 1.0 | ug/L | A | | t-Butylbenzene | < 1.0 | ug/L | A | |
| 1,2,4-Trimethylbenzene | < 1.0 | ug/L | A | | s-Butylbenzene | < 1.0 | ug/L | A | |
| 4-Isopropyltoluene | < 1.0 | ug/L | A | | 1,3-Dichlorobenzene | < 1.0 | ug/L | A | |
| 1,4-Dichlorobenzene | < 1.0 | ug/L | A | | 1,2,3-Trimethylbenzene | < 1.0 | ug/L | U | |
| n-Butylbenzene | < 1.0 | ug/L | A | | 1,2-Dichlorobenzene | < 1.0 | ug/L | A | |
| 1,2-Dibromo-3-Chloropropane | < 2.0 | ug/L | A | | 1,2,4-Trichlorobenzene | < 2.0 | ug/L | A | |
| 1,3,5-Trichlorobenzene | < 2.0 | ug/L | N | | Hexachlorobutadiene | < 0.5 | ug/L | A | |
| Naphthalene | < 0.5 | ug/L | A | | 1,2,3-Trichlorobenzene | < 0.5 | ug/L | A | |
| Surr. 1 (Dibromofluoromethane) | 100 | % | A | | Surr. 2 (Toluene d8) | 99 | % | A | |
| Surr. 3 (4-Bromofluorobenzene) | 99 | % | A | | Unidentified Peaks | 4 | | U | |

CLIENT: KAS, Inc
PROJECT: Brandon Landfill

WORK ORDER: 2305-14321
DATE RECEIVED: 05/26/2023

TEST METHOD: EPA 8260C

| 004 Site: MW-5 | | Sampled: 5/22/23 10:35 | | Test Date: 6/2/23 | | W TRP | | | |
|--------------------------------|--------|------------------------|-------|-------------------|---------------------------|--------|------|-------|------|
| Parameter | Result | Unit | Nelac | Qual | Parameter | Result | Unit | Nelac | Qual |
| Dichlorodifluoromethane | < 5.0 | ug/L | A | | Chloromethane | < 3.0 | ug/L | A | |
| Vinyl chloride | < 0.5 | ug/L | A | | Bromomethane | < 0.5 | ug/L | A | |
| Chloroethane | < 5.0 | ug/L | A | | Trichlorofluoromethane | < 2.0 | ug/L | A | |
| Diethyl ether | < 5.0 | ug/L | N | | 1,1-Dichloroethene | < 0.7 | ug/L | A | |
| Acetone | < 10.0 | ug/L | A | | Carbon disulfide | < 5.0 | ug/L | A | |
| Methylene chloride | < 5.0 | ug/L | A | | t-Butanol | < 20.0 | ug/L | N | QA- |
| Methyl-t-butyl ether (MTBE) | < 2.0 | ug/L | A | | trans-1,2-Dichloroethene | < 1.0 | ug/L | A | |
| Di-isopropyl ether (DIPE) | < 2.0 | ug/L | N | | 1,1-Dichloroethane | < 1.0 | ug/L | A | |
| Ethyl-t-butyl ether (ETBE) | < 2.0 | ug/L | N | | 2-Butanone | < 10.0 | ug/L | A | |
| 2,2-Dichloropropane | < 1.0 | ug/L | N | | cis-1,2-Dichloroethene | < 1.0 | ug/L | A | |
| Bromochloromethane | < 0.8 | ug/L | A | | Chloroform | < 1.0 | ug/L | A | |
| Tetrahydrofuran | < 10.0 | ug/L | N | | 1,1,1-Trichloroethane | < 1.0 | ug/L | A | |
| Carbon tetrachloride | < 0.5 | ug/L | A | | 1,1-Dichloropropene | < 1.0 | ug/L | N | |
| Benzene | < 0.5 | ug/L | A | | t-Amylmethyl ether (TAME) | < 2.0 | ug/L | N | |
| 1,2-Dichloroethane | < 0.5 | ug/L | A | | Trichloroethene | < 0.5 | ug/L | A | |
| 1,2-Dichloropropane | < 0.5 | ug/L | A | | Dibromomethane | < 2.0 | ug/L | A | |
| Bromodichloromethane | < 0.5 | ug/L | A | | cis-1,3-Dichloropropene | < 1.0 | ug/L | A | |
| 4-Methyl-2-pentanone (MIBK) | < 10.0 | ug/L | A | | Toluene | < 1.0 | ug/L | A | |
| trans-1,3-Dichloropropene | < 1.0 | ug/L | A | | 1,1,2-Trichloroethane | < 1.0 | ug/L | A | |
| Tetrachloroethene | < 0.5 | ug/L | A | | 1,3-Dichloropropane | < 1.0 | ug/L | N | |
| 2-Hexanone | < 10.0 | ug/L | A | | Dibromochloromethane | < 1.0 | ug/L | A | |
| 1,2-Dibromoethane | < 2.0 | ug/L | A | | Chlorobenzene | < 1.0 | ug/L | A | |
| Ethylbenzene | < 1.0 | ug/L | A | | 1,1,1,2-Tetrachloroethane | < 2.0 | ug/L | A | |
| Xylenes, Total | < 2.0 | ug/L | A | | Styrene | < 1.0 | ug/L | A | |
| Bromoform | < 2.0 | ug/L | A | | Isopropylbenzene | < 1.0 | ug/L | A | |
| 1,1,2,2-Tetrachloroethane | < 2.0 | ug/L | A | | Bromobenzene | < 1.0 | ug/L | A | |
| n-Propylbenzene | < 1.0 | ug/L | A | | 1,2,3-Trichloropropane | < 2.0 | ug/L | A | |
| 2-Chlorotoluene | < 1.0 | ug/L | A | | 1,3,5-Trimethylbenzene | < 1.0 | ug/L | A | |
| 4-Chlorotoluene | < 1.0 | ug/L | A | | t-Butylbenzene | < 1.0 | ug/L | A | |
| 1,2,4-Trimethylbenzene | < 1.0 | ug/L | A | | s-Butylbenzene | < 1.0 | ug/L | A | |
| 4-Isopropyltoluene | < 1.0 | ug/L | A | | 1,3-Dichlorobenzene | < 1.0 | ug/L | A | |
| 1,4-Dichlorobenzene | < 1.0 | ug/L | A | | 1,2,3-Trimethylbenzene | < 1.0 | ug/L | U | |
| n-Butylbenzene | < 1.0 | ug/L | A | | 1,2-Dichlorobenzene | < 1.0 | ug/L | A | |
| 1,2-Dibromo-3-Chloropropane | < 2.0 | ug/L | A | | 1,2,4-Trichlorobenzene | < 2.0 | ug/L | A | |
| 1,3,5-Trichlorobenzene | < 2.0 | ug/L | N | | Hexachlorobutadiene | < 0.5 | ug/L | A | |
| Naphthalene | < 0.5 | ug/L | A | | 1,2,3-Trichlorobenzene | < 0.5 | ug/L | A | |
| Surr. 1 (Dibromofluoromethane) | 100 | % | A | | Surr. 2 (Toluene d8) | 100 | % | A | |
| Surr. 3 (4-Bromofluorobenzene) | 99 | % | A | | Unidentified Peaks | 0 | | U | |

Laboratory Report

REPORT DATE: 6/26/2023

CLIENT: KAS, Inc
PROJECT: Brandon Landfill

WORK ORDER: 2305-14321
DATE RECEIVED: 05/26/2023

TEST METHOD: EPA 8260C

| 005 Site: Duplicate | | Sampled: 5/22/23 | | 10:35 | | Test Date: 6/3/23 | | W TRP | |
|--------------------------------|--------|------------------|-------|-------|---------------------------|-------------------|------|-------|------|
| Parameter | Result | Unit | Nelac | Qual | Parameter | Result | Unit | Nelac | Qual |
| Dichlorodifluoromethane | < 5.0 | ug/L | A | | Chloromethane | < 3.0 | ug/L | A | |
| Vinyl chloride | < 0.5 | ug/L | A | | Bromomethane | < 0.5 | ug/L | A | |
| Chloroethane | < 5.0 | ug/L | A | | Trichlorofluoromethane | < 2.0 | ug/L | A | |
| Diethyl ether | < 5.0 | ug/L | N | | 1,1-Dichloroethene | < 0.7 | ug/L | A | |
| Acetone | < 10.0 | ug/L | A | | Carbon disulfide | < 5.0 | ug/L | A | |
| Methylene chloride | < 5.0 | ug/L | A | | t-Butanol | < 20.0 | ug/L | N | QA- |
| Methyl-t-butyl ether (MTBE) | < 2.0 | ug/L | A | | trans-1,2-Dichloroethene | < 1.0 | ug/L | A | |
| Di-isopropyl ether (DIPE) | < 2.0 | ug/L | N | | 1,1-Dichloroethane | < 1.0 | ug/L | A | |
| Ethyl-t-butyl ether (ETBE) | < 2.0 | ug/L | N | | 2-Butanone | < 10.0 | ug/L | A | |
| 2,2-Dichloropropane | < 1.0 | ug/L | N | | cis-1,2-Dichloroethene | < 1.0 | ug/L | A | |
| Bromochloromethane | < 0.8 | ug/L | A | | Chloroform | < 1.0 | ug/L | A | |
| Tetrahydrofuran | < 10.0 | ug/L | N | | 1,1,1-Trichloroethane | < 1.0 | ug/L | A | |
| Carbon tetrachloride | < 0.5 | ug/L | A | | 1,1-Dichloropropene | < 1.0 | ug/L | N | |
| Benzene | < 0.5 | ug/L | A | | t-Amylmethyl ether (TAME) | < 2.0 | ug/L | N | |
| 1,2-Dichloroethane | < 0.5 | ug/L | A | | Trichloroethene | < 0.5 | ug/L | A | |
| 1,2-Dichloropropane | < 0.5 | ug/L | A | | Dibromomethane | < 2.0 | ug/L | A | |
| Bromodichloromethane | < 0.5 | ug/L | A | | cis-1,3-Dichloropropene | < 1.0 | ug/L | A | |
| 4-Methyl-2-pentanone (MIBK) | < 10.0 | ug/L | A | | Toluene | < 1.0 | ug/L | A | |
| trans-1,3-Dichloropropene | < 1.0 | ug/L | A | | 1,1,2-Trichloroethane | < 1.0 | ug/L | A | |
| Tetrachloroethene | < 0.5 | ug/L | A | | 1,3-Dichloropropane | < 1.0 | ug/L | N | |
| 2-Hexanone | < 10.0 | ug/L | A | | Dibromochloromethane | < 1.0 | ug/L | A | |
| 1,2-Dibromoethane | < 2.0 | ug/L | A | | Chlorobenzene | < 1.0 | ug/L | A | |
| Ethylbenzene | < 1.0 | ug/L | A | | 1,1,1,2-Tetrachloroethane | < 2.0 | ug/L | A | |
| Xylenes, Total | < 2.0 | ug/L | A | | Styrene | < 1.0 | ug/L | A | |
| Bromoform | < 2.0 | ug/L | A | | Isopropylbenzene | < 1.0 | ug/L | A | |
| 1,1,2,2-Tetrachloroethane | < 2.0 | ug/L | A | | Bromobenzene | < 1.0 | ug/L | A | |
| n-Propylbenzene | < 1.0 | ug/L | A | | 1,2,3-Trichloropropane | < 2.0 | ug/L | A | |
| 2-Chlorotoluene | < 1.0 | ug/L | A | | 1,3,5-Trimethylbenzene | < 1.0 | ug/L | A | |
| 4-Chlorotoluene | < 1.0 | ug/L | A | | t-Butylbenzene | < 1.0 | ug/L | A | |
| 1,2,4-Trimethylbenzene | < 1.0 | ug/L | A | | s-Butylbenzene | < 1.0 | ug/L | A | |
| 4-Isopropyltoluene | < 1.0 | ug/L | A | | 1,3-Dichlorobenzene | < 1.0 | ug/L | A | |
| 1,4-Dichlorobenzene | < 1.0 | ug/L | A | | 1,2,3-Trimethylbenzene | < 1.0 | ug/L | U | |
| n-Butylbenzene | < 1.0 | ug/L | A | | 1,2-Dichlorobenzene | < 1.0 | ug/L | A | |
| 1,2-Dibromo-3-Chloropropane | < 2.0 | ug/L | A | | 1,2,4-Trichlorobenzene | < 2.0 | ug/L | A | |
| 1,3,5-Trichlorobenzene | < 2.0 | ug/L | N | | Hexachlorobutadiene | < 0.5 | ug/L | A | |
| Naphthalene | < 0.5 | ug/L | A | | 1,2,3-Trichlorobenzene | < 0.5 | ug/L | A | |
| Surr. 1 (Dibromofluoromethane) | 100 | % | A | | Surr. 2 (Toluene d8) | 100 | % | A | |
| Surr. 3 (4-Bromofluorobenzene) | 99 | % | A | | Unidentified Peaks | 0 | | U | |

