



# September 2022 ANNUAL/SEMI-ANNUAL GROUNDWATER MONITORING REPORT NIAGARA WIND POWER, LLC STEEL WINDS I FACILITY (Site No. C915205) LACKAWANNA, NEW YORK

November 2022 (Revised January 2023) File No. 03.0033579.15



# **PREPARED FOR:**

Niagara Wind Power, LLC 200 Liberty Street, 14<sup>th</sup> Fl. NY, NY 10281

# **GZA GeoEnvironmental of New York**

300 Pearl Street Suite 700 | Buffalo, NY 14202 716-685-2300

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November 16, 2022, Revised January 9, 2023 File No. 03.0033579.15

Mr. Cris Basden **Brookfield Renewable** c/o Niagara Wind Power, LLC 200 Liberty Street, 14th Floor New York, NY 10281

Re: September 2022 Annual/Semi-Annual Groundwater Monitoring Report

Steel Winds I Site (Site No. C915205)

Lackawanna, NY

Dear Mr. Basden:

GZA GeoEnvironmental (GZA) is pleased to submit this annual/semi-annual groundwater monitoring report to Niagara Wind Power, LLC (NWP) summarizing the analytical results of the groundwater sampling event conducted in September 2022 at the above referenced Site. The objective of the sampling event was to collect and analyze groundwater samples from the on-site monitoring wells in accordance with the Site Management Plan, dated September 2007, prepared by Benchmark Environmental Engineering and Science, PLLC (Benchmark) and approved by the New York State Department of Environmental Conservation (NYSDEC).

Should you have any questions or require additional information following your review, please contact Daniel Troy at (716) 570-6673 or Ed Summerly at (401) 427-2707.

Richard A. Carlone, P.E

Consultant Reviewer

Sincerely,

GZA GEOENVIRONMENTAL OF NEW YORK

Senior Project Manager

Edward A. Summerly, P.G.

Principal / District Office Manager

Matt Carson (Brookfield Renewable) cc:

Megan Kuczka (NYSDEC)

Attachments: Report



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#### 1.00 INTRODUCTION

In accordance with our March 16, 2022 proposal, GZA GeoEnvironmental, Inc. (GZA) collected and analyzed groundwater samples at the nine (9) annual site-wide groundwater monitoring well locations (designated the Long-Term Groundwater Monitoring Plan (LTGWM)) and the six (6) semi-annual WT-1 vicinity groundwater monitoring well locations at the Steel Winds I facility located in Lackawanna, New York (Site). A Locus Plan and Site Plan are attached as **Figures 1** and **2**, respectively.

#### 1.10 BACKGROUND AND SITE HISTORY

Tecumseh Redevelopment, Inc. (Tecumseh) owns approximately 1,100 acres of land at 1951 Hamburg Turnpike, as shown on attached **Figure 1**. The property was formerly used for the production of steel, coke and related products by Bethlehem Steel Corporation (BSC). Steel production on the Tecumseh property was discontinued in 1983 and the coke ovens ceased activity in 2000. Tecumseh acquired the property, along with other BSC assets, out of bankruptcy in 2003.

In September 2006, BQ Energy entered into a long-term lease agreement with Tecumseh to construct and operate wind turbines and supporting power generation equipment and infrastructure on an approximately 29-acre parcel of the Tecumseh property, referred to as the Steel Winds I Site. BQ Energy and the NYSDEC also entered into a Brownfield Cleanup Agreement for the Steel Winds Site. The Site is wholly contained within the Slag Fill Area (SFA) Zones 3 and 4 of the Tecumseh property bordered by Lake Erie to the west, Smoke Creek to the south, and former industrial lands of BSC to the north and east. Niagara Wind Power, LLC (NWP) an affiliate of Brookfield Renewable, operates the eight wind turbines installed at the Site. In accordance with an October 30, 2020 letter to NYSDEC, Niagara Wind Power, LLC assumed the Remedial Party status for the Site.

The Brownfield Cleanup Program (BCP) was successful in achieving the remedial objectives for the Steel Winds Site. The Site Management Plan (SMP) and Final Engineering Report (FER) were approved by NYSDEC in December 2007. NYSDEC issued a Certificate of Completion (COC) for the Site on December 18, 2007.

The remedial activities conducted at the Site include:

- Excavation and off-site disposal of impacted slag fill from the eight wind turbine foundations and interconnecting utility trenches;
- In-situ enhanced biodegradation of residual volatile organic compounds (VOCs), including benzene, toluene, total xylenes, and naphthalene, using oxygen release compound (ORC®) socks within the saturated soil and groundwater in the vicinity of monitoring well WT-01 and associated groundwater monitoring; and,
- Completion of a soil cover system (cap).



As a requirement of the SMP, LTGWM is being performed at nine (9) wells across the Site. Additional groundwater monitoring was also performed to monitor the effectiveness of the ORC in-situ treatment in the vicinity of wind turbine WT-01. During 2011, both the LTGWM and WT-01 vicinity groundwater monitoring programs were performed on an annual basis and were done on July 13 and 14, 2011. The five ORC in-situ treatment wells were to be monitored semi-annually, in accordance with the SMP. However, only one ORC monitoring event (on May 4, 2011) was conducted because of the ineffectiveness of the remedy.

An *Operation, Monitoring and Maintenance Request for Modification* report, dated November 2011, was submitted to NYSDEC by Benchmark. This report proposed ceasing operation of ORC® groundwater remedy for the WT-01 Vicinity because the remedy was not effective in reducing VOC concentrations, due primarily to the geochemical conditions (i.e., high baseline chemical oxygen demand, highly negative oxidation reduction potential and high pH) of the Site. NYSDEC provided comments to this report on April 10, 2012 and GZA provided a response letter on May 9, 2012. Based on this letter and correspondence with NYSDEC, the ORC® remedy has been terminated (i.e., the ORC socks have been removed from the five treatment wells and disposed of as a solid waste).

In accordance with a letter from GZA to NYSDEC, dated June 22, 2012<sup>1</sup>, semi-annual/annual groundwater monitoring will continue at the Site until a Technical Impracticability Waiver (TI Waiver) for groundwater treatment at the Site is submitted to, and approved by NYSDEC.

On September 30, 2013, GZA submitted a *Technical Impracticability Waiver Supplemental Field Studies Work Plan* for the Site, detailing sampling, laboratory analysis, data evaluation and reporting to be conducted in support of a TI Waiver request for the Site. This *Work Plan* was approved by NYSDEC on February 24, 2014. Sampling and analysis described in the *Work Plan* was conducted by GZA in summer 2014 and a TI Waiver application was submitted to NYSDEC on November 5, 2014, with a supplemental *Endangered Species Review* letter submitted to NYSDEC on January 28, 2015. Based on the remedial evaluation presented in the application, it is GZA's opinion that active remediation is not warranted or feasible, would not result in significant benefit to the environment relative to the cost, and is technically impracticable. The application recommended limited additional sampling to evaluate risk to ecological receptors. NYSDEC verbally approved the additional recommended field work on April 27, 2015. GZA submitted a *Work Plan* to NYSDEC on August 5, 2015 describing the proposed additional field work, which was implemented in September 2015. The *TI Waiver Supplemental Report* was submitted to NYSDEC on April 24, 2018.

Due to the length of cold days experienced during the winter of 2014-2015 the semi-annual sampling event, originally scheduled for January 2015, was not able to be completed until March 2015. In order to reduce negative impacts and delays associated from freezing weather conditions, NYSDEC approved rescheduling the future semi-annual and annual sampling events to occur during the months of March and September, respectively.

<sup>&</sup>lt;sup>1</sup>GZA's June 22, 2012 letter was prepared in response to NYSDEC's comments on GZA's May 9, 2012 Responses to NYSDEC's April 10, 2012 Comments on the November 2011 *Operation, Monitoring and Maintenance Request for Modification*, prepared by Benchmark.



#### 2.00 PURPOSE AND SCOPE OF WORK

The purpose of the September 2022 annual/semi-annual sampling event was to collect groundwater samples from the nine (9) annual site-wide and six (6) semi-annual WT-1 vicinity groundwater monitoring wells, respectively, in accordance with the routine monitoring protocol described in the September 2007 SMP. To accomplish this, the following activities were completed by GZA:

- Collected one (1) groundwater sample from each annual/semi-annual monitoring well location for laboratory analysis (conducted by Alpha Analytical of Westborough, Massachusetts) in accordance with the analytical testing summary provided in **Table 1**. Test parameters included the following:
  - STARS list VOCs via EPA Method 8260B;
  - o Base-Neutral semi-volatile organic compounds (SVOCs) via EPA Method 8270C; and
  - Arsenic, barium, chromium, and/or manganese via EPA Method 6010B (select annual groundwater monitoring wells only).
- Prepared this report, which summarizes the data collected during each sampling event and compared the current results to historic data and assessed contaminant concentration trends.

This report presents GZA's field observations, results, and opinions and is subject to the limitations presented in **Appendix A** and modifications if subsequent information is developed by GZA or any other party.

#### 3.00 FIELD STUDIES

This section describes the field studies conducted as part of GZA's groundwater annual/semi-annual sampling event.

#### 3.10 GROUNDWATER DATA COLLECTION

GZA collected groundwater samples from the nine (9) annual Site-wide monitoring wells (MWN-01, MWN-01B, MWN-02, MWN-02B, MWN-02D, MWN-03, MWN-03B, MWN-03D, and MWN-04), and six (6) WT-1 vicinity semi-annual monitoring wells (MWN-01, MWN-01B, WT1-02, WT1-04, WT1-05, and BCP-ORC-1). Samples were collected on September 13 through 15, 2022. Note, when the two monitoring programs included the same wells, only one sample was collected, and that analysis was used for both programs.



The following tables show the volume of water purged in gallons and the number of well volumes removed from the respective well after a constant head was established. In general, groundwater purge rates were 500(±) millimeter per minute (ml/min). We note that due to complications experienced with the downhole pump in monitoring wells MWN-2D, MWN-3, MW-3B, MWN-03D and MWN-4, alternative sampling methods using a dedicated bailer to remove three well volumes were required for sample collection. The groundwater samples collected using this method were generally observed to have increased turbidity, which required laboratory filtration (from unpreserved samples) prior to inorganic analysis. Well development forms for each monitoring well sampled are included in **Appendix D**.