CLIENT: KAS, Inc
PROJECT: Brandon Landfill

WORK ORDER: 2305-14321
DATE RECEIVED: 05/26/2023

TEST METHOD: EPA 8260C

| 006 | Site: Trip Blank | Sampled: 5/22/23 11:05 | | Test Date: 6/3/23 | | W | TRP | | |
|--------------------------------|------------------|------------------------|-------|-------------------|---------------------------|--------|------|-------|------|
| Parameter | Result | Unit | Nelac | Qual | Parameter | Result | Unit | Nelac | Qual |
| Dichlorodifluoromethane | < 5.0 | ug/L | A | | Chloromethane | < 3.0 | ug/L | A | |
| Vinyl chloride | < 0.5 | ug/L | A | | Bromomethane | < 0.5 | ug/L | A | |
| Chloroethane | < 5.0 | ug/L | A | | Trichlorofluoromethane | < 2.0 | ug/L | A | |
| Diethyl ether | < 5.0 | ug/L | N | | 1,1-Dichloroethene | < 0.7 | ug/L | A | |
| Acetone | < 10.0 | ug/L | A | | Carbon disulfide | < 5.0 | ug/L | A | |
| Methylene chloride | < 5.0 | ug/L | A | | t-Butanol | < 20.0 | ug/L | N | QA- |
| Methyl-t-butyl ether (MTBE) | < 2.0 | ug/L | A | | trans-1,2-Dichloroethene | < 1.0 | ug/L | A | |
| Di-isopropyl ether (DIPE) | < 2.0 | ug/L | N | | 1,1-Dichloroethane | < 1.0 | ug/L | A | |
| Ethyl-t-butyl ether (ETBE) | < 2.0 | ug/L | N | | 2-Butanone | < 10.0 | ug/L | A | |
| 2,2-Dichloropropane | < 1.0 | ug/L | N | | cis-1,2-Dichloroethene | < 1.0 | ug/L | A | |
| Bromochloromethane | < 0.8 | ug/L | A | | Chloroform | < 1.0 | ug/L | A | |
| Tetrahydrofuran | < 10.0 | ug/L | N | | 1,1,1-Trichloroethane | < 1.0 | ug/L | A | |
| Carbon tetrachloride | < 0.5 | ug/L | A | | 1,1-Dichloropropene | < 1.0 | ug/L | N | |
| Benzene | < 0.5 | ug/L | A | | t-Amylmethyl ether (TAME) | < 2.0 | ug/L | N | |
| 1,2-Dichloroethane | < 0.5 | ug/L | A | | Trichloroethene | < 0.5 | ug/L | A | |
| 1,2-Dichloropropane | < 0.5 | ug/L | A | | Dibromomethane | < 2.0 | ug/L | A | |
| Bromodichloromethane | < 0.5 | ug/L | A | | cis-1,3-Dichloropropene | < 1.0 | ug/L | A | |
| 4-Methyl-2-pentanone (MIBK) | < 10.0 | ug/L | A | | Toluene | < 1.0 | ug/L | A | |
| trans-1,3-Dichloropropene | < 1.0 | ug/L | A | | 1,1,2-Trichloroethane | < 1.0 | ug/L | A | |
| Tetrachloroethene | < 0.5 | ug/L | A | | 1,3-Dichloropropane | < 1.0 | ug/L | N | |
| 2-Hexanone | < 10.0 | ug/L | A | | Dibromochloromethane | < 1.0 | ug/L | A | |
| 1,2-Dibromoethane | < 2.0 | ug/L | A | | Chlorobenzene | < 1.0 | ug/L | A | |
| Ethylbenzene | < 1.0 | ug/L | A | | 1,1,1,2-Tetrachloroethane | < 2.0 | ug/L | A | |
| Xylenes, Total | < 2.0 | ug/L | A | | Styrene | < 1.0 | ug/L | A | |
| Bromoform | < 2.0 | ug/L | A | | Isopropylbenzene | < 1.0 | ug/L | A | |
| 1,1,2,2-Tetrachloroethane | < 2.0 | ug/L | A | | Bromobenzene | < 1.0 | ug/L | A | |
| n-Propylbenzene | < 1.0 | ug/L | A | | 1,2,3-Trichloropropane | < 2.0 | ug/L | A | |
| 2-Chlorotoluene | < 1.0 | ug/L | A | | 1,3,5-Trimethylbenzene | < 1.0 | ug/L | A | |
| 4-Chlorotoluene | < 1.0 | ug/L | A | | t-Butylbenzene | < 1.0 | ug/L | A | |
| 1,2,4-Trimethylbenzene | < 1.0 | ug/L | A | | s-Butylbenzene | < 1.0 | ug/L | A | |
| 4-Isopropyltoluene | < 1.0 | ug/L | A | | 1,3-Dichlorobenzene | < 1.0 | ug/L | A | |
| 1,4-Dichlorobenzene | < 1.0 | ug/L | A | | 1,2,3-Trimethylbenzene | < 1.0 | ug/L | U | |
| n-Butylbenzene | < 1.0 | ug/L | A | | 1,2-Dichlorobenzene | < 1.0 | ug/L | A | |
| 1,2-Dibromo-3-Chloropropane | < 2.0 | ug/L | A | | 1,2,4-Trichlorobenzene | < 2.0 | ug/L | A | |
| 1,3,5-Trichlorobenzene | < 2.0 | ug/L | N | | Hexachlorobutadiene | < 0.5 | ug/L | A | |
| Naphthalene | < 0.5 | ug/L | A | | 1,2,3-Trichlorobenzene | < 0.5 | ug/L | A | |
| Surr. 1 (Dibromofluoromethane) | 99 | % | A | | Surr. 2 (Toluene d8) | 99 | % | A | |
| Surr. 3 (4-Bromofluorobenzene) | 99 | % | A | | Unidentified Peaks | 0 | | U | |

Report Summary of Qualifiers and Notes

QA-: QA/QC associated with this analysis did not meet laboratory acceptance limits indicating the results may be biased low.

B: Blank contamination was observed at levels that could affect analytical results.

June 23, 2023

Clare Santos
KAS Environmental
589 Avenue D
Williston, VT 05495

Project Location: Brandon, VT
Client Job Number:
Project Number: 609210052
Laboratory Work Order Number: 23F0005

Enclosed are results of analyses for samples as received by the laboratory on May 31, 2023. If you have any questions concerning this report, please feel free to contact me.

Sincerely,



Kaitlyn A. Feliciano
Project Manager

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KAS Environmental
589 Avenue D
Williston, VT 05495
ATTN: Clare Santos

REPORT DATE: 6/23/2023

PURCHASE ORDER NUMBER:

PROJECT NUMBER: 609210052

ANALYTICAL SUMMARY

WORK ORDER NUMBER: 23F0005

The results of analyses performed on the following samples submitted to CON-TEST, a Pace Analytical Laboratory, are found in this report.

PROJECT LOCATION: Brandon, VT

| FIELD SAMPLE # | LAB ID: | MATRIX | SAMPLE DESCRIPTION | TEST | SUB LAB |
|----------------|------------|-----------------------|--------------------|--------------|---------|
| MW-1 | 23F0005-01 | Ground Water | | SOP-454 PFAS | |
| MW-2C | 23F0005-02 | Ground Water | | SOP-454 PFAS | |
| MW-3 | 23F0005-03 | Ground Water | | SOP-454 PFAS | |
| MW-5 | 23F0005-04 | Ground Water | | SOP-454 PFAS | |
| Brandon ERB | 23F0005-05 | Equipment Blank Water | | SOP-454 PFAS | |

CASE NARRATIVE SUMMARY

All reported results are within defined laboratory quality control objectives unless listed below or otherwise qualified in this report.

SOP-454 PFAS**Qualifications:**

PF-17

Extracted Internal Standard recovery is outside of control limits. Data is not significantly affected since associated analyte is not detected and bias is on the high side.

Analyte & Samples(s) Qualified:**M2-4:2FTS**

23F0005-01[MW-1], 23F0005-02[MW-2C], 23F0005-04[MW-5]

M2-6:2FTS

23F0005-01[MW-1], 23F0005-02[MW-2C], 23F0005-04[MW-5]

M2-8:2FTS

23F0005-01[MW-1], 23F0005-02[MW-2C]

PF-18

Duplicate analysis confirmed Extracted Internal Standard failure due to matrix effects.

Analyte & Samples(s) Qualified:**D3-NMeFOSAA**

23F0005-03[MW-3]

D5-NEtFOSAA

23F0005-03[MW-3], 23F0005-04[MW-5]

M2PF₇A

23F0005-02[MW-2C], 23F0005-03[MW-3], 23F0005-04[MW-5]

M7PF₁₀UnA

23F0005-03[MW-3], 23F0005-04[MW-5]

MPFBA

23F0005-02[MW-2C]

MPFDoA

23F0005-02[MW-2C], 23F0005-03[MW-3], 23F0005-04[MW-5]

PF-20

Quantifying ion signal to noise ratio is <10. Detection is suspect.

Analyte & Samples(s) Qualified:**Perfluorobutanoic acid (PFBA)**

23F0005-02[MW-2C], 23F0005-04[MW-5]

Perfluoropentanoic acid (PFPeA)

23F0005-02[MW-2C]

S-29

Extracted Internal Standard is outside of control limits.

Analyte & Samples(s) Qualified:**M2PF₇A**

23F0005-01[MW-1]

The results of analyses reported only relate to samples submitted to Con-Test, a Pace Analytical Laboratory, for testing.

I certify that the analyses listed above, unless specifically listed as subcontracted, if any, were performed under my direction according to the approved methodologies listed in this document, and that based upon my inquiry of those individuals immediately responsible for obtaining the information, the material contained in this report is, to the best of my knowledge and belief, accurate and complete.



Lisa A. Worthington
Technical Representative

Project Location: Brandon, VT

Sample Description:

Work Order: 23F0005

Date Received: 5/31/2023

Field Sample #: MW-1

Sampled: 5/22/2023 12:25

Sample ID: 23F0005-01

Sample Matrix: Ground Water

Semivolatile Organic Compounds by - LC/MS-MS

| Analyte | Results | RL | Units | Dilution | Flag/Qual | Method | Date Prepared | Date/Time Analyzed | Analyst |
|---|---------|-----|-------|----------|-----------|--------------|---------------|--------------------|---------|
| Perfluorobutanoic acid (PFBA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorobutanesulfonic acid (PFBS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluoropentanoic acid (PFPeA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorohexanoic acid (PFHxA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| 11Cl-PF3OUdS (F53B Major) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| 9Cl-PF3ONS (F53B Minor) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Hexafluoropropylene oxide dimer acid (HFPO-DA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| 8:2 Fluorotelomersulfonic acid (8:2FTS A) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorodecanoic acid (PFDA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorododecanoic acid (PFDoA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluoroheptanesulfonic acid (PFHpS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| N-EtFOSAA (NEtFOSAA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| N-MeFOSAA (NMeFOSAA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorotetradecanoic acid (PFTA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorotridecanoic acid (PFTrDA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| 4:2 Fluorotelomersulfonic acid (4:2FTS A) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorodecanesulfonic acid (PFDS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorooctanesulfonamide (FOSA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorononanesulfonic acid (PFNS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluoro-1-hexanesulfonamide (FHxSA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluoro-1-butanesulfonamide (FBSA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorohexanesulfonic acid (PFHxS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluoro-4-oxapentanoic acid (PFMPA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluoro-5-oxahexanoic acid (PFMBA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| 6:2 Fluorotelomersulfonic acid (6:2FTS A) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluoropentanesulfonic acid (PFPeS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluoroundecanoic acid (PFUnA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Nonfluoro-3,6-dioxaheptanoic acid (NFDHA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluoroheptanoic acid (PFHpA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorooctanoic acid (PFOA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorooctanesulfonic acid (PFOS) | 2.4 | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |
| Perfluorononanoic acid (PFNA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:41 | RRB |

Project Location: Brandon, VT

Sample Description:

Work Order: 23F0005

Date Received: 5/31/2023

Field Sample #: MW-2C

Sampled: 5/22/2023 09:30

Sample ID: 23F0005-02

Sample Matrix: Ground Water

Semivolatile Organic Compounds by - LC/MS-MS

| Analyte | Results | RL | Units | Dilution | Flag/Qual | Method | Date Prepared | Date/Time Analyzed | Analyst |
|---|---------|-----|-------|----------|-----------|--------------|---------------|--------------------|---------|
| Perfluorobutanoic acid (PFBA) | 4.6 | 1.8 | ng/L | 1 | PF-20 | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorobutanesulfonic acid (PFBS) | 5.3 | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluoropentanoic acid (PFPeA) | 20 | 1.8 | ng/L | 1 | PF-20 | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorohexanoic acid (PFHxA) | 25 | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| 11Cl-PF3OUdS (F53B Major) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| 9Cl-PF3ONS (F53B Minor) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Hexafluoropropylene oxide dimer acid (HFPO-DA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| 8:2 Fluorotelomersulfonic acid (8:2FTS A) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorodecanoic acid (PFDA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorododecanoic acid (PFDoA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluoroheptanesulfonic acid (PFHpS) | 5.0 | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| N-EtFOSAA (NEtFOSAA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| N-MeFOSAA (NMeFOSAA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorotetradecanoic acid (PFTA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorotridecanoic acid (PFTrDA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| 4:2 Fluorotelomersulfonic acid (4:2FTS A) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorodecanesulfonic acid (PFDS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorooctanesulfonamide (FOSA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorononanesulfonic acid (PFNS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluoro-1-hexanesulfonamide (FHxSA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluoro-1-butanesulfonamide (FBSA) | 4.5 | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorohexanesulfonic acid (PFHxS) | 36 | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluoro-4-oxapentanoic acid (PFMPA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluoro-5-oxahexanoic acid (PFMBA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| 6:2 Fluorotelomersulfonic acid (6:2FTS A) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluoropentanesulfonic acid (PFPeS) | 4.3 | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluoroundecanoic acid (PFUnA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluoroheptanoic acid (PFHpA) | 18 | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorooctanoic acid (PFOA) | 97 | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorooctanesulfonic acid (PFOS) | 150 | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |
| Perfluorononanoic acid (PFNA) | 4.7 | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:48 | RRB |

Project Location: Brandon, VT

Sample Description:

Work Order: 23F0005

Date Received: 5/31/2023

Field Sample #: MW-3

Sampled: 5/22/2023 11:46

Sample ID: 23F0005-03

Sample Matrix: Ground Water

Semivolatile Organic Compounds by - LC/MS-MS

| Analyte | Results | RL | Units | Dilution | Flag/Qual | Method | Date Prepared | Date/Time Analyzed | Analyst |
|---|---------|-----|-------|----------|-----------|--------------|---------------|--------------------|---------|
| Perfluorobutanoic acid (PFBA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorobutanesulfonic acid (PFBS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluoropentanoic acid (PFPeA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorohexanoic acid (PFHxA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| 11Cl-PF3OUdS (F53B Major) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| 9Cl-PF3ONS (F53B Minor) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Hexafluoropropylene oxide dimer acid (HFPO-DA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| 8:2 Fluorotelomersulfonic acid (8:2FTS A) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorodecanoic acid (PFDA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorododecanoic acid (PFDoA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluoroheptanesulfonic acid (PFHpS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| N-EtFOSAA (NEtFOSAA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| N-MeFOSAA (NMeFOSAA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorotetradecanoic acid (PFTA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorotridecanoic acid (PFTrDA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| 4:2 Fluorotelomersulfonic acid (4:2FTS A) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorodecanesulfonic acid (PFDS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorooctanesulfonamide (FOSA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorononanesulfonic acid (PFNS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluoro-1-hexanesulfonamide (FHxSA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluoro-1-butanesulfonamide (FBSA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorohexanesulfonic acid (PFHxS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluoro-4-oxapentanoic acid (PFMPA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluoro-5-oxahexanoic acid (PFMBA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| 6:2 Fluorotelomersulfonic acid (6:2FTS A) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluoropentanesulfonic acid (PFPeS) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluoroundecanoic acid (PFUnA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Nonfluoro-3,6-dioxaheptanoic acid (NFDHA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluoroheptanoic acid (PFHpA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorooctanoic acid (PFOA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorooctanesulfonic acid (PFOS) | 5.2 | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |
| Perfluorononanoic acid (PFNA) | ND | 1.8 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 18:56 | RRB |

Project Location: Brandon, VT

Sample Description:

Work Order: 23F0005

Date Received: 5/31/2023

Field Sample #: MW-5

Sampled: 5/22/2023 10:35

Sample ID: 23F0005-04

Sample Matrix: Ground Water

Semivolatile Organic Compounds by - LC/MS-MS

| Analyte | Results | RL | Units | Dilution | Flag/Qual | Method | Date Prepared | Date/Time Analyzed | Analyst |
|---|---------|-----|-------|----------|-----------|--------------|---------------|--------------------|---------|
| Perfluorobutanoic acid (PFBA) | 3.1 | 1.9 | ng/L | 1 | PF-20 | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorobutanesulfonic acid (PFBS) | 2.1 | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluoropentanoic acid (PFPeA) | 3.1 | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorohexanoic acid (PFHxA) | 6.5 | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| 11Cl-PF3OUdS (F53B Major) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| 9Cl-PF3ONS (F53B Minor) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Hexafluoropropylene oxide dimer acid (HFPO-DA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| 8:2 Fluorotelomersulfonic acid (8:2FTS A) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorodecanoic acid (PFDA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorododecanoic acid (PFDoA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluoroheptanesulfonic acid (PFHpS) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| N-EtFOSAA (NEtFOSAA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| N-MeFOSAA (NMeFOSAA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorotetradecanoic acid (PFTA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorotridecanoic acid (PFTTrDA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| 4:2 Fluorotelomersulfonic acid (4:2FTS A) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorodecanesulfonic acid (PFDS) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorooctanesulfonamide (FOSA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorononanesulfonic acid (PFNS) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluoro-1-hexanesulfonamide (FHxSA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluoro-1-butanesulfonamide (FBSA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorohexanesulfonic acid (PFHxS) | 17 | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluoro-4-oxapentanoic acid (PFMPA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluoro-5-oxahexanoic acid (PFMBA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| 6:2 Fluorotelomersulfonic acid (6:2FTS A) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluoropentanesulfonic acid (PFPeS) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluoroundecanoic acid (PFUnA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluoroheptanoic acid (PFHpA) | 5.1 | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorooctanoic acid (PFOA) | 30 | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorooctanesulfonic acid (PFOS) | 14 | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |
| Perfluorononanoic acid (PFNA) | ND | 1.9 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:03 | RRB |