Annual Site-Wide Monitoring Well ID	Cumulative Volume Purged (gallons)	Well Volumes (#)
MWN-01	25	10
MWN-01B	10	3.7
MWN-02	10	3.0
MWN-02B	10	2.3
MWN-02D	24*	3.0
MWN-03	19*	3.0
MWN-03B	15*	3.1
MWN-03D	39*	3.0
MWN-04	6**	1.3

WT-1 Vicinity Semi-Annual Monitoring Well ID	Cumulative Volume Purged (gallons)	Well Volumes (#)
MWN-01	25	10
MWN-01B	10	3.7
WT1-02	4	0.6
WT1-04	2.5	1.3
WT1-05	19	11.2
BCP-ORC-1	2.0	0.8

Note: wells highlighted in yellow are included in both programs.

As part of the Annual/Semi-annual groundwater monitoring round, static groundwater level measurements were made from top of riser of the monitoring wells listed in the table below prior to purging. Monitoring point elevation data was available from previous groundwater monitoring reports completed by Benchmark, and/or field survey work conducted by GZA. From this data, groundwater flow directions were estimated and are shown on **Figure 2**. Based on the available information, groundwater flow is generally in a westerly direction towards Lake Erie or south toward Smoke Creek (in the immediate vicinity of Smoke Creek only).

<sup>\*</sup>Well was unable to be purged via low flow methods and had 3 well volumes removed with a dedicated bailer.

<sup>\*\*</sup>Well bailed dry and allowed to recharge for 1 hour prior to sample collection.



Monitoring Well	Top of Riser	Groundwater Depth	Groundwater
Location	Elevation (ft.)	(ft.)	Elevation (ft.)
MWN-01	585.14	15.22	569.92
MWN-01B	587.03	15.22	571.81
MWN-02	601.01	28.49	572.52
MWN-02B	601.28	28.73	572.55
MWN-02D	602.95	29.30	573.65
MWN-03	611.96	39.49	572.47
MWN-03B	612.29	39.64	572.65
MWN-03D	613.51	39.49	574.02
MWN-04	623.45	51.36	572.09
WT1-02	600.78	27.82	572.96
WT1-04	586.45	13.76	572.69
WT1-05	584.41	12.48	571.93
BCP-ORC-1	591.97	19.23	572.74

#### 4.00 ANALYTICAL LABORATORY TESTING

Thirteen (13) annual/semi-annual groundwater samples were submitted for analytical testing as part of the September 2022 sampling event. The samples were packed in an ice-filled cooler and, following chain-of-custody procedures, sent to Alpha Analytical for analysis. **Table 1** presents a summary of the samples collected and the analyses completed. As noted above, the samples from monitoring wells MWN-2D, MWN-3, MW-3B, MWN-03D and MWN-4 required laboratory filtering prior to metals analysis as samples were collected with a dedicated bailer and samples from MWN-01 and MWN-01B were included for both semi-annual and annual monitoring programs.

### **5.00 ANALYTICAL TEST RESULTS**

A discussion of the laboratory results for the groundwater samples is presented below. The laboratory reports are provided in **Appendix B** and the analytical test results are summarized on **Tables 2 and 3**.

The analytical test results for the groundwater samples were compared to NYSDEC Class GA criteria presented in the *Division of Water Technical and Operational Guidance Series* (TOGS 1.1.1), dated October 1993, revised June 1998, errata January 1999 and amended April 2000.

The analytical data generated as part of this sampling event has also been provided to NYSDEC electronically for their Environmental Information Management System (EIMS). The data was provided in a standardized electronic data deliverable (EDD) format that uses the database software application EQuIS<sup>™</sup> (EQuIS) from EarthSoft® Inc. The laboratory data and required information were imported into the EQuIS Data Processor (EDP) and submitted to NYSDEC on October 25, 2022.



#### 5.10 ANNUAL SITE-WIDE MONITORING WELLS

- o <u>MWN-01</u> (screen depth: 9.2' 19.2'): Eight (8) VOCs were detected above method reporting limits of which four (4) exceeded their respective NYSDEC Class GA criteria and guidance values, as follows.
  - Benzene at 12 parts per billion (ppb);
  - o m,p-Xylene at 6.0 ppb;
  - o Total Xylene at 11.0 ppb; and

Naphthalene was detected as a VOC at a concentration of 240 ppb, which exceeds its guidance value of 10 ppb.

Sixteen (16) SVOCs were detected above their method reporting limits of which six (6) exceeded their respective NYSDEC Class GA criteria and guidance values, as follows.

- Naphthalene at 91.9 ppb;
- o Phenanthrene at 69.9 ppb;
- o Biphenyl at 6.48 ppb;
- Benz [a] Anthracene at 0.380 (estimated value, i.e., J detect);
- Benzo [b] Fluoranthene at 0.079 (estimated value, i.e., J detect); and
- o Chrysene at 0.214 ppb (estimated value, i.e., J detect).
- MWN-01B (screen depth: 22.2' 32.2'): Ten (10) VOCs were detected above method reporting limits, of which eight (8) exceeded their respective NYSDEC Class GA criteria and guidance values, as follows.
  - Benzene at 55 ppb;
  - Toluene at 20 ppb;
  - o m,p-Xylene at 15 ppb;
  - o -Xylene at 11 ppb;
  - Total Xylene at 26 ppb;
  - o 1,3,5-Trimethylbenzene at 5.2 ppb; and
  - o 1,2,4-Trimethylbenzene at 7.4 ppb.

Naphthalene was detected at a concentration of 1,500 ppb, which exceeds its guidance value of 10 ppb.

Twelve (12) SVOCs were detected above their method reporting limits of which three (3) exceeded their respective NYSDEC Class GA criteria and guidance values, as follows.

- Naphthalene at 742 ppb;
- Phenanthrene at 69.5 ppb; and
- o Biphenyl at 7.84 ppb (estimated value, i.e., J detect).



- MWN-02 (screen depth: 23.6' 33.6'): Two (2) VOCs were detected above method reporting limits of which one (1) exceeded its respective NYSDEC Class GA criteria and guidance values, as follows.
  - o Benzene at 1.5 ppb.

Thirteen (13) SVOCs were detected above their method reporting limits, but all below their respective NYSDEC Class GA criteria or guidance values.

- o <u>MWN-02B (screen depth: 46.3' 56.3')</u>: Seven (7) VOCs were detected above method reporting limits of which six (6) exceeded their respective NYSDEC Class GA criteria and guidance values, as follows.
  - Benzene at 62 ppb;
  - Toluene at 10 ppb;
  - o m,p-Xylene at 7.2 ppb;
  - o o-Xylene at 10 ppb; and
  - o Total Xylene at 17.2 ppb.

Naphthalene was detected at a concentration of 320 ppb, which exceeds its guidance value of 10 ppb.

Fourteen (14) SVOCs were detected above method reporting limits, but below their respective NYSDEC Class GA criteria or guidance values, except for Naphthalene. Naphthalene was detected at a concentration of 146 ppb, which exceeds its guidance value of 10 ppb.

One (1) metal, arsenic, was detected at a concentration of 37.9 ppb, which exceeds its Class GA criteria of 25 ppb.

 MWN-02D (screen depth: 74.3' - 79.3'): One (1) metal was detected above its method reporting limits, but below its respective NYSDEC Class GA criteria.

Note: Monitoring well MWN-02D was unable to be low-flow sampled with a submersible pump and the sample was collected via a dedicated bailer. Due to potentially elevated turbidity resulting from the sampling technique, metal samples were filtered using a 0.45-micron filter by the laboratory.

- o <u>MWN-03 (screen depth: 39.2' 49.2')</u>: Seven (7) VOCs were detected above method reporting limits of which two (2) exceeded their respective NYSDEC Class GA criteria and guidance value, as follows.
  - o Benzene at 11 ppb.

Naphthalene was detected at a concentration of 25 ppb, which exceeds its guidance value of 10 ppb.

Fourteen (14) SVOCs were detected above method reporting limits, but below their respective NYSDEC Class GA criteria or guidance values, except for Naphthalene. Naphthalene was detected at a concentration of 15.0 ppb, which exceeds its guidance value of 10 ppb.



Note: Monitoring well MWN-03 was unable to be low-flow sampled with a submersible pump and the sample was collected via a dedicated bailer.

- o MWN-03B (screen depth: 60.7' 70.7'): Three (3) metals were detected above method reporting limits of which one (1) exceeded its respective NYSDEC Class GA criteria, as follows.
  - o Barium at 1,320 ppb.

Note: Monitoring well MWN-03B was unable to be low-flow sampled with a submersible pump and the sample was collected via a dedicated bailer. Due to potentially elevated turbidity resulting from the sampling technique, metal samples were filtered using a 0.45-micron filter by the laboratory.

o <u>MWN-03D</u> (screen depth: 111.3' - 121.3'): No VOCs were detected above method reporting limits. Four (4) SVOCs were detected above method reporting limits all of which were below their respective NYSDEC Class GA criteria.

Two (2) metals were detected above method reporting limits of which one (1) exceeded its respective NYSDEC Class GA criteria, as follows.

Manganese at 333 ppb.

Note: Monitoring well MWN-03D was unable to be low-flow sampled with a submersible pump and the sample was collected via a dedicated bailer. Due to potentially elevated turbidity resulting from the sampling technique, metal samples were filtered using a 0.45-micron filter by the laboratory.

o <u>MWN-04 (screen depth: 48.5' - 58.5')</u>: Two (2) VOCs were detected above method reporting limits of which only naphthalene was detected at 16.0 ppb which exceeds its guidance value of 10 ppb.

Fifteen (15) SVOCs were detected above method reporting limits of which three (3) exceeded their respective NYSDEC Class GA criteria and guidance values, as follows

- o Naphthalene at 11.2 ppb;
- o Benzo [b] Fluoranthene at 0.125 ppb (estimated value, i.e., J detect); and
- o Benzo [a] Pyrene at 0.076 ppb (estimated value, i.e., J detect).

Note: Monitoring well MWN-04 was unable to be low-flow sampled with a submersible pump and the sample was collected via a dedicated bailer. Sample results from MWN-04 were elevated compared to previous rounds, potentially due to the elevated turbidity of the samples.

In general, contaminant concentrations were consistent with historical data collected during previous sampling events completed at the Site. A more detailed discussion, including trend analysis, is provided in Section 6.00 of



this report. Bis(2-Ethylhexyl)Phthalate was detected in MWN-03D at 44.9 ug/l in 2020 and was significantly lower in 2021 (7.15 ug/l) and 2022 (0.376 J ug/L). This trend will continue to be evaluated.

#### 5.20 SEMI-ANNUAL WT-1 VICINITY MONITORING WELLS

Monitoring well locations MWN-01 and MWN-01B are included in both annual and semi-annual sampling schedules. The analytical results for these monitoring locations are discussed above in Section 5.10. Results from the remaining semi-annual wells are discussed below.

- WT1-02 (screen depth: 27.8' 37.8'): Eight (8) VOCs were detected above method reporting limits of which two (2) exceeded their respective NYSDEC Class GA criteria and guidance values, as follows.
  - o Benzene at 8.7 ppb.

Naphthalene was detected at a concentration of 27 ppb, which exceeds its guidance value of 10 ppb.

Thirteen (13) SVOCs were detected above their method reporting limits, but below their respective NYSDEC Class GA criteria or guidance values with the exception of Naphthalene. Naphthalene was detected at a concentration of 17.2 ppb, which exceeds its guidance value of 10 ppb.

- WT1-04 (screen depth: 15.5' 25.5'): Eight (8) VOCs were detected above method reporting limits of which three (3) exceeded their respective NYSDEC Class GA criteria and guidance values, as follows.
  - o Benzene at 14 ppb; and
  - o Total Xylene at 6.1 ppb.

Naphthalene was detected at a concentration of 66 ppb, which exceeds its respective guidance value of 10 ppb.

Sixteen (16) SVOCs were detected above their method reporting limits and four (4) exceeded their respective NYSDEC Class GA guidance values, as follows.

- Naphthalene at 32.6 ppb;
- o Benzo [a] Anthracene at 0.342 ppb (estimated value, i.e., J detect);
- o Benzo [b] Fluoranthene at 0.093 ppb (estimated value, i.e., J detect); and
- Chrysene at 0.250 ppb (estimated value, i.e., J detect).
- o <u>WT1-05 (screen depth: 13.3' 23.3')</u>: Eight (8) VOCs were detected above method reporting limits of which four (4) exceeded their respective NYSDEC Class GA criteria and guidance values, as follows.
  - Benzene at 9.7 ppb;
  - o m,p-Xylene at 5.4 ppb; and
  - o Total Xylene at 9.4 ppb (estimated value, i.e., J detect).



Naphthalene was detected at a concentration of 220 ppb which exceeds its guidance value of 10 ppb.

Thirteen (13) SVOCs were detected above method reporting limits, of which two (2) exceeded their respective NYSDEC Class GA guidance values as follows.

- o Naphthalene at 106 ppb; and
- o Biphenyl at 5.70 ppb.
- o <u>BCP-ORC-1</u> (screen depth: 24.7′ 34.7′): Six (6) VOCs were detected above method reporting limits of which three (3) exceeded their respective NYSDEC Class GA criteria and guidance values, as follows.
  - o Benzene at 25 ppb; and
  - o Total Xylene at 8.2 ppb (estimated value, i.e., J detect).

Naphthalene was detected at a concentration of 460 ppb, which exceeds its guidance value of 10 ppb.

Fifteen (15) SVOCs were detected above method reporting limits, of which three (3) exceeded their respective NYSDEC Class GA guidance values as follows.

- o Naphthalene at 198 ppb;
- o Benzo [a] Anthracene at 0.214 ppb (estimated value, i.e., J detect); and
- o Chrysene at 0.145 ppb (estimated value, i.e., J detect).

In general, VOC and SVOC concentrations were consistent with historical data collected during previous sampling events. A more detailed discussion, including a trend analysis, is provided in Section 6.00 of this report.

#### **6.00 STATISTICAL ANALYSIS**

As stated in Section 2.4 of Attachment A4 (LTGWM Plan) of the September 2007 *Site Management Plan*, a statistical analysis is required for all detected constituents (in groundwater) that are observed at concentrations above NYSDEC Class GA criteria or guidance values. In lieu of performing moving trend analysis, as described in the LTGWM Plan, GZA generated time series plots for parameters which exceeded the NYSDEC Class GA criteria, either during this monitoring round or in previous routine monitoring rounds (routine monitoring started in 2008). These plots were evaluated for trends over the routine monitoring period time, which started in 2008 (approximately 14 years) at a 95% confidence interval, and outliers. Sen's Test for trends were performed to evaluate statistically significant trends in the data with respect to time. Time series plots were generated on a well-by-well basis and are presented in **Appendix C**. During future monitoring rounds, the time series plots may be evaluated over the most recent five-year period, rather than the entire routine monitoring period.



Thirty-one statistically significantly decreasing trends in contaminant concentrations were identified by the Sen's Tests:

- BCP-ORC-1: benzene and biphenyl;
- MWN-01: 1,2,4-trimethylbenzene, benzene, biphenyl, fluorene, m, p-xylene, naphthalene, o-xylene, phenanthrene, toluene and xylenes;
- MWN-01B: benzene;
- MWN-02: benzene and xylenes;
- MWN-02B: benzene and toluene;
- MWN-03B: manganese;
- WT1-02: 1,3,5-trimethylbenzene, benzene, m,p-xylene, o-xylene, toluene and xylenes; and
- WT1-04: 1,2,4-trimethylbenzene, 1,3,5-trimethylbenzene, benzene, m, p-xylene, o-xylene, toluene and xylenes.

The Sen's Tests also identified two statistically significant increasing trends:

- BCP-ORC-1: o-xylene; and
- WT1-05: phenanthrene.

Time series plots were also evaluated for seasonality and outliers. There does not appear to be significant seasonal fluctuations of contaminant concentrations or outliers in the current monitoring data.

#### 7.00 SUMMARY

GZA was retained to collect and analyze groundwater samples from thirteen (13) annual/semi-annual monitoring wells at the Steel Winds I facility in accordance with the *Site Management Plan*. A summary of our findings follows.

## **Annual Well Locations**

- Static groundwater level measurements indicate that groundwater flows predominantly in a westerly direction at the Site, toward Lake Erie. Groundwater in the vicinity of WT-01 was observed to flow south-southwesterly towards Smoke Creek and Lake Erie.
- VOCs were detected at concentrations above NYSDEC Class GA criteria in the groundwater samples collected from LTGWM wells MWN-01, MWN-01B, MWN-02, MWN-02B, MWN-03 and MWN-04.
- VOCs were not detected above method reporting limits (and therefore not above NYSDEC Class GA criteria) in LTGWM monitoring well location MWN-03D.



- SVOCs were detected at concentrations above NYSDEC Class GA or their respective guidance criteria in the groundwater samples collected from LTGWM wells MWN-01, MWN-01B, MWN-02B, MWN-03, and MWN-04.
- Arsenic was detected at concentrations above NYSDEC Class GA criteria in LTGWM well MWN-02B.
- Barium was detected at concentrations above NYSDEC Class GA criteria in LTGWM well MWN-03B.
- Manganese detected at concentrations above NYSDEC Class GA criteria in LTGWM well MWN-03D.

# <u>Semi-Annual Well Locations</u>

- VOCs were detected at concentrations above NYSDEC Class GA criteria in the groundwater samples collected from the semi-annual WT1 vicinity wells MWN-01, MWN-01B, WT1-02, WT1-04, WT1-05 and BCP-ORC-1.
- SVOCs were detected at concentrations above NYSDEC Class GA or their respective guidance criteria in the groundwater samples collected from the semi-annual WT1 vicinity wells MWN-01, MWN-01B, WT1-02, WT1-04, WT1-05, and BCP-ORC-1.
- Based on our review of the historic and current analytical data, the analytical test results from the September 2022 round of sampling are generally consistent with historical data. Statistically significant trends in contaminant concentrations were identified as noted in Section 6.00.



**TABLES** 

# TABLE 1

# September 2022 Analytical Testing Program Summary Steel Winds I Facility Lackawanna, New York

Well Designation	Sample ID	Date Collected	Screened Interval (TOR)	STARS VOCs	SVOCs (BN)	Total Arsenic	Total Barium	Total Chromium	Total Manganese
<b>Annual Monitoring</b>	g Well Sample Location	s (LTGWM Netwo	ork)						
MWN-01	MWN-01-091322	9/13/2022	9.2 - 19.2	Х	X				
MWN-01B	MWN-01B-091322	9/13/2022	22.2 - 32.2	X	X				
MWN-02	MWN-02-091422	9/14/2022	23.6 - 33.6	Х	X				
MWN-02B	MWN-02B-091422	9/14/2022	46.3 - 56.3	X	X	X			
MWN-02D	MWN-02D-091522	9/15/2022	74.3 - 79.3			Х	Х	X	
MWN-03	MWN-03-091522	9/15/2022	39.2 - 49.2	Х	Х				
MWN-03B	MWN-03B-091522	9/15/2022	60.7 - 70.7			Х	Х	Х	Χ
MWN-03D	MWN-03D-091522	9/15/2022	111.3 - 121.3	Х	X		Х		Χ
MWN-04	MWN-04-091522	9/15/2022	48.5 - 58.5	X	X				
Semi-Annual Mon	itoring Well Sample Lo	cations (WT-1 Vi	cinity Network)						
MWN-01	MWN-01-091322	9/13/2022	9.2 - 19.2	X	X				
MWN-01B	MWN-01B-091322	9/13/2022	22.2 - 32.2	X	X				
WT1-02	WT1-02-091322	9/13/2022	27.8 - 37.8	X	X				
WT1-04	WT1-04-091322	9/13/2022	15.5 - 25.5	Х	Х				
WT1-05	WT1-05-091322	9/13/2022	13.3 - 23.3	Х	Х				
BCP-ORC-1	BCP-ORC-1-091322	9/13/2022	24.7 - 34.7	Х	Х				

- 1. VOCs = Volatile Organic Compounds STARS list via EPA Method 8260C.
- 2. SVOCs = Semi-Volatile Organic Compounds Base-Neutrals list via EPA Method 8270D.
- 3. Arsenic, Barium, Chromium, and Manganese via EPA Method 6010D.
- 4. "WT", "MWN", and "BCP-ORC" monitoring well information provided in Table 1 was referenced from Turnkey Environmental Restoration, LLC's 2009 Annual LTGWM & First Semi-Annual WT-1 Vicinity Monitoring Report.
- 5. TOR = measurement recorded in feet below top-of-well riser.