Project Location: Brandon, VT

Sample Description:

Work Order: 23F0005

Date Received: 5/31/2023

 Field Sample #: **Brandon ERB**

Sampled: 5/22/2023 12:18

 Sample ID: **23F0005-05**

Sample Matrix: Equipment Blank Water

Semivolatile Organic Compounds by - LC/MS-MS

| Analyte | Results | RL | Units | Dilution | Flag/Qual | Method | Date Prepared | Date/Time Analyzed | Analyst |
|---|---------|-----|-------|----------|-----------|--------------|---------------|--------------------|---------|
| Perfluorobutanoic acid (PFBA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorobutanesulfonic acid (PFBS) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluoropentanoic acid (PFPeA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorohexanoic acid (PFHxA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| 11Cl-PF3OUdS (F53B Major) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| 9Cl-PF3ONS (F53B Minor) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Hexafluoropropylene oxide dimer acid (HFPO-DA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| 8:2 Fluorotelomersulfonic acid (8:2FTS A) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorodecanoic acid (PFDA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorododecanoic acid (PFDoA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluoroheptanesulfonic acid (PFHpS) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| N-EtFOSAA (NEtFOSAA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| N-MeFOSAA (NMeFOSAA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorotetradecanoic acid (PFTA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorotridecanoic acid (PFTTrDA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| 4:2 Fluorotelomersulfonic acid (4:2FTS A) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorodecanesulfonic acid (PFDS) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorooctanesulfonamide (FOSA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorononanesulfonic acid (PFNS) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluoro-1-hexanesulfonamide (FHxSA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluoro-1-butanefulfonamide (FBSA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorohexanesulfonic acid (PFHxS) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluoro-4-oxapentanoic acid (PFMPA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluoro-5-oxahexanoic acid (PFMBA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| 6:2 Fluorotelomersulfonic acid (6:2FTS A) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluoropentanesulfonic acid (PFPeS) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluoroundecanoic acid (PFUnA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Nonfluoro-3,6-dioxaheptanoic acid (NFDHA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluoroheptanoic acid (PFHpA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorooctanoic acid (PFOA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorooctanesulfonic acid (PFOS) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |
| Perfluorononanoic acid (PFNA) | ND | 2.1 | ng/L | 1 | | SOP-454 PFAS | 6/14/23 | 6/15/23 19:10 | RRB |

Sample Extraction Data

Prep Method:SOP 454-PFAAS Analytical Method:SOP-454 PFAS

| Lab Number [Field ID] | Batch | Initial [mL] | Final [mL] | Date |
|--------------------------|---------|--------------|------------|----------|
| 23F0005-01 [MW-1] | B343017 | 271 | 1.00 | 06/14/23 |
| 23F0005-02 [MW-2C] | B343017 | 272 | 1.00 | 06/14/23 |
| 23F0005-03 [MW-3] | B343017 | 275 | 1.00 | 06/14/23 |
| 23F0005-04 [MW-5] | B343017 | 269 | 1.00 | 06/14/23 |
| 23F0005-05 [Brandon ERB] | B343017 | 233 | 1.00 | 06/14/23 |

QUALITY CONTROL
Semivolatile Organic Compounds by - LC/MS-MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|

Batch B343017 - SOP 454-PFAAS
Blank (B343017-BLK1)

Prepared: 06/14/23 Analyzed: 06/15/23

| | | | | | | | | | | |
|--|----|-----|------|--|--|--|--|--|--|--|
| Perfluorobutanoic acid (PFBA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorobutanesulfonic acid (PFBS) | ND | 2.0 | ng/L | | | | | | | |
| Perfluoropentanoic acid (PFPeA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorohexanoic acid (PFHxA) | ND | 2.0 | ng/L | | | | | | | |
| 11Cl-PF3OUdS (F53B Major) | ND | 2.0 | ng/L | | | | | | | |
| 9Cl-PF3ONS (F53B Minor) | ND | 2.0 | ng/L | | | | | | | |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA) | ND | 2.0 | ng/L | | | | | | | |
| Hexafluoropropylene oxide dimer acid (HFPO-DA) | ND | 2.0 | ng/L | | | | | | | |
| 8:2 Fluorotelomersulfonic acid (8:2FTS A) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorodecanoic acid (PFDA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorododecanoic acid (PFDoA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluoro(2-ethoxyethane)sulfonic acid (PFEEESA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluoroheptanesulfonic acid (PFHpS) | ND | 2.0 | ng/L | | | | | | | |
| N-EtFOSAA (NEtFOSAA) | ND | 2.0 | ng/L | | | | | | | |
| N-MeFOSAA (NMeFOSAA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorotetradecanoic acid (PFTA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorotridecanoic acid (PFTrDA) | ND | 2.0 | ng/L | | | | | | | |
| 4:2 Fluorotelomersulfonic acid (4:2FTS A) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorodecanesulfonic acid (PFDS) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorooctanesulfonamide (FOSA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorononanesulfonic acid (PFNS) | ND | 2.0 | ng/L | | | | | | | |
| Perfluoro-1-hexanesulfonamide (FHxSA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluoro-1-butanesulfonamide (FBSA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorohexanesulfonic acid (PFHxS) | ND | 2.0 | ng/L | | | | | | | |
| Perfluoro-4-oxapentanoic acid (PFMPA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluoro-5-oxahexanoic acid (PFMBA) | ND | 2.0 | ng/L | | | | | | | |
| 6:2 Fluorotelomersulfonic acid (6:2FTS A) | ND | 2.0 | ng/L | | | | | | | |
| Perfluoropentanesulfonic acid (PFPeS) | ND | 2.0 | ng/L | | | | | | | |
| Perfluoroundecanoic acid (PFUnA) | ND | 2.0 | ng/L | | | | | | | |
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluoroheptanoic acid (PFHpA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorooctanoic acid (PFOA) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorooctanesulfonic acid (PFOS) | ND | 2.0 | ng/L | | | | | | | |
| Perfluorononanoic acid (PFNA) | ND | 2.0 | ng/L | | | | | | | |

LCS (B343017-BS1)

Prepared: 06/14/23 Analyzed: 06/15/23

| | | | | | | |
|--|------|-----|------|------|------|----------|
| Perfluorobutanoic acid (PFBA) | 8.23 | 2.0 | ng/L | 9.98 | 82.5 | 73-129 |
| Perfluorobutanesulfonic acid (PFBS) | 7.28 | 2.0 | ng/L | 8.83 | 82.4 | 72-130 |
| Perfluoropentanoic acid (PFPeA) | 8.11 | 2.0 | ng/L | 9.98 | 81.2 | 72-129 |
| Perfluorohexanoic acid (PFHxA) | 8.31 | 2.0 | ng/L | 9.98 | 83.3 | 72-129 |
| 11Cl-PF3OUdS (F53B Major) | 7.66 | 2.0 | ng/L | 9.40 | 81.5 | 55.1-141 |
| 9Cl-PF3ONS (F53B Minor) | 8.12 | 2.0 | ng/L | 9.30 | 87.3 | 59.6-146 |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA) | 8.23 | 2.0 | ng/L | 9.40 | 87.5 | 60.3-131 |
| Hexafluoropropylene oxide dimer acid (HFPO-DA) | 8.53 | 2.0 | ng/L | 9.98 | 85.5 | 37.6-167 |
| 8:2 Fluorotelomersulfonic acid (8:2FTS A) | 7.76 | 2.0 | ng/L | 9.58 | 81.0 | 67-138 |
| Perfluorodecanoic acid (PFDA) | 8.30 | 2.0 | ng/L | 9.98 | 83.2 | 71-129 |
| Perfluorododecanoic acid (PFDoA) | 8.77 | 2.0 | ng/L | 9.98 | 87.8 | 72-134 |
| Perfluoro(2-ethoxyethane)sulfonic acid (PFEEESA) | 7.71 | 2.0 | ng/L | 8.88 | 86.8 | 49.4-154 |

QUALITY CONTROL
Semivolatile Organic Compounds by - LC/MS-MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|

Batch B343017 - SOP 454-PFAAS
LCS (B343017-BS1)