# Table 2

# September 2022 Annual Groundwater Analytical Data Summary Steel Winds I Facility Lackawanna, New York

	NYSDEC			MWN-01					MWN-01B					MWN-02		
Parameter	Class GA	9/17/2020	4/2/2021	9/2/2021	3/30/2022	9/13/2022	9/17/2020	4/2/2021	9/2/2021	3/30/2022	9/13/2022	9/18/2018	9/25/2019	9/17/2020	9/3/2021	9/14/2022
2 1111111111111111111111111111111111111	Criteria	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result
Water Quality Field Measurements																
pH (units)	6.5 - 8.5	7.81	7.66	11.53	13.19	11.81	7.83	8.01	11.1	13.03	11.46	12.28	11.94	8.31	11.7	11.85
Temperature (°C)	NV	14.4	10.5	10.8	9.6	12.0	10.9	11	9.8	9.4	10.6	12.6	11.3	12.35	12.6	12.6
Specific Conductance (mS/cm)	NV	1.450	1.380	1.212	1.170	1.258	0.991	1.010	0.831	0.808	0.891	1.886	1.763	2.04	1.776	1.965
Turbidity (NTU)	5	2.9	2.4	2.61	1.08	2.80	7.3	5.4	7.67	22.3	22.18	3.0	38.6	6.8	2.51	2.54
Dissolved Oxygen (mg/L)	NV	116.7	132.3	1.2	2.2	5.9	134.7	115.9	0.8	20.7	11.3	1.51	0.060	97.2	2.8	13.6
Oxygen Reduction Potential (mV)	NV	-237	-231	-159.2	-347.1	-104.5	-247	-204	-214.2	-244.3	-118.8	-87.2	-121.0	-281	-115.1	137.8
<b>Volatile Organic Compounds - EPA</b>	Method 8260C	(ug/L)														
Benzene	1	17	14	14	14	12	59	57	55	54	55	2.5	2.2	1	5.1	1.5
Toluene	5	4.2	4.0 J	3.6 J	3.1 J	2.8 J	18 J	20 J	19 J	16 J	20	<	<	<	1.4 J	<
Ethylbenzene	5	0.98 J	<	<	<	<	<	<	<	<	0.95 J	<	<	<	<	<
m,p-Xylene	5	10	9.3	8.7	7.9	6.0	13 J	15 J	12 J	12 J	15	1.3 J	1.1 J	0.76 J	2.4 J	<
o-Xylene	5	8	7.1	6.5	5.8	5.0	9.1 J	10 J	9.0 J	8.9 J	11	1.2 J	1.1 J	<	2.1 J	<
Xylene (Total)	5	18.0	16.0	15.2	14	11.0	22.1	25 J	21 J	21 J	26	2.5	2.2	0.76 J	4.5 J	<
Isopropylbenzene	5	<	<	<	<	<	<	<	<	<	1.4 J	<	<	<	<	<
1,3,5-Trimethylbenzene	5	4	4.5 J	4.2 J	3.9 J	2.8 J	<	<	<	<	5.2	1.8 J	1.4 J	0.91 J	1.8 J	<
1,2,4-Trimethylbenzene	5	4.3	4.8 J	4.6 J	4.1 J	3.0 J	<	7.6 J	7.1 J	<	7.4	<	<	<	1.2 J	<
Naphthalene*	10	240	310	270	290	240	1,500	1,800	1,500	1,700	1,500	7.3	9.4	20	20	4.2
Semi-Volatile Organic Compounds -	EPA Method 8	270D (ug/L)														
Acetophenone	NV	<	<	<	<	0.570 J	<	<	<	<	<	0.368 J	<	<	<	0.246 J
Acenaphthylene	NV	22.4	34	22.3	30.3	23.5	54.6	44	44.0	33.8	54.3	0.815	1.36	0.727	1.98	1.03
Naphthalene*	10	139.0	140	96.2	141	91.9	1,030	910	962	970	742	1.19	2.87	2.38	5.23	3.44
2-Methylnaphthalene	NV	27.1	35	21.9	40.0	27.8	48.0	41	35.8	46.2	52.4	0.302 J	1.02	0.552	1.78	1.01
Acenaphthene*	20	8.34	13	8.66	11.9	10.1	11.2	10	12.0	10.5	11.8	0.538	0.758	0.431 J	1.20	0.603
Dibenzofuran	NV	25.9	44	28.9	39.6	29.7	29.4	23	30.3	24.8	30.6	0.156J	0.922	0.584	2.35	0.967
Fluorene*	50	38.3	70	41.9	58.8	44.4	43.9	38	43.7	35.7	42.3	0.92	2.98	1.52	4.76	2.26
Phenanthrene*	50	45.3	110	71.0	81.5	69.9	64.3	55	61.9	53.6	69.5	<	2.55	1.46	4.14	1.76
Dibenzo (a,h)Anthracene	NV	<	0.05 J	<	<	<	<	<	<	<	<	<	<	<	<	<
Carbazole	NV	20.30	26	19.6	24.1	19.7	62.4	52	60.0	55.4	61.3	0.598	1.34	0.702	3.67	1.28
Anthracene*	50	5.81	19	7.74	11.9	12.2	11.00	5.3	8.19	6.46	11.8	0.212J	0.635	0.467 J	0.983	0.588
Fluoranthene*	50	5.72	24	9.44	10.6	12.3	10.3	10	8.97	8.33	10.8	1.86	1.4	1.14	1.56	0.971
Biphenyl	5	6.41	8.8	5.85	7.86	6.48	8.19	6.5	7.45	6.09	7.84 J	0.199 J	0.412 J	0.198 J	0.732	0.332 J
Pyrene*	50	4.47	14	6.16	6.38	6.81	6.62	5.9	6.44	4.95	5.57 J	1.28	1.26	1.41	1.56	1.70
Butyl benzyl phthalate*	50	<	<	0.104 J	<	<	<	<	<	<	<	<	<	<	0.093 J	<
Benzo [a] Anthracene	0.002	<	1.4	<	0.372 J	0.380 J	<	0.38	0.461 J	0.316 J	<	<	<	<	<	<
Benzo [b] Fluoranthene*	0.002	<	0.4	<	<	0.079 J	<	0.10 J	0.105 J	0.105 J	<	<	<	<	<	<
Benzo [k] Fluoranthene*	0.002	<	0.14	<	<	<	<	0.04 J	< .	<	<	<	<	<	<	<
Benzo [a] Pyrene	ND 0.002	<	0.26	<	<	<	<	0.05 J	0.072 J	0.079 J	<	<	<	<	<	<
Indeno [1,2,3-cd] Pyrene*	0.002	<	0.11	<	<	<	<	0.04 J	<	<	<	<	<	<	<	<
Benzo (g,h,i) Perylene	NV	<	0.09 J	0.216 I	< 0.197 T	< 0.214 T	<	0.03 J	< 0.256 I	( ) 190 T	<	<	<	<	<	<
Chrysene bis(2-Ethylhexyl)Phthalate	0.002 5	< 0.456 J	0.82	0.216 J	0.187 J	0.214 J	1.10	0.22	0.256 J	0.180 J	< <	<	0.098 J	0.602	< <	<
Metals - EPA Method 6010D (ug/L)	<u>J</u>	0.430 J		<	<	<	1.10	<	<	<			0.036 J	0.002		<
. 6	25	NIT	NIT	NIT	NIT	NIT	NIT	NIT	NIT	NIT	NTT	NIT	NIT	NIT	NIT	NIT
Arsenic	25	NT NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
Barium	1,000	NT NT	NT	NT	NT NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT NT	NT
Chromium	50	NT NT	NT NT	NT NT	NT NT	NT NT	NT NT	NT NT	NT	NT	NT NT	NT	NT	NT NT	NT NT	NT
Manganese	300	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT

- 1. Compounds detected in one or more sample are presented on this table. Refer to Appendix B for list of all compounds included in analysis.
- 2. Analytical testing completed by Alpha Analytical, Westborough, Massachusetts.
- 3. NYSDEC Groundwater Class GA criteria obtained from Division of Water Technical and Operational Guidance Series (TOGS 1.1.1),

dated October 1993, revised June 1998, errata January 1999 and amended April 2000 (Class GA).

- 5. < indicates compound was not detected above method detection limits.</li>
  6. "J" qualifier = Analyte detected below quantitation limits.
- 7. Value shown in **bold** indicate exceedance of respective Class GA Criteria or guidance value.
- 8. NV = no value, NT = not tested, ND = Not detected above method detection limit
- 9. \* = value shown is a guidance value rather than a groundwater standard.
- 10. The equipment used to collect water quality data was calibrated prior to and during use in accordance with the manufacturer's recommendations.
- 11. DO and pH measurements are routinely made using the same model water quality meter, however the measurements made on 9/2020 and 4/2021 appear erroneous.

<sup>4.</sup> ug/L = part per billion (ppb).

Table 2

# September 2022 Annual Groundwater Analytical Data Summary Steel Winds I Facility Lackawanna, New York

	NYSDEC			MWN-02B					MWN-02D					MWN-03		
Parameter	Class GA	9/17/2018	9/25/2019	9/17/2020	9/3/2021	9/14/2022	9/18/2018	9/24/2019	9/18/2020	9/3/2021	9/15/2022	9/18/2018	9/25/2019	9/17/2020	9/2/2021	9/15/2022
	Criteria	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result
<b>Water Quality Field Measurements</b>	<u> </u>					•										
pH (units)	6.5 - 8.5	11.57	11.34	8.21	11.30	11.17	7.10	7.00	6.99	6.61	7.86	12.39	12.32	8.53	12.00	12.49
Temperature (°C)	NV	13.1	12.1	12.92	12.6	13.9	13.7	12.6	13.61	12.9	13.8	13.5	12.8	13.57	13.3	14.3
Specific Conductance (mS/cm)	NV	0.942	0.958	1.13	0.910	0.902	1.864	1.890	1.970	1.354	2.027	2.825	2.724	2.89	2.729	3.058
Turbidity (NTU)	5	2.8	1.9	6.9	2.52	2.57	3.8	15.1	7.2	5.15	189.3	3.1	3.9	3.9	4.82	4.06
Dissolved Oxygen (mg/L)	NV	0.14	0.15	95.5	1.2	6.5	0.15	0.09	6.1	1.5	29.9	0.17	0.11	115.2	2.1	43.0
Oxygen Reduction Potential (mV)	NV	-284.9	-220.6	-256	-202.6	-56.1	-70.5	-96.3	-72	-51.6	49.8	-352.6	-412.1	-361	-267.3	-39.3
Volatile Organic Compounds - EPA	Method 8260C	(ug/L)														
Benzene	1	71	64	69	61	62	NT	NT	NT	NT	NT	12	8.0	10	7.1	11
Toluene	5	12	11	11	11	10	NT	NT	NT	NT	NT	2.7	2.0 J	2.2 J	1.8 J	2.4 J
Ethylbenzene	5	<	0.76 J	<	<	<	NT	NT	NT	NT	NT	<	<	<	<	<
m,p-Xylene	5	8.2	8.2	8.5	9.2	7.2	NT	NT	NT	NT	NT	2.4 J	1.4 J	1.5 J	1.3 J	1.6 J
o-Xylene	5	12	12	13.0	13	10	NT	NT	NT	NT	NT	2.7	1.5 J	1.8 J	1.4 J	1.7 J
Xylene (Total)	5	20.2	20.2	21.5	22.2	17.2	NT	NT	NT	NT	NT	5.1	2.9	3.3	2.7 J	3.3 J
1,3,5-Trimethylbenzene	5	<	1.5 J	1.5 J	2.0 J	<	NT	NT	NT	NT	NT	1.4 J	0.90 J	0.97 J	0.93 J	0.97 J
1,2,4-Trimethylbenzene	5	2.8 J	2.5	2.6 J	3.5 J	1.9 J	NT	NT	NT	NT	NT	<	<	<	<	<
Naphthalene*	10	340	240	270	280	320	NT	NT	NT	NT	NT	20	23	26	19	25
<b>Semi-Volatile Organic Compounds</b>	- EPA Method 8	3270D (ug/L)														
Acetophenone	NV	0.451 J	<	<	<	0.770 J	NT	NT	NT	NT	NT	0.271 J	<	<	<	0.308 J
Acenaphthylene	NV	3.77	4.58	3.90	3.18	2.83	NT	NT	NT	NT	NT	1.78	1.73	0.980	1.23	2.70
1,2-Dichlorobenzene	3	0.178J	0.171 J	0.168 J	0.162 J	0.200 J	NT	NT	NT	NT	NT	0.104 J	0.099 J	0.121 J	0.102 J	0.115 J
Naphthalene*	10	150	217	205	183	146	NT	NT	NT	NT	NT	16.5	17.1	18.1	11.2	15.0
2-Methylnaphthalene	NV	9.6	8.05	8.83	6.89	8.48	NT	NT	NT	NT	NT	3.26	2.7	3.10	1.93	3.03
Acenaphthene*	20	7.39	7.09	7.47	7.46	6.20	NT	NT	NT	NT	NT	1.63	1.3	1.45	1.11	1.54
Dibenzofuran	NV	5.58	5.76	6.24	6.32	4.50	NT	NT	NT	NT	NT	3.12	2.34	2.81	1.99	2.92
Fluorene*	50	8.44	10.7	11.40	10.2	7.72	NT	NT	NT	NT	NT	5.47	4.5	4.82	3.48	5.10
Phenanthrene*	50	15.4	17.5	18.30	18.0	13.7	NT	NT	NT	NT	NT	8.69	8.23	8.29	7.54	9.37
Carbazole	NV	21.9	23.2	24.40	23.1	21.2	NT	NT	NT	NT	NT	4.88	4.30	4.58	3.26	5.17
Anthracene*	50	2.33	2.32	2.35	1.67	1.88	NT	NT	NT	NT	NT	0.86	1.00	0.612	0.884	1.38
Fluoranthene*	50	3.96	3.32	4.13	3.34	3.51	NT	NT	NT	NT	NT	2.48	2.7	2.53	2.18	3.19
Biphenyl	5	1.41	1.64	1.62	1.52	1.11	NT	NT	NT	NT	NT	0.868	0.707	0.792	0.512	0.715
Pyrene*	50	2.5	2.22	2.82	2.49	2.00	NT	NT	NT	NT	NT	1.76	1.66	1.63	1.78	1.91
Butylbenzylphthalate*	50	<	<	<	0.124 J	<	NT	NT	NT	NT	NT	<	<	<	<	<
bis(2-Ethylhexyl)Phthalate	5	<	<	<	<	<	NT	NT	NT	NT	NT	<	<	0.336 J	<	<
Benzo [b] Fluoranthene	0.002	<	<	<	<	<	NT	NT	NT	NT	NT	0.171J	<	<	<	<
Metals - EPA Method 6010D (ug/L)																
Arsenic	25	30.81	32	28.44	27.68	37.9	0.72	0.60	0.63	0.62	<	NT	NT	NT	NT	NT
Barium	1,000	NT	NT	NT	NT	NT	957	931.9	912.8	922.5	860	NT	NT	NT	NT	NT
Chromium	50	NT	NT	NT	NT	NT	0.19 J	<	0.30 J	0.60 J	<	NT	NT	NT	NT	NT
Manganese	300	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT

# Notes:

- 1. Compounds detected in one or more sample are presented on this table. Refer to Appendix B for list of all compounds included in analysis.
- 2. Analytical testing completed by Alpha Analytical, Westborough, Massachusetts.
- 3. NYSDEC Groundwater Class GA criteria obtained from Division of Water Technical and Operational Guidance Series (TOGS 1.1.1),

dated October 1993, revised June 1998, errata January 1999 and amended April 2000 (Class GA).

- 4. ug/L = part per billion (ppb).
- 5. < indicates compound was not detected above method detection limits.
- 6. "J" qualifier = Analyte detected below quantitation limits.
- 7. Value shown in **bold** indicate exceedance of respective Class GA Criteria or guidance value.
- 8. NV = no value, NT = not tested, ND = Not detected above method detection limit
- 9. \* = value shown is a guidance value rather than a groundwater standard.
- 10. The equipment used to collect water quality data was calibrated prior to and during use in accordance with the manufacturer's recommendations.
- 11. DO and pH measurements are routinely made using the same model water quality meter, however the measurements made on 9/2020 and 4/2021 appear erroneous.
- 12. WellS MWN-02D and MWN-03 were unable to be low flow sampled. Hand bailing techniques were required. Metals analysis required laboratory filtration.

# Table 2

# September 2022 Annual Groundwater Analytical Data Summary Steel Winds I Facility Lackawanna, New York

	NYSDEC			MWN-3B					MWN-03D					MWN-04		
Parameter	Class GA	9/18/2018	9/25/2019	10/1/2020	9/3/2021	9/15/2022	9/18/2018	9/25/2019	9/24/2020	9/3/2021	9/15/2022	9/18/2018	9/25/2019	9/17/2020	9/2/2021	9/15/2022
	Criteria	Result	Result	Result	Result	Result	Result	Result	Result	Result <sup>11</sup>	Result	Result	Result	Result	Result	Result
Water Quality Field Measurements																
pH (units)	6.5 - 8.5	7.21	7.80	7.2	7.29	6.62	6.52	6.17	6.25	7.31	7.26	11.71	12.05	7.98	11.57	11.35
Temperature (°C)	NV	14.3	13.7	13.9	14.7	14.2	13.5	12.9	14.4	13.5	13.5	16.7	16.0	15.97	15.7	17.3
Specific Conductance (mS/cm)	NV	3.126	3.139	2.413	2.586	27.710	26.69	24.662	25.881	24.410	26.110	2.490	2.311	2.35	2.313	3.540
Turbidity (NTU)	5	6.2	25.6	38.04	16.44	40.12	4.4	29.4	14.31	35.83	165.2	2.2	2.6	2.4	1.98	33.47
Dissolved Oxygen (mg/L)	NV	0.17	0.15	49.7	2.9	25.3	0.12	0.56	36.5	5.5	16.2	4.11	5.56	107.4	3.0	69.6
Oxygen Reduction Potential (mV)	NV	-188.8	-188.8	-63.7	-146.7	97.7	-32.0	-32.4	-45.3	41.6	50.8	-101.3	-99.7	-65	-81.2	35.4
<b>Volatile Organic Compounds - EPA</b>	Method 8260C	(ug/L)														
Benzene	1	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	<	<	0.51
Toluene	5	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	<	<	<
Ethylbenzene	5	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	<	<	<
m,p-Xylene	5	NT	NT	NT	NT	NT	1.1 J	<	<	<	<	<	<	<	<	<
o-Xylene	5	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	<	<	<
Xylene (Total)	5	NT	NT	NT	NT	NT	1.1	<	<	<	<	<	<	<	<	<
1,3,5-Trimethylbenzene	5	NT	NT	NT	NT	NT	<	<	0.73 J	<	<	<	<	<	<	<
1,2,4-Trimethylbenzene	5	NT	NT	NT	NT	NT	0.74 J	<	<	<	<	<	<	<	<	<
Naphthalene*	10	NT	NT	NT	NT	NT	<	<	<	<	<	1.0 J	<	1.4 J	<	16
<b>Semi-Volatile Organic Compounds</b>	- EPA Method 8	270D (ug/L)														
Acetophenone	NV	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	<	<	0.967 J
Acenaphthylene	NV	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	<	<	0.167 J
Naphthalene*	10	NT	NT	NT	NT	NT	0.420 J	0.196 J	<	0.121 J	<	0.090 J	<	0.163 J	<	11.2
2-Methylnaphthalene	NV	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	<	<	2.49
Acenaphthene*	20	NT	NT	NT	NT	NT	<	<	<	<	0.536	<	<	0.377 J	<	5.26
Dibenzofuran	NV	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	0.107 J	<	2.54
Fluorene*	50	NT	NT	NT	NT	NT	<	<	<	<	0.187 J	<	<	0.304 J	<	4.37
Phenanthrene*	50	NT	NT	NT	NT	NT	<	<	<	<	0.434 J	0.212 J	<	0.302 J	<	7.31
Carbazole	NV	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	<	<	8.59
Anthracene*	50	NT	NT	NT	NT	NT	<	<	<	<	<	0.156 J	<	<	<	1.39
Fluoranthene*	50	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	0.168 J	<	1.55
Biphenyl	5	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	<	<	0.394 J
Pyrene*	50	NT	NT	NT	NT	NT	<	<	<	<	<	0.536	0.640	0.447 J	0.459 J	1.90
Benzo [b] Fluoranthene*	0.002	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	<	<	0.125 J
Benzo [a] Pyrene*	0.002	NT	NT	NT	NT	NT	<	<	<	<	<	<	<	<	<	0.076 J
Di-n-octylphthalate*	50	NT	NT	NT	NT	NT	<	<	0.690 J	<	<	<	<	<	<	<
Butylbenzylphthalate*	50	NT	NT	NT	NT	NT	<	0.211 J	0.091 J	0.137 J	<	<	<	<	<	<
Diethylphthalate*	50	NT	NT	NT	NT	NT	<	<	0.518	0.549	<	<	<	<	<	<
bis(2-Ethylhexyl)Phthalate	5	NT	NT	NT	NT	NT	0.232 J	0.514	44.9	7.15	0.376 J	0.083J	0.123 J	0.342 J	<	<
Metals - EPA Method 6010D (ug/L)		I	1	I		ı			·							
Arsenic	25	38	36.12	2.73	86.97	<	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
Barium	1,000	1,348	1,291	837.3	1,049	1,320	1,404	1,286	1,234	1,318	779	NT	NT	NT	NT	NT
Chromium	50	1.89	1.74	0.28 J	5.10	3.2 J	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
Manganese	300	471.8	267.4	336.7	400.2	178	245.6	38.19	41.49	24.52	333	NT	NT	NT	NT	NT