Prepared: 06/14/23 Analyzed: 06/15/23

| | | | | | | | | | | |
|---|------|-----|------|------|--|------|----------|--|--|--|
| Perfluoroheptanesulfonic acid (PFHpS) | 9.00 | 2.0 | ng/L | 9.53 | | 94.5 | 69-134 | | | |
| N-EtFOSAA (NEtFOSAA) | 9.15 | 2.0 | ng/L | 9.98 | | 91.7 | 61-135 | | | |
| N-MeFOSAA (NMeFOSAA) | 10.3 | 2.0 | ng/L | 9.98 | | 103 | 65-136 | | | |
| Perfluorotetradecanoic acid (PFTA) | 8.49 | 2.0 | ng/L | 9.98 | | 85.0 | 71-132 | | | |
| Perfluorotridecanoic acid (PFTTrDA) | 8.88 | 2.0 | ng/L | 9.98 | | 89.0 | 65-144 | | | |
| 4:2 Fluorotelomersulfonic acid (4:2FTS A) | 8.31 | 2.0 | ng/L | 9.33 | | 89.1 | 63-143 | | | |
| Perfluorodecanesulfonic acid (PFDS) | 6.55 | 2.0 | ng/L | 9.63 | | 68.0 | 53-142 | | | |
| Perfluorooctanesulfonamide (FOSA) | 8.35 | 2.0 | ng/L | 9.98 | | 83.7 | 67-137 | | | |
| Perfluorononanesulfonic acid (PFNS) | 8.22 | 2.0 | ng/L | 9.58 | | 85.8 | 69-127 | | | |
| Perfluoro-1-hexanesulfonamide (FHxSA) | 8.19 | 2.0 | ng/L | 9.98 | | 82.0 | 61.7-156 | | | |
| Perfluoro-1-butanefulfonamide (FBSA) | 8.58 | 2.0 | ng/L | 9.98 | | 86.0 | 61.3-145 | | | |
| Perfluorohexanesulfonic acid (PFHxS) | 7.01 | 2.0 | ng/L | 9.13 | | 76.7 | 68-131 | | | |
| Perfluoro-4-oxapentanoic acid (PFMPA) | 9.02 | 2.0 | ng/L | 9.98 | | 90.4 | 59.8-147 | | | |
| Perfluoro-5-oxahexanoic acid (PFMBA) | 8.61 | 2.0 | ng/L | 9.98 | | 86.3 | 59.5-146 | | | |
| 6:2 Fluorotelomersulfonic acid (6:2FTS A) | 8.97 | 2.0 | ng/L | 9.48 | | 94.5 | 64-140 | | | |
| Perfluoropentanesulfonic acid (PFPeS) | 7.51 | 2.0 | ng/L | 9.38 | | 80.1 | 71-127 | | | |
| Perfluoroundecanoic acid (PFUnA) | 8.30 | 2.0 | ng/L | 9.98 | | 83.2 | 69-133 | | | |
| Nonafluoro-3,6-dioxahexanoic acid (NFDHA) | 9.33 | 2.0 | ng/L | 9.98 | | 93.4 | 58.5-143 | | | |
| Perfluoroheptanoic acid (PFHpA) | 8.35 | 2.0 | ng/L | 9.98 | | 83.6 | 72-130 | | | |
| Perfluorooctanoic acid (PFOA) | 8.61 | 2.0 | ng/L | 9.98 | | 86.3 | 71-133 | | | |
| Perfluorooctanesulfonic acid (PFOS) | 7.96 | 2.0 | ng/L | 9.23 | | 86.3 | 65-140 | | | |
| Perfluorononanoic acid (PFNA) | 8.16 | 2.0 | ng/L | 9.98 | | 81.7 | 69-130 | | | |

LCS Dup (B343017-BS1)

Prepared: 06/14/23 Analyzed: 06/15/23

| | | | | | | | | | | |
|---|------|-----|------|------|--|------|----------|-------|----|--|
| Perfluorobutanoic acid (PFBA) | 7.84 | 1.9 | ng/L | 9.52 | | 82.3 | 73-129 | 4.89 | 30 | |
| Perfluorobutanesulfonic acid (PFBS) | 6.93 | 1.9 | ng/L | 8.43 | | 82.2 | 72-130 | 4.90 | 30 | |
| Perfluoropentanoic acid (PFPeA) | 7.73 | 1.9 | ng/L | 9.52 | | 81.1 | 72-129 | 4.83 | 30 | |
| Perfluorohexanoic acid (PFHxA) | 7.76 | 1.9 | ng/L | 9.52 | | 81.5 | 72-129 | 6.81 | 30 | |
| 11Cl-PF3OUdS (F53B Major) | 7.61 | 1.9 | ng/L | 8.97 | | 84.8 | 55.1-141 | 0.668 | 30 | |
| 9Cl-PF3ONS (F53B Minor) | 8.05 | 1.9 | ng/L | 8.88 | | 90.7 | 59.6-146 | 0.873 | 30 | |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA) | 7.95 | 1.9 | ng/L | 8.97 | | 88.6 | 60.3-131 | 3.45 | 30 | |
| Hexafluoropropylene oxide dimer acid (HFPO-DA) | 8.57 | 1.9 | ng/L | 9.52 | | 90.0 | 37.6-167 | 0.481 | 30 | |
| 8:2 Fluorotelomersulfonic acid (8:2FTS A) | 7.85 | 1.9 | ng/L | 9.14 | | 85.9 | 67-138 | 1.09 | 30 | |
| Perfluorodecanoic acid (PFDA) | 7.87 | 1.9 | ng/L | 9.52 | | 82.6 | 71-129 | 5.37 | 30 | |
| Perfluorododecanoic acid (PFDoA) | 7.71 | 1.9 | ng/L | 9.52 | | 80.9 | 72-134 | 12.9 | 30 | |
| Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA) | 7.25 | 1.9 | ng/L | 8.48 | | 85.5 | 49.4-154 | 6.17 | 30 | |
| Perfluoroheptanesulfonic acid (PFHpS) | 8.52 | 1.9 | ng/L | 9.09 | | 93.7 | 69-134 | 5.53 | 30 | |
| N-EtFOSAA (NEtFOSAA) | 8.94 | 1.9 | ng/L | 9.52 | | 93.9 | 61-135 | 2.32 | 30 | |
| N-MeFOSAA (NMeFOSAA) | 9.07 | 1.9 | ng/L | 9.52 | | 95.2 | 65-136 | 12.8 | 30 | |
| Perfluorotetradecanoic acid (PFTA) | 8.23 | 1.9 | ng/L | 9.52 | | 86.4 | 71-132 | 3.08 | 30 | |
| Perfluorotridecanoic acid (PFTTrDA) | 8.15 | 1.9 | ng/L | 9.52 | | 85.6 | 65-144 | 8.59 | 30 | |
| 4:2 Fluorotelomersulfonic acid (4:2FTS A) | 7.69 | 1.9 | ng/L | 8.90 | | 86.3 | 63-143 | 7.80 | 30 | |
| Perfluorodecanesulfonic acid (PFDS) | 5.81 | 1.9 | ng/L | 9.19 | | 63.3 | 53-142 | 12.0 | 30 | |
| Perfluorooctanesulfonamide (FOSA) | 8.54 | 1.9 | ng/L | 9.52 | | 89.7 | 67-137 | 2.18 | 30 | |
| Perfluorononanesulfonic acid (PFNS) | 7.39 | 1.9 | ng/L | 9.14 | | 80.8 | 69-127 | 10.6 | 30 | |
| Perfluoro-1-hexanesulfonamide (FHxSA) | 8.12 | 1.9 | ng/L | 9.52 | | 85.2 | 61.7-156 | 0.841 | 30 | |
| Perfluoro-1-butanefulfonamide (FBSA) | 7.91 | 1.9 | ng/L | 9.52 | | 83.1 | 61.3-145 | 8.13 | 30 | |
| Perfluorohexanesulfonic acid (PFHxS) | 7.03 | 1.9 | ng/L | 8.71 | | 80.7 | 68-131 | 0.313 | 30 | |
| Perfluoro-4-oxapentanoic acid (PFMPA) | 8.73 | 1.9 | ng/L | 9.52 | | 91.7 | 59.8-147 | 3.22 | 30 | |
| Perfluoro-5-oxahexanoic acid (PFMBA) | 8.30 | 1.9 | ng/L | 9.52 | | 87.2 | 59.5-146 | 3.68 | 30 | |

QUALITY CONTROL
Semivolatile Organic Compounds by - LC/MS-MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|

Batch B343017 - SOP 454-PFAAS
LCS Dup (B343017-BSD1)

Prepared: 06/14/23 Analyzed: 06/15/23

| | | | | | | | | | | |
|--|------|-----|------|------|--|------|----------|--------|----|--|
| 6:2 Fluorotelomersulfonic acid (6:2FTS A) | 6.88 | 1.9 | ng/L | 9.05 | | 76.0 | 64-140 | 26.4 | 30 | |
| Perfluoropentanesulfonic acid (PFPeS) | 7.16 | 1.9 | ng/L | 8.95 | | 80.0 | 71-127 | 4.78 | 30 | |
| Perfluoroundecanoic acid (PFUnA) | 7.58 | 1.9 | ng/L | 9.52 | | 79.6 | 69-133 | 9.07 | 30 | |
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA) | 9.17 | 1.9 | ng/L | 9.52 | | 96.3 | 58.5-143 | 1.71 | 30 | |
| Perfluoroheptanoic acid (PFHpA) | 8.07 | 1.9 | ng/L | 9.52 | | 84.7 | 72-130 | 3.42 | 30 | |
| Perfluorooctanoic acid (PFOA) | 8.61 | 1.9 | ng/L | 9.52 | | 90.5 | 71-133 | 0.0320 | 30 | |
| Perfluorooctanesulfonic acid (PFOS) | 7.40 | 1.9 | ng/L | 8.81 | | 84.0 | 65-140 | 7.37 | 30 | |
| Perfluorononanoic acid (PFNA) | 7.74 | 1.9 | ng/L | 9.52 | | 81.3 | 69-130 | 5.27 | 30 | |

FLAG/QUALIFIER SUMMARY

| | |
|-------|--|
| * | QC result is outside of established limits. |
| † | Wide recovery limits established for difficult compound. |
| ‡ | Wide RPD limits established for difficult compound. |
| # | Data exceeded client recommended or regulatory level |
| ND | Not Detected |
| RL | Reporting Limit is at the level of quantitation (LOQ) |
| DL | Detection Limit is the lower limit of detection determined by the MDL study |
| MCL | Maximum Contaminant Level |
| | Percent recoveries and relative percent differences (RPDs) are determined by the software using values in the calculation which have not been rounded. |
| | No results have been blank subtracted unless specified in the case narrative section. |
| PF-17 | Extracted Internal Standard recovery is outside of control limits. Data is not significantly affected since associated analyte is not detected and bias is on the high side. |
| PF-18 | Duplicate analysis confirmed Extracted Internal Standard failure due to matrix effects. |
| PF-20 | Quantifying ion signal to noise ratio is <10. Detection is suspect. |
| S-29 | Extracted Internal Standard is outside of control limits. |