- 1. Compounds detected in one or more sample are presented on this table. Refer to Appendix B for list of all compounds included in analysis.
- 2. Analytical testing completed by Alpha Analytical, Westborough, Massachusetts.
- 3. NYSDEC Groundwater Class GA criteria obtained from Division of Water Technical and Operational Guidance Series (TOGS 1.1.1),
- dated October 1993, revised June 1998, errata January 1999 and amended April 2000 (Class GA).
- 4. ug/L = part per billion (ppb).
- 5. < indicates compound was not detected above method detection limits.
- 6. "J" qualifier = Analyte detected below quantitation limits.
- 7. Value shown in **bold** indicate exceedance of respective Class GA Criteria or guidance value.
- 8. NV = no value, NT = not tested, ND = Not detected above method detection limit
- 9. \* = value shown is a guidance value rather than a groundwater standard.
- 10. The equipment used to collect water quality data was calibrated prior to and during use in accordance with the manufacturer's recommendations.
- 11. DO and pH measurements are routinely made using the same model water quality meter, however the measurements made on 9/2020 and 4/2021 appear erroneous.
- 12. Wells MWN-03B, MWN-03D and MWN-04 were unable to be low flow sampled. Hand bailing techniques were required. Metals analysis required laboratory filtration.

# Table 3

# September 2022 Semi-Annual Groundwater Analytical Data Summary Steel Winds I Facility Lackawanna, New York

	NYSDEC			MWN-01					MWN-01B					WT1-02		
Parameter	Class GA	9/17/2020	4/2/2021	9/2/2021	3/30/2022	9/13/2022	9/17/2020	4/2/2021	9/2/2021	3/30/2022	9/13/2022	9/18/2020	4/2/2021	9/2/2021	3/30/2022	9/14/2022
	Criteria	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result
Water Quality Field Measurements																
pH (units)	6.5 - 8.5	7.81	7.66	11.53	13.19	11.81	7.83	8.01	11.1	13.03	11.46	9.07	7.68	11.85	13.45	11.98
Temperature (°C)	NV	14.4	10.5	10.8	9.6	12.0	10.9	11.0	9.8	9.4	10.6	13.08	11.5	12.3	11.6	13.1
Specific Conductance (mS/cm)	NV	1.450	1.380	1.212	1.170	1.258	0.991	1.01	0.831	0.808	0.891	2.090	1.84	1.770	1.746	1.592
Turbidity (NTU)	5	2.9	2.4	2.61	1.08	2.80	7.30	5.4	7.67	22.3	22.18	16	8.6	2.7	1.37	1.43
Dissolved Oxygen (mg/L)	NV	116.7	132.3	1.2	2.2	5.9	134.7	115.9	0.8	20.7	11.3	28.3	33.6	4.7	3.9	7.6
Oxygen Reduction Potential (mV)	NV	-237	-231	-159.2	-347.1	-104.5	-247	-204	-214.2	-244.3	-118.8	-200	-177	-160.7	-271.7	-41.2
Volatile Organic Compounds - EPA M	Tethod 8260C (ug/L)									•						
Benzene	1	17	14	14	14	12	59	57	55	54	55	7.6	6.0	12	11.0	8.7
Toluene	5	4.2	4.0 J	3.6 J	3.1 J	2.8 J	18 J	20 J	19 J	16 J	20	1.6 J	1.3 J	2.4 J	2.1 J	1.7 J
Ethylbenzene	5	0.98 J	<	<	<	<	<	<	<	<	0.95 J	<	<	<	<	<
m,p-Xylene	5	10	9.3	8.7	7.9	6.0	13 J	15 J	12 J	12 J	15	3.1	2.2 J	4.2	4	2.6
o-Xylene	5	8	7.1	6.5	5.8	5.0	9.1 J	10 J	9.0 J	8.9 J	11	2.6	1.6 J	3.0	2.9	1.9 J
Xylene (Total)	5	18.0	16	15.2	14	11.0	22.1	25 J	21 J	21 J	26	5.7	3.8 J	7.2	6.9	4.5 J
Isopropylbenzene	5	<	<	<	<	<	<	<	<	<	1.4 J	<	<	<	<	<
1,3,5-Trimethylbenzene	5	4	4.5 J	4.2 J	3.9 J	2.8 J	<	<	<	<	5.2	1.5 J	1.3 J	2.0 J	2.0 J	1.2 J
1,2,4-Trimethylbenzene	5	4.3	4.8 J	4.6 J	4.1 J	3.0 J	<	7.6 J	7.1 J	<	7.4	1.1 J	0.86 J	1.5 J	1.5 J	0.84 J
Naphthalene*	10	240	310	270	290	240	1,500	1,800	1,500	1,700	1,500	36	22	43	45	27
Semi-Volatile Organic Compounds - I	EPA Method 8270D (ug	;/L)														
Acetophenone	NV	<	<	<	<	0.570 J	<	<	<	<	<	<	<	<	<	0.317 J
Acenaphthylene	NV	22.4	34	22.3	30.3	23.5	54.6	44	44.0	33.8	54.3	1.08	1.1	0.651	1.30	1.16
Naphthalene*	10	139.0	140	96.2	141	91.9	1,030	910	962	970	742	16.90	14	9.38	16.8	17.2
2-Methylnaphthalene	NV	27.1	35	21.9	40.0	27.8	48.0	41	35.8	46.2	52.4	3.57	3.3	2.11	4.05	4.62
Acenaphthene*	20	8.34	13	8.66	11.9	10.1	11.2	10	12.0	10.5	11.8	1.08	1.3	0.710	1.51	1.47
Dibenzofuran	NV	25.9	44	28.9	39.6	29.7	29.4	23.0	30.3	24.8	30.6	3.94	3.7	2.47	4.92	4.92
Fluorene*	50	38.30	70	41.9	58.8	44.4	43.9	38	43.7	35.7	42.3	6.14	7.3	3.50	7.51	7.48
Phenanthrene*	50	45.30	110	71.0	81.5	69.9	64.3	55	61.9	53.6	69.5	13.30	17	8.10	14.1	13.7
Dibenzo (a,h)Anthracene	NV	<	0.05 J	<	<	<	<	<	<	<	<	<	<	<	<	<
Carbazole	NV	20.30	26	19.6	24.1	19.7	62.4	52.0	60.0	55.4	61.3	3.75	2.9	2.88	4.80	6.02
Anthracene*	50	5.81	19	7.74	11.9	12.2	11.00	5.3	8.19	6.46	11.8	2.40	2.8	1.44	2.52	2.74
Fluoranthene*	50	5.72	24	9.44	10.6	12.3	10.3	10	8.97	8.33	10.8	4.02	6.6	3.18	5.42	4.61
Biphenyl	5	6.41	8.8	5.85	7.86	6.48	8.19	6.50	7.45	6.09	7.84 J	0.99	0.82 J	0.548	1.02	1.13
Pyrene*	50	4.47	14	6.16	6.38	6.81	6.62	5.90	6.44	4.95	5.57 J	3.45	4.8	2.39	3.57	2.93
Butyl benzyl phthalate*	50	<	<	0.104 J	<	<	<	<	<	<	<	<	<	<	<	<
Benz [a] Anthracene*	0.002	<	1.4	<	0.372 J	0.380 J	<	0.38	0.461 J	0.316 J	<	<	0.24	<	0.202 J	<
Benzo [b] Fluoranthene*	0.002	<	0.40	<	<	0.079 J	<	0.10 J	0.105 J	0.105 J	<	<	0.03 J	<	<	<
Benzo [k] Fluoranthene*	0.002	<	0.14	<	<	<	<	0.04 J	<	<	<	<	0.01 J	<	<	<
Benzo [a] Pyrene	ND	<	0.26	<	<	<	<	0.05 J	0.072 J	0.079 J	<	<	<	<	<	<
Indeno [1,2,3-cd] Pyrene*	0.002	<	0.11	<	<	<	<	0.04 J	<	<	<	<	<	<	<	<
Benzo (g,h,i) Perylene	NV	<	0.09 J	<	<	<	<	0.03 J	<	<	<	<	<	<	<	<
Chrysene*	0.002	<	0.82	0.216 J	0.187 J	0.214 J	<	0.22	0.256 J	0.180 J	<	<	0.17	<	0.146 J	<
bis(2-Ethylhexyl)phthalate	5	0.456 J	<	<	<	<	1.10	<	<	<	<	0.334 JB	<	<	<	<

- 1. Compounds detected in one or more sample for the past five sampling events are presented on this table. Refer to Appendix B for list of all compounds included in analysis.
- 2. Analytical testing completed by Alpha Analytical in Westborough, MA.
- 3. NYSDEC Groundwater Class GA criteria obtained from Division of Water Technical and Operational Guidance Series (TOGS 1.1.1), dated October 1993, revised June 1998, errata January 1999 and amended April 2000 (Class GA).
- 4. ug/L = part per billion (ppb).
- 5. < indicates compound was not detected above method detection limits.
- 6. "J" qualifier = Analyte detected below quantitation limits.
- 7. "B" qualifier = indicates compound was detected in the method blank sample.
- 8. "D" qualifier = indicates the compound concentration was obtained from a secondary dilution analysis.
- 7. Value shown in **bold** indicates exceedance of respective Class GA Criteria or guidance value.
- 8. NV = no value, NT = not tested, ND = Not detected above method detection limit
- 9. \* = value shown is a guidance value rather than a groundwater standard.
- 10. The equipment used to collect water quality data was calibrated prior to and during use in accordance with the manufacturer's recommendations.
- 11. DO and pH measurements are routinely made using the same model water quality meter, however the measurements made on 9/2020 and 4/2021 appear erroneous.