INTERNAL STANDARD AREA AND RT SUMMARY
SOP-454 PFAS

| Internal Standard | Response | RT | Reference Response | Reference RT | Area % | Area % Limits | RT Diff | RT Diff Limit | Q |
|--|----------|----------|--------------------|--------------|--------|---------------|---------|---------------|---|
| MW-1 (23F0005-01) | | | | | | | | | |
| Lab File ID: 23F0005-01.d Analyzed: 06/15/23 18:41 | | | | | | | | | |
| M8FOSA | 162454.2 | 3.980567 | 227,522.00 | 3.980567 | 71 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-4:2FTS | 46432.68 | 2.5543 | 27,028.00 | 2.5543 | 172 | 50 - 150 | 0.0000 | +/-0.50 | * |
| M2PFTA | 86505.05 | 4.329667 | 479,880.00 | 4.329667 | 18 | 50 - 150 | 0.0000 | +/-0.50 | * |
| M2-8:2FTS | 50290.09 | 3.794817 | 33,212.00 | 3.794817 | 151 | 50 - 150 | 0.0000 | +/-0.50 | * |
| MPFBA | 210934.5 | 1.058467 | 256,957.00 | 1.058467 | 82 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3HFPO-DA | 82571.76 | 2.872033 | 114,902.00 | 2.872033 | 72 | 50 - 150 | 0.0000 | +/-0.50 | |
| M6PFDA | 434129.1 | 3.795333 | 445,919.00 | 3.795333 | 97 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PFBS | 113722.9 | 1.928117 | 104,197.00 | 1.9281 | 109 | 50 - 150 | 0.0000 | +/-0.50 | |
| M7PFUnA | 394559.8 | 3.946033 | 453,308.00 | 3.946033 | 87 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-6:2FTS | 65097.14 | 3.437283 | 27,565.00 | 3.437283 | 236 | 50 - 150 | 0.0000 | +/-0.50 | * |
| M5PFPeA | 297566.3 | 1.749417 | 276,869.00 | 1.749417 | 107 | 50 - 150 | 0.0000 | +/-0.50 | |
| M5PFHxA | 484621.1 | 2.638533 | 458,596.00 | 2.638533 | 106 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PFHxS | 69047.38 | 3.21025 | 68,806.00 | 3.21025 | 100 | 50 - 150 | 0.0000 | +/-0.50 | |
| M4PFHpA | 508716.6 | 3.186933 | 461,168.00 | 3.186933 | 110 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOA | 550241.1 | 3.445817 | 508,809.00 | 3.445817 | 108 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOS | 66070.52 | 3.636183 | 76,995.00 | 3.636183 | 86 | 50 - 150 | 0.0000 | +/-0.50 | |
| M9PFNA | 480064.4 | 3.637217 | 526,406.00 | 3.637217 | 91 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPFDoA | 229361 | 4.088634 | 386,713.00 | 4.088634 | 59 | 50 - 150 | 0.0000 | +/-0.50 | |
| D5-NEtFOSAA | 101208.3 | 3.9535 | 101,789.00 | 3.9535 | 99 | 50 - 150 | 0.0000 | +/-0.50 | |
| D3-NMeFOSAA | 110600.4 | 3.873767 | 116,586.00 | 3.873767 | 95 | 50 - 150 | 0.0000 | +/-0.50 | |

INTERNAL STANDARD AREA AND RT SUMMARY
SOP-454 PFAS

| Internal Standard | Response | RT | Reference Response | Reference RT | Area % | Area % Limits | RT Diff | RT Diff Limit | Q |
|----------------------------|----------|----------|---------------------------|--------------|--------|--------------------------|---------|---------------|---|
| MW-2C (23F0005-02) | | | Lab File ID: 23F0005-02.d | | | Analyzed: 06/15/23 18:48 | | | |
| M8FOSA | 150695 | 3.980567 | 227,522.00 | 3.980567 | 66 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-4:2FTS | 203238 | 2.529667 | 27,028.00 | 2.5543 | 752 | 50 - 150 | -0.0246 | +/-0.50 | * |
| M2PF _n TA | 26303.48 | 4.329667 | 479,880.00 | 4.329667 | 05 | 50 - 150 | 0.0000 | +/-0.50 | * |
| M2-8:2FTS | 89108.22 | 3.794817 | 33,212.00 | 3.794817 | 268 | 50 - 150 | 0.0000 | +/-0.50 | * |
| MPFBA | 67793.53 | 1.04185 | 256,957.00 | 1.058467 | 26 | 50 - 150 | -0.0166 | +/-0.50 | * |
| M3HFPO-DA | 83848.29 | 2.86385 | 114,902.00 | 2.872033 | 73 | 50 - 150 | -0.0082 | +/-0.50 | |
| M6PFDA | 443991.9 | 3.79535 | 445,919.00 | 3.795333 | 100 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PFBS | 94608.18 | 1.911533 | 104,197.00 | 1.9281 | 91 | 50 - 150 | -0.0166 | +/-0.50 | |
| M7PFU _n A | 289061.9 | 3.946033 | 453,308.00 | 3.946033 | 64 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-6:2FTS | 134425.8 | 3.4293 | 27,565.00 | 3.437283 | 488 | 50 - 150 | -0.0080 | +/-0.50 | * |
| M5PFPeA | 216588.3 | 1.731383 | 276,869.00 | 1.749417 | 78 | 50 - 150 | -0.0180 | +/-0.50 | |
| M5PFHxA | 501152.8 | 2.621617 | 458,596.00 | 2.638533 | 109 | 50 - 150 | -0.0169 | +/-0.50 | |
| M3PFHxS | 71434.09 | 3.21025 | 68,806.00 | 3.21025 | 104 | 50 - 150 | 0.0000 | +/-0.50 | |
| M4PFHpA | 502491.8 | 3.17885 | 461,168.00 | 3.186933 | 109 | 50 - 150 | -0.0081 | +/-0.50 | |
| M8PFOA | 444929.9 | 3.445817 | 508,809.00 | 3.445817 | 87 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOS | 60936.63 | 3.636183 | 76,995.00 | 3.636183 | 79 | 50 - 150 | 0.0000 | +/-0.50 | |
| M9PFNA | 403638.6 | 3.637217 | 526,406.00 | 3.637217 | 77 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPFDoA | 85726.39 | 4.088634 | 386,713.00 | 4.088634 | 22 | 50 - 150 | 0.0000 | +/-0.50 | * |
| D5-NEtFOSAA | 74701.45 | 3.9535 | 101,789.00 | 3.9535 | 73 | 50 - 150 | 0.0000 | +/-0.50 | |
| D3-NMeFOSAA | 83922.8 | 3.873767 | 116,586.00 | 3.873767 | 72 | 50 - 150 | 0.0000 | +/-0.50 | |

INTERNAL STANDARD AREA AND RT SUMMARY
SOP-454 PFAS

| Internal Standard | Response | RT | Reference Response | Reference RT | Area % | Area % Limits | RT Diff | RT Diff Limit | Q |
|--|----------|----------|--------------------|--------------|--------|---------------|---------|---------------|---|
| MW-3 (23F0005-03) | | | | | | | | | |
| Lab File ID: 23F0005-03.d Analyzed: 06/15/23 18:56 | | | | | | | | | |
| M8FOSA | 144129.3 | 3.980567 | 227,522.00 | 3.980567 | 63 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-4:2FTS | 31890.78 | 2.5543 | 27,028.00 | 2.5543 | 118 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2PF _T A | 4088.255 | 4.329667 | 479,880.00 | 4.329667 | 01 | 50 - 150 | 0.0000 | +/-0.50 | * |
| M2-8:2FTS | 19103.5 | 3.794817 | 33,212.00 | 3.794817 | 58 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPFBA | 247818.5 | 1.058467 | 256,957.00 | 1.058467 | 96 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3HFPO-DA | 106199.2 | 2.872033 | 114,902.00 | 2.872033 | 92 | 50 - 150 | 0.0000 | +/-0.50 | |
| M6PFDA | 257111.8 | 3.795333 | 445,919.00 | 3.795333 | 58 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PFBS | 116013.5 | 1.9364 | 104,197.00 | 1.9281 | 111 | 50 - 150 | 0.0083 | +/-0.50 | |
| M7PFU _n A | 100458.5 | 3.946033 | 453,308.00 | 3.946033 | 22 | 50 - 150 | 0.0000 | +/-0.50 | * |
| M2-6:2FTS | 34820.13 | 3.437283 | 27,565.00 | 3.437283 | 126 | 50 - 150 | 0.0000 | +/-0.50 | |
| M5PFPeA | 293698.8 | 1.749417 | 276,869.00 | 1.749417 | 106 | 50 - 150 | 0.0000 | +/-0.50 | |
| M5PFH _x A | 485765.5 | 2.646767 | 458,596.00 | 2.638533 | 106 | 50 - 150 | 0.0082 | +/-0.50 | |
| M3PFH _x S | 73167.49 | 3.21025 | 68,806.00 | 3.21025 | 106 | 50 - 150 | 0.0000 | +/-0.50 | |
| M4PFH _p A | 506378.7 | 3.186933 | 461,168.00 | 3.186933 | 110 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOA | 532779.1 | 3.445817 | 508,809.00 | 3.445817 | 105 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOS | 50709.96 | 3.636183 | 76,995.00 | 3.636183 | 66 | 50 - 150 | 0.0000 | +/-0.50 | |
| M9PFNA | 420872.5 | 3.637217 | 526,406.00 | 3.637217 | 80 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPFDoA | 18804.53 | 4.088634 | 386,713.00 | 4.088634 | 05 | 50 - 150 | 0.0000 | +/-0.50 | * |
| D5-NEtFOSAA | 24872.35 | 3.9535 | 101,789.00 | 3.9535 | 24 | 50 - 150 | 0.0000 | +/-0.50 | * |
| D3-NMeFOSAA | 41892.87 | 3.873767 | 116,586.00 | 3.873767 | 36 | 50 - 150 | 0.0000 | +/-0.50 | * |