Table 3

# September 2021 Semi-Annual Groundwater Analytical Data Summary Steel Winds I Facility

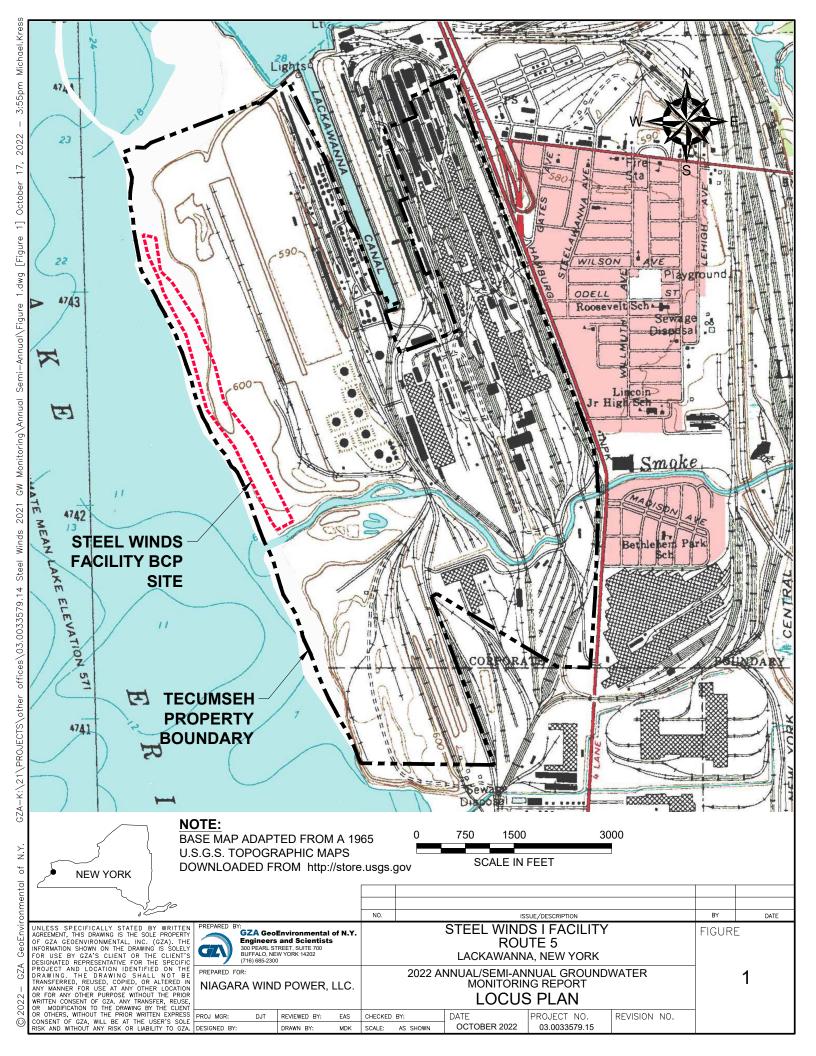
# Lackawanna, New York

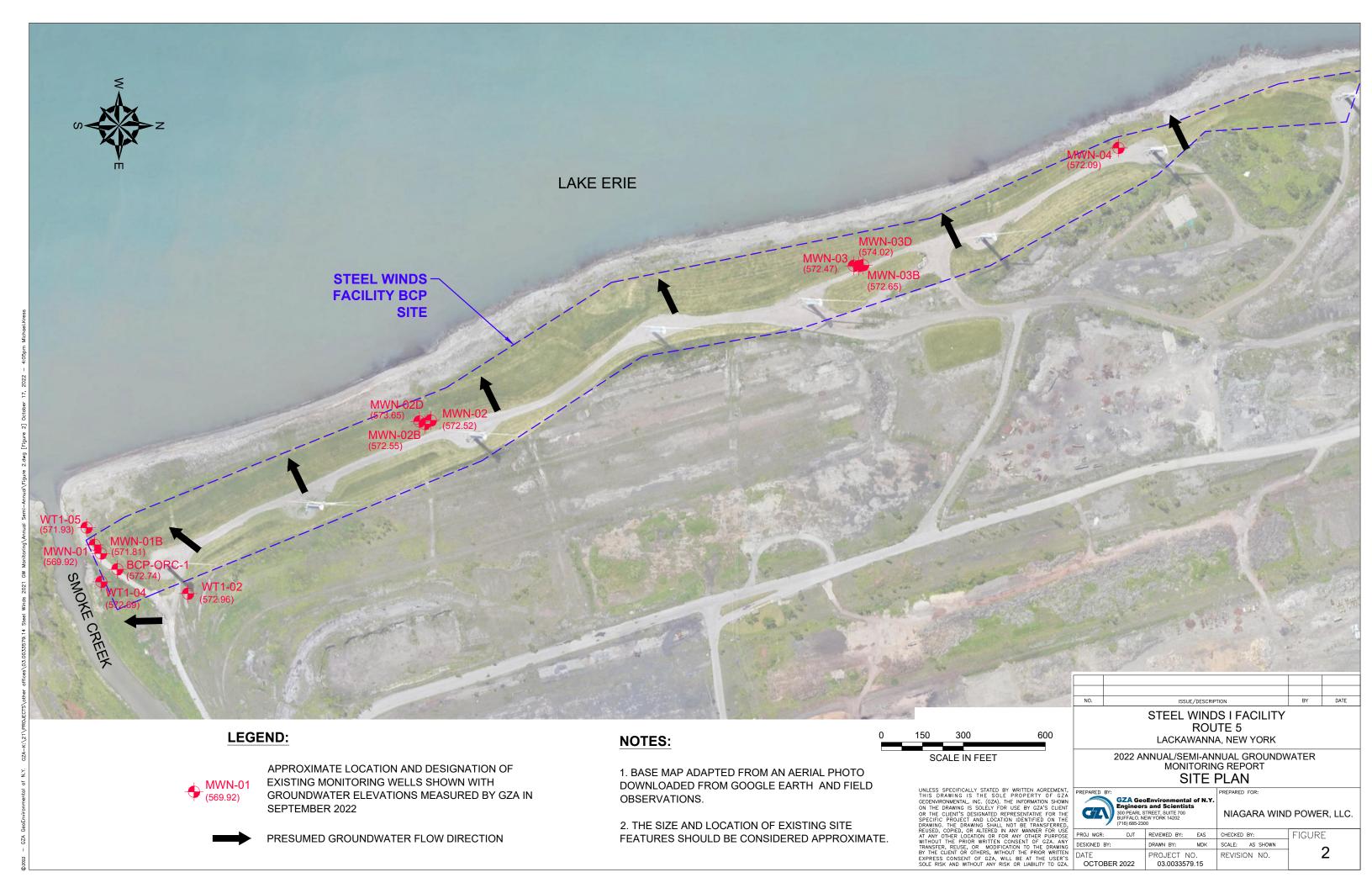
	NYSDEC WT1-04						WT1-05					BCP-ORC-1				
Parameter	Class GA	9/18/2020	4/2/2021	9/2/2021	3/30/2022	9/13/2022	9/18/2020	4/2/2021	9/2/2021	3/30/2022	9/13/2022	9/18/2020	4/2/2021	9/2/2021	3/30/2022	9/13/2022
	Criteria	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result	Result
Water Quality Field Measurements																
pH (units)	6.5 - 8.5	8.28	8.27	11.51	13.81	11.75	8.6	7.83	11.46	12.99	11.61	8.64	7.85	11.21	13.47	11.6
Temperature (°C)	NV	13.27	10.8	11.1	8.4	13.3	12.57	10.9	11.2	9.2	13.0	12.02	11.1	10.0	9.0	11.5
Specific Conductance (mS/cm)	NV	1.410	1.550	1.326	1.294	1.326	1.340	1.490	1.200	1.182	1.292	1.230	1.360	0.957	1.00	1.060
Turbidity (NTU)	5	8.3	6.1	3.76	0.41	3.8	4.6	5.3	1.74	2.48	0.98	2.1	6.1	2.17	0.11	1.56
Dissolved Oxygen (mg/L)	NV	1.9	4.3	1.0	1.5	5.5	100.90	77.30	1.2	10.3	5.7	4.2	16.5	4.7	36.2	11.0
Oxygen Reduction Potential (mV)	NV	-288	-223	-172.4	-327.3	-118.5	-190	-175	-157.2	-261.8	-68.7	-248	-207	-188.1	-181.1	20.7
<b>Volatile Organic Compounds - EPA</b>	Method 8260C (ug/L)															
Benzene	1	16	7.2	14	9.6	14	10	9.0	9.3	13	9.7	22	29	27	11	25
Toluene	5	3.1	1.7 J	2.3 J	1.9 J	2.1 J	2.6	3.0	2.6 J	3.2 J	2.3 J	2.8 J	4.3 J	4.0 J	1.4 J	3.2 J
Ethylbenzene	5	<	<	<	<	<	<	0.78 J	<	<	<	<	<	<	<	<
m,p-Xylene	5	6.1	3.7	4.1	4.4	3.5	6.3	7.8	6.7	8.8	5.4	2.1 J	5.2 J	3.9 J	1.4 J	3.4 J
o-Xylene	5	5.2	2.9	3.2	3.3	2.6	5.4	5.7	5.1	6.3	4.0 J	4.2 J	6.8 J	6.1 J	2.2 J	4.8 J
Xylene (Total)	5	11.3	6.6	7.3	7.7	6.1	11.7	14.0	11.8	15	9.4 J	6.3	12 J	10.0 J	3.6 J	8.2 J
1,3,5-Trimethylbenzene	5	2.6	2.3 J	2.2 J	2.3 J	1.7 J	2.4 J	3.1	3.1 J	3.8 J	2.7 J	1.8 J	<	<	1.1 J	<
1,2,4-Trimethylbenzene	5	2.2 J	1.8 J	1.7 J	1.8 J	1.4 J	2.8	3.6	3.5 J	4.3 J	2.7 J	2.4 J	3.4 J	3.0 J	1.2 J	<
Naphthalene*	10	93	60	54	66	66	230	160	200	270	220	490	500	460	190	460
Semi-Volatile Organic Compounds -	EPA Method 8270D (ug	g/L)														
Acetophenone	NV	<	<	<	<	0.413 J	<	0.58 J	<	<	0.561 J	<	<	<	<	0.492 J
Acenaphthylene	NV	3.28	3.50	2.66	1.95	3.24	20.4	30.0	19.8	28.4	22.1	15.7	26	19.3	7.61	17.0
Naphthalene*	10	43.6	36.00	31.1	21.8	32.6	108	150	111	141	106	198	240	246	63.3	198
2-Methylnaphthalene	NV	7.04	8.50	6.14	6.77	8.39	17.4	29.0	18.2	30.8	27.0	14.9	24	22.7	6.86	23.2
Acenaphthene*	20	3.58	3.90	3.24	2.39	3.42	6.04	8.80	6.44	10.2	8.69	4.83	6.5	7.06	2.21	5.68
Dibenzofuran	NV	10.9	9.40	9.20	6.80	10.1	20.1	28.0	19.7	32.0	24.5	9.82	16	18.2	4.24	13.8
Fluorene*	50	17.2	19.00	14.3	10.4	15.2	27.3	42.0	27.0	46.7	34.7	17.7	30	29.0	7.45	21.4
Phenanthrene*	50	53.1	42.00	42.8	25.3	36.3	27.4	56.0	20.6	33.8	30.7	26.9	38	44.5	8.84	30.0
Dibenzo (a,h)Anthracene*	NV	<	0.04 J	<	<	<	<	1.6	<	<	<	<	<	<	<	<
Carbazole	NV	8.82	5.60	6.64	4.44	8.48	18.7	20.0	15.9	18.8	19.8	29.1	31	37.6	9.37	26.2
Anthracene*	50	6.19	6.00	5.10	4.04	7.70	2.47	13.00	2.44	4.46	4.93	1.91	3.8	3.59	1.56	3.76
Fluoranthene*	50	11.6	11.00	9.41	5.78	10.9	2.63	39.00	2.03	2.78	3.38	4.69	7.3	5.95	2.44	6.32
Biphenyl	5	1.86	1.9 J	1.67	1.17	1.96	4.31	5.90	4.39	7.74	5.70	2.42	3.9	4.03	1.07	3.06
Pyrene*	50	8.10	7.00	6.28	3.51	6.39	2.50	33.00	1.90	2.64	2.59	4.02	4.9	4.90	1.84	4.06
Butyl benzyl phthalate*	50	<	<	0.083 J	<	<	<	<	<	<	<	<	<	<	<	<
Benz [a] Anthracene*	0.002	0.590	0.61	0.402 J	0.226 J	0.342 J	0.242 J	14	<	<	<	0.298 J	0.28	0.295 J	<	0.214 J
Benzo [b] Fluoranthene*	0.002	0.255 J	0.37	0.136 J	0.073 J	0.093 J	0.140 J	17	<	0.076 J	<	0.111 J	0.06 J	<	<	<
Benzo [k] Fluoranthene*	0.002	<	0.15	<	<	<	<	5.6	<	<	<	<	0.01 J	<	<	<
Benzo [a] Pyrene	ND	0.156 J	0.27	0.091 J	<	<	0.092 J	12	<	<	<	0.065 J	0.02 J	<	<	<
Indeno [1,2,3-cd] Pyrene	0.002	0.110 J	0.19	<	<	<	<	8.6	<	<	<	<	<	<	<	<
Benzo (g,h,i) Perylene	NV	0.114 J	0.17	<	<	<	<	7.6	<	<	<	<	<	<	<	<
Chrysene*	0.002	0.461 J	0.48	0.331 J	0.166 J	0.250 J	0.198 J	14	<	<	<	0.208 J	0.19	0.225 J	<	0.145 J
bis(2-Ethylhexyl)Phthalate	5	0.086 JB	<	<	<	<	0.094 JB	4.0	<	<	<	<	<	<	<	<