INTERNAL STANDARD AREA AND RT SUMMARY
SOP-454 PFAS

| Internal Standard | Response | RT | Reference Response | Reference RT | Area % | Area % Limits | RT Diff | RT Diff Limit | Q |
|--|----------|----------|--------------------|--------------|--------|---------------|---------|---------------|---|
| MW-5 (23F0005-04) | | | | | | | | | |
| Lab File ID: 23F0005-04.d Analyzed: 06/15/23 19:03 | | | | | | | | | |
| M8FOSA | 126189.5 | 3.980567 | 227,522.00 | 3.980567 | 55 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-4:2FTS | 51181.77 | 2.5543 | 27,028.00 | 2.5543 | 189 | 50 - 150 | 0.0000 | +/-0.50 | * |
| M2PF _{TA} | 13962.42 | 4.329667 | 479,880.00 | 4.329667 | 03 | 50 - 150 | 0.0000 | +/-0.50 | * |
| M2-8:2FTS | 32987.29 | 3.794817 | 33,212.00 | 3.794817 | 99 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPFBA | 147895.7 | 1.058467 | 256,957.00 | 1.058467 | 58 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3HFPO-DA | 87246.78 | 2.872033 | 114,902.00 | 2.872033 | 76 | 50 - 150 | 0.0000 | +/-0.50 | |
| M6PFDA | 342832.7 | 3.795333 | 445,919.00 | 3.795333 | 77 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PFBS | 107472.1 | 1.928117 | 104,197.00 | 1.9281 | 103 | 50 - 150 | 0.0000 | +/-0.50 | |
| M7PFU _{nA} | 170601.8 | 3.946033 | 453,308.00 | 3.946033 | 38 | 50 - 150 | 0.0000 | +/-0.50 | * |
| M2-6:2FTS | 53191.43 | 3.4373 | 27,565.00 | 3.437283 | 193 | 50 - 150 | 0.0000 | +/-0.50 | * |
| M5PFPeA | 277750.1 | 1.749417 | 276,869.00 | 1.749417 | 100 | 50 - 150 | 0.0000 | +/-0.50 | |
| M5PFH _x A | 485465.7 | 2.638533 | 458,596.00 | 2.638533 | 106 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PFH _x S | 72394.59 | 3.218333 | 68,806.00 | 3.21025 | 105 | 50 - 150 | 0.0081 | +/-0.50 | |
| M4PFH _p A | 512351.3 | 3.186933 | 461,168.00 | 3.186933 | 111 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOA | 557249.2 | 3.445817 | 508,809.00 | 3.445817 | 110 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOS | 60354.33 | 3.636183 | 76,995.00 | 3.636183 | 78 | 50 - 150 | 0.0000 | +/-0.50 | |
| M9PFNA | 483919 | 3.637217 | 526,406.00 | 3.637217 | 92 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPFDoA | 50278.58 | 4.088634 | 386,713.00 | 4.088634 | 13 | 50 - 150 | 0.0000 | +/-0.50 | * |
| D5-NEtFOSAA | 38104.12 | 3.9535 | 101,789.00 | 3.9535 | 37 | 50 - 150 | 0.0000 | +/-0.50 | * |
| D3-NMeFOSAA | 58924.64 | 3.873767 | 116,586.00 | 3.873767 | 51 | 50 - 150 | 0.0000 | +/-0.50 | |

INTERNAL STANDARD AREA AND RT SUMMARY

SOP-454 PFAS

| Internal Standard | Response | RT | Reference Response | Reference RT | Area % | Area % Limits | RT Diff | RT Diff Limit | Q |
|---------------------------------|----------|----------|---------------------------|--------------|--------|--------------------------|---------|---------------|---|
| Brandon ERB (23F0005-05) | | | | | | | | | |
| | | | Lab File ID: 23F0005-05.d | | | Analyzed: 06/15/23 19:10 | | | |
| M8FOSA | 215138.4 | 3.980567 | 227,522.00 | 3.980567 | 95 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-4:2FTS | 27811.8 | 2.5543 | 27,028.00 | 2.5543 | 103 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2PF _{TA} | 520295.8 | 4.329667 | 479,880.00 | 4.329667 | 108 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-8:2FTS | 40282.16 | 3.794817 | 33,212.00 | 3.794817 | 121 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPF _{BA} | 246271.6 | 1.058467 | 256,957.00 | 1.058467 | 96 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3HFPO-DA | 98507.22 | 2.872033 | 114,902.00 | 2.872033 | 86 | 50 - 150 | 0.0000 | +/-0.50 | |
| M6PF _{DA} | 473234.7 | 3.795333 | 445,919.00 | 3.795333 | 106 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PF _{BS} | 110777.4 | 1.9281 | 104,197.00 | 1.9281 | 106 | 50 - 150 | 0.0000 | +/-0.50 | |
| M7PF _{UnA} | 547202.7 | 3.938033 | 453,308.00 | 3.946033 | 121 | 50 - 150 | -0.0080 | +/-0.50 | |
| M2-6:2FTS | 37692.93 | 3.437283 | 27,565.00 | 3.437283 | 137 | 50 - 150 | 0.0000 | +/-0.50 | |
| M5PF _{PeA} | 271501.3 | 1.749417 | 276,869.00 | 1.749417 | 98 | 50 - 150 | 0.0000 | +/-0.50 | |
| M5PF _{HxA} | 463810.7 | 2.638533 | 458,596.00 | 2.638533 | 101 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PF _{HxS} | 71209.7 | 3.21025 | 68,806.00 | 3.21025 | 103 | 50 - 150 | 0.0000 | +/-0.50 | |
| M4PF _{HpA} | 482628.4 | 3.186933 | 461,168.00 | 3.186933 | 105 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOA | 543119.2 | 3.445817 | 508,809.00 | 3.445817 | 107 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOS | 76222.82 | 3.636183 | 76,995.00 | 3.636183 | 99 | 50 - 150 | 0.0000 | +/-0.50 | |
| M9PF _{NA} | 526302.8 | 3.637217 | 526,406.00 | 3.637217 | 100 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPF _{DoA} | 395622.5 | 4.08065 | 386,713.00 | 4.088634 | 102 | 50 - 150 | -0.0080 | +/-0.50 | |
| D5-NEtFOSAA | 134570.4 | 3.9455 | 101,789.00 | 3.9535 | 132 | 50 - 150 | -0.0080 | +/-0.50 | |
| D3-NMeFOSAA | 121108.1 | 3.8656 | 116,586.00 | 3.873767 | 104 | 50 - 150 | -0.0082 | +/-0.50 | |

INTERNAL STANDARD AREA AND RT SUMMARY
SOP-454 PFAS

| Internal Standard | Response | RT | Reference Response | Reference RT | Area % | Area % Limits | RT Diff | RT Diff Limit | Q |
|-----------------------------|----------|----------|-----------------------------|--------------|--------|--------------------------|---------|---------------|---|
| Blank (B343017-BLK1) | | | Lab File ID: B343017-BLK1.d | | | Analyzed: 06/15/23 18:34 | | | |
| M8FOSA | 215016.1 | 3.980567 | 227,522.00 | 3.980567 | 95 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-4:2FTS | 30333.71 | 2.5543 | 27,028.00 | 2.5543 | 112 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2PFTA | 436933.4 | 4.329667 | 479,880.00 | 4.329667 | 91 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-8:2FTS | 36123.36 | 3.794817 | 33,212.00 | 3.794817 | 109 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPFBA | 279710.7 | 1.058467 | 256,957.00 | 1.058467 | 109 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3HFPO-DA | 104600.4 | 2.872033 | 114,902.00 | 2.872033 | 91 | 50 - 150 | 0.0000 | +/-0.50 | |
| M6PFDA | 473304.3 | 3.795333 | 445,919.00 | 3.795333 | 106 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PFBS | 123107.8 | 1.9364 | 104,197.00 | 1.9281 | 118 | 50 - 150 | 0.0083 | +/-0.50 | |
| M7PFUnA | 457712 | 3.946033 | 453,308.00 | 3.946033 | 101 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-6:2FTS | 35864.22 | 3.437283 | 27,565.00 | 3.437283 | 130 | 50 - 150 | 0.0000 | +/-0.50 | |
| M5PFPeA | 310102.7 | 1.749417 | 276,869.00 | 1.749417 | 112 | 50 - 150 | 0.0000 | +/-0.50 | |
| M5PFHxA | 520505 | 2.638533 | 458,596.00 | 2.638533 | 113 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PFHxS | 78885.7 | 3.21025 | 68,806.00 | 3.21025 | 115 | 50 - 150 | 0.0000 | +/-0.50 | |
| M4PFHpA | 542553.6 | 3.186933 | 461,168.00 | 3.186933 | 118 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOA | 588328.2 | 3.445817 | 508,809.00 | 3.445817 | 116 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOS | 81873.97 | 3.636183 | 76,995.00 | 3.636183 | 106 | 50 - 150 | 0.0000 | +/-0.50 | |
| M9PFNA | 555643.1 | 3.637217 | 526,406.00 | 3.637217 | 106 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPFDoA | 346001 | 4.088634 | 386,713.00 | 4.088634 | 89 | 50 - 150 | 0.0000 | +/-0.50 | |
| D5-NEtFOSAA | 106083.9 | 3.9535 | 101,789.00 | 3.9535 | 104 | 50 - 150 | 0.0000 | +/-0.50 | |
| D3-NMeFOSAA | 120334.9 | 3.873767 | 116,586.00 | 3.873767 | 103 | 50 - 150 | 0.0000 | +/-0.50 | |

INTERNAL STANDARD AREA AND RT SUMMARY
SOP-454 PFAS

| Internal Standard | Response | RT | Reference Response | Reference RT | Area % | Area % Limits | RT Diff | RT Diff Limit | Q |
|---------------------------|----------|----------|----------------------------|--------------|--------|--------------------------|---------|---------------|---|
| LCS (B343017-BS1) | | | Lab File ID: B343017-BS1.d | | | Analyzed: 06/15/23 18:20 | | | |
| M8FOSA | 200169.3 | 3.980567 | 227,522.00 | 3.980567 | 88 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-4:2FTS | 27979.2 | 2.5543 | 27,028.00 | 2.5543 | 104 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2PFTA | 422371.8 | 4.329667 | 479,880.00 | 4.329667 | 88 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-8:2FTS | 31719.93 | 3.794817 | 33,212.00 | 3.794817 | 96 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPFBA | 257064.5 | 1.058467 | 256,957.00 | 1.058467 | 100 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3HFPO-DA | 110701.5 | 2.872033 | 114,902.00 | 2.872033 | 96 | 50 - 150 | 0.0000 | +/-0.50 | |
| M6PFDA | 431297 | 3.795333 | 445,919.00 | 3.795333 | 97 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PFBS | 111061.8 | 1.9364 | 104,197.00 | 1.9281 | 107 | 50 - 150 | 0.0083 | +/-0.50 | |
| M7PFUnA | 426879 | 3.946033 | 453,308.00 | 3.946033 | 94 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-6:2FTS | 29169.89 | 3.437283 | 27,565.00 | 3.437283 | 106 | 50 - 150 | 0.0000 | +/-0.50 | |
| M5PFPeA | 289508.5 | 1.749417 | 276,869.00 | 1.749417 | 105 | 50 - 150 | 0.0000 | +/-0.50 | |
| M5PFHxA | 469620.1 | 2.646767 | 458,596.00 | 2.638533 | 102 | 50 - 150 | 0.0082 | +/-0.50 | |
| M3PFHxS | 72817.55 | 3.218333 | 68,806.00 | 3.21025 | 106 | 50 - 150 | 0.0081 | +/-0.50 | |
| M4PFHpA | 489111.9 | 3.186933 | 461,168.00 | 3.186933 | 106 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOA | 521309.5 | 3.453817 | 508,809.00 | 3.445817 | 102 | 50 - 150 | 0.0080 | +/-0.50 | |
| M8PFOS | 74040.57 | 3.636183 | 76,995.00 | 3.636183 | 96 | 50 - 150 | 0.0000 | +/-0.50 | |
| M9PFNA | 527230 | 3.637217 | 526,406.00 | 3.637217 | 100 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPFDoA | 344025.2 | 4.088634 | 386,713.00 | 4.088634 | 89 | 50 - 150 | 0.0000 | +/-0.50 | |
| D5-NEtFOSAA | 93220.06 | 3.9535 | 101,789.00 | 3.9535 | 92 | 50 - 150 | 0.0000 | +/-0.50 | |
| D3-NMeFOSAA | 108673.3 | 3.873767 | 116,586.00 | 3.873767 | 93 | 50 - 150 | 0.0000 | +/-0.50 | |

INTERNAL STANDARD AREA AND RT SUMMARY
SOP-454 PFAS

| Internal Standard | Response | RT | Reference Response | Reference RT | Area % | Area % Limits | RT Diff | RT Diff Limit | Q |
|-------------------------------|----------|----------|-----------------------------|--------------|--------|--------------------------|---------|---------------|---|
| LCS Dup (B343017-BSD1) | | | Lab File ID: B343017-BSD1.d | | | Analyzed: 06/15/23 18:27 | | | |
| M8FOSA | 194266.3 | 3.980567 | 227,522.00 | 3.980567 | 85 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-4:2FTS | 27529.69 | 2.5543 | 27,028.00 | 2.5543 | 102 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2PFTA | 426555.5 | 4.329667 | 479,880.00 | 4.329667 | 89 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-8:2FTS | 32444.97 | 3.794817 | 33,212.00 | 3.794817 | 98 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPFBA | 258494.3 | 1.058467 | 256,957.00 | 1.058467 | 101 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3HFPO-DA | 103649.6 | 2.872033 | 114,902.00 | 2.872033 | 90 | 50 - 150 | 0.0000 | +/-0.50 | |
| M6PFDA | 420816.1 | 3.795333 | 445,919.00 | 3.795333 | 94 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PFBS | 113307.8 | 1.9281 | 104,197.00 | 1.9281 | 109 | 50 - 150 | 0.0000 | +/-0.50 | |
| M7PFUnA | 409909.9 | 3.946033 | 453,308.00 | 3.946033 | 90 | 50 - 150 | 0.0000 | +/-0.50 | |
| M2-6:2FTS | 32701.71 | 3.437283 | 27,565.00 | 3.437283 | 119 | 50 - 150 | 0.0000 | +/-0.50 | |
| M5PFPeA | 289122.8 | 1.749417 | 276,869.00 | 1.749417 | 104 | 50 - 150 | 0.0000 | +/-0.50 | |
| M5PFHxA | 477646.3 | 2.638533 | 458,596.00 | 2.638533 | 104 | 50 - 150 | 0.0000 | +/-0.50 | |
| M3PFHxS | 74887.8 | 3.21025 | 68,806.00 | 3.21025 | 109 | 50 - 150 | 0.0000 | +/-0.50 | |
| M4PFHpA | 485196.5 | 3.186933 | 461,168.00 | 3.186933 | 105 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOA | 502057.2 | 3.445817 | 508,809.00 | 3.445817 | 99 | 50 - 150 | 0.0000 | +/-0.50 | |
| M8PFOS | 70558.66 | 3.636183 | 76,995.00 | 3.636183 | 92 | 50 - 150 | 0.0000 | +/-0.50 | |
| M9PFNA | 493344.9 | 3.637217 | 526,406.00 | 3.637217 | 94 | 50 - 150 | 0.0000 | +/-0.50 | |
| MPFDoA | 344756.5 | 4.088634 | 386,713.00 | 4.088634 | 89 | 50 - 150 | 0.0000 | +/-0.50 | |
| D5-NEtFOSAA | 97125.26 | 3.9535 | 101,789.00 | 3.9535 | 95 | 50 - 150 | 0.0000 | +/-0.50 | |
| D3-NMeFOSAA | 104454.8 | 3.873767 | 116,586.00 | 3.873767 | 90 | 50 - 150 | 0.0000 | +/-0.50 | |

CERTIFICATIONS
Certified Analyses included in this Report

| Analyte | Certifications |
|--|----------------|
| <i>SOP-454 PFAS in Water</i> | |
| Perfluorobutanoic acid (PFBA) | NH-P,PA |
| Perfluorobutanesulfonic acid (PFBS) | NH-P,PA |
| Perfluoropentanoic acid (PFPeA) | NH-P,PA |
| Perfluorohexanoic acid (PFHxA) | NH-P,PA |
| 11Cl-PF3OUdS (F53B Major) | NH-P,PA |
| 9Cl-PF3ONS (F53B Minor) | NH-P,PA |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA) | NH-P,PA |
| Hexafluoropropylene oxide dimer acid (HFPO-DA) | NH-P,PA |
| 8:2 Fluorotelomersulfonic acid (8:2FTS A) | NH-P,PA |
| Perfluorodecanoic acid (PFDA) | NH-P,PA |
| Perfluorododecanoic acid (PFDoA) | NH-P,PA |
| Perfluoro(2-ethoxyethane)sulfonic acid (PFEEESA) | NH-P,PA |
| Perfluoroheptanesulfonic acid (PFHpS) | NH-P,PA |
| N-EtFOSAA (NEtFOSAA) | NH-P,PA |
| N-MeFOSAA (NMeFOSAA) | NH-P,PA |
| Perfluorotetradecanoic acid (PFTA) | NH-P,PA |
| Perfluorotridecanoic acid (PFTrDA) | NH-P,PA |
| 4:2 Fluorotelomersulfonic acid (4:2FTS A) | NH-P,PA |
| Perfluorodecanesulfonic acid (PFDS) | NH-P,PA |
| Perfluorooctanesulfonamide (FOSA) | NH-P,PA |
| Perfluorononanesulfonic acid (PFNS) | NH-P,PA |
| Perfluoro-1-hexanesulfonamide (FHxSA) | NH-P,PA |
| Perfluoro-1-butanesulfonamide (FBSA) | NH-P,PA |
| Perfluorohexanesulfonic acid (PFHxS) | NH-P,PA |
| Perfluoro-4-oxapentanoic acid (PFMPA) | NH-P,PA |
| Perfluoro-5-oxahexanoic acid (PFMBA) | NH-P,PA |
| 6:2 Fluorotelomersulfonic acid (6:2FTS A) | NH-P,PA |
| Perfluoropentanesulfonic acid (PFPeS) | NH-P,PA |
| Perfluoroundecanoic acid (PFUnA) | NH-P,PA |
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA) | NH-P,PA |
| Perfluoroheptanoic acid (PFHpA) | NH-P,PA |
| Perfluorooctanoic acid (PFOA) | NH-P,PA |
| Perfluorooctanesulfonic acid (PFOS) | NH-P,PA |
| Perfluorononanoic acid (PFNA) | NH-P,PA |

Con-Test, a Pace Environmental Laboratory, operates under the following certifications and accreditations:

| Code | Description | Number | Expires |
|------|----------------------------------|------------|------------|
| NH-P | New Hampshire Environmental Lab | 2557 NELAP | 09/6/2023 |
| PA | Commonwealth of Pennsylvania DEP | 68-05812 | 06/30/2024 |

39 Spruce St.
East Longmeadow, MA. 01028
P: 413-525-2332
F: 413-525-6405
www.pacelabs.com

ENV-FRM-ELON-0001 V05__ Sample Receiving Checklist

Log In Back-Sheet

Login Sample Receipt Checklist - (Rejection Criteria Listing
- Using Acceptance Policy) Any False statement will be
brought to the attention of the Client - True or False



Client Acis Inc
Project Brankon Landfill
MCP/RCP Required N/A
Deliverable Package Requirement N/A
Location Brankon, VT
PWSID# (When Applicable) N/A
Arrival Method:
Courier Fed Ex Walk In Other
Received By / Date / Time ME M 5/31/23 14:25
Back Sheet By / Date / Time LA 5/11/23 8:56
Temperature Method gun # 5
Temp < 6° C Actual Temperature 5.3/5.7
Rush Samples: Yes / No Notify _____
Short Hold: Yes / No Notify _____

| | True | False |
|---|--|--|
| Received on Ice | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Received in Cooler | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Custody Seal: DATE TIME | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| COC Relinquished | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| COC/Samples Labels Agree | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| All Samples in Good Condition | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Samples Received within Holding Time | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Is there enough Volume | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Proper Media/Container Used | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Splitting Samples Required | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| MS/MSD | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| Trip Blanks | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| Lab to Filters | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| COC Legible | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| COC Included: (Check all included) | | |
| Client <input checked="" type="checkbox"/> | Analysis <input checked="" type="checkbox"/> | Sampler Name <input checked="" type="checkbox"/> |
| Project <input checked="" type="checkbox"/> | IDs <input checked="" type="checkbox"/> | Collection Date/Time <input checked="" type="checkbox"/> |
| All Samples Proper pH: | <input checked="" type="checkbox"/> N/A | <input type="checkbox"/> |

Notes regarding Samples/COC outside of SOP:

Additional Container Notes