- 1. Compounds detected in one or more sample for the past five sampling events are presented on this table. Refer to Appendix B for list of all compounds included in analysis.
- 2. Analytical testing completed by Alpha Analytical in Westborough, MA.
- 3. NYSDEC Groundwater Class GA criteria obtained from Division of Water Technical and Operational Guidance Series (TOGS 1.1.1), dated October 1993, revised June 1998, errata January 1999 and amended April 2000 (Class GA).
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- 8. "D" qualifier = indicates the compound concentration was obtained from a secondary dilution analysis.
- 7. Value shown in **bold** indicates exceedance of respective Class GA Criteria or guidance value.
- 8. NV = no value, NT = not tested, ND = Not detected above method detection limit
- 9. \* = value shown is a guidance value rather than a groundwater standard.
- 10. The equipment used to collect water quality data was calibrated prior to and during use in accordance with the manufacturer's recommendations.
- 11. DO and pH measurements are routinely made using the same model water quality meter, however the measurements made on 9/2020 and 4/2021 appear erroneous.



**FIGURES** 







**APPENDIX A** 

**LIMITATIONS** 

# GZN

#### GEOHYDROLOGICAL LIMITATIONS

# Use of Report

1. GZA GeoEnvironmental, Inc. (GZA) prepared this report on behalf of, and for the exclusive use of our Client for the stated purpose(s) and location(s) identified in the Proposal for Services and/or Report. Use of this report, in whole or in part, at other locations, or for other purposes, may lead to inappropriate conclusions; and we do not accept any responsibility for the consequences of such use(s). Further, reliance by any party not expressly identified in the agreement, for any use, without our prior written permission, shall be at that party's sole risk, and without any liability to GZA.

# Standard of Care

- 2. GZA's findings and conclusions are based on the work conducted as part of the Scope of Services set forth in the Proposal for Services and/or Report and reflect our professional judgment. These findings and conclusions must be considered not as scientific or engineering certainties, but rather as our professional opinions concerning the limited data gathered during the course of our work. Conditions other than described in this report may be found at the subject location(s).
- 3. GZA's services were performed using the degree of skill and care ordinarily exercised by qualified professionals performing the same type of services, at the same time, under similar conditions, at the same or a similar property. No warranty, expressed or implied, is made. Specifically, GZA does not and cannot represent that the Site contains no hazardous material, oil, or other latent condition beyond that observed by GZA during its study. Additionally, GZA makes no warranty that any response action or recommended action will achieve all of its objectives or that the findings of this study will be upheld by a local, state or federal agency.
- 4. In conducting our work, GZA relied upon certain information made available by public agencies, Client and/or others. GZA did not attempt to independently verify the accuracy or completeness of that information. Inconsistencies in this information which we have noted, if any, are discussed in the Report.

# **Subsurface Conditions**

5. The generalized soil profile(s) provided in our Report are based on widely-spaced subsurface explorations and are intended only to convey trends in subsurface conditions. The boundaries between strata are approximate and idealized, and were based on our assessment of subsurface conditions. The composition of strata, and the transitions between strata, may be more variable and more complex than indicated. For more specific information on soil conditions at a specific location refer to the exploration logs. The nature and extent of variations between these explorations may not become evident until further exploration or construction. If variations or other latent conditions then become evident, it will be necessary to reevaluate the conclusions and recommendations of this report.

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6. Water level readings have been made, as described in this Report, in and monitoring wells at the specified times and under the stated conditions. These data have been reviewed and interpretations have been made in this report. Fluctuations in the level of the groundwater however occur due to temporal or spatial variations in areal recharge rates, soil heterogeneities, the presence of subsurface utilities, and/or natural or artificially induced perturbations. The observed water table may be other than indicated in the Report.

# Compliance with Codes and Regulations

7. We used reasonable care in identifying and interpreting applicable codes and regulations necessary to execute our scope of work. These codes and regulations are subject to various, and possibly contradictory, interpretations. Interpretations and compliance with codes and regulations by other parties is beyond our control.

# Screening and Analytical Testing

- 8. GZA collected environmental samples at the locations identified in the Report. These samples were analyzed for the specific parameters identified in the report. Additional constituents, for which analyses were not conducted, may be present in soil, groundwater, surface water, sediment and/or air. Future Site activities and uses may result in a requirement for additional testing.
- 9. Our interpretation of field screening and laboratory data is presented in the Report. Unless otherwise noted, we relied upon the laboratory's QA/QC program to validate these data.
- 10. Variations in the types and concentrations of contaminants observed at a given location or time may occur due to release mechanisms, disposal practices, changes in flow paths, and/or the influence of various physical, chemical, biological or radiological processes. Subsequently observed concentrations may be other than indicated in the Report.

### Interpretation of Data

11. Our opinions are based on available information as described in the Report, and on our professional judgment. Additional observations made over time, and/or space, may not support the opinions provided in the Report.

# Additional Information

12. In the event that the Client or others authorized to use this report obtain additional information on environmental or hazardous waste issues at the Site not contained in this report, such information shall be brought to GZA's attention forthwith. GZA will evaluate such information and, on the basis of this evaluation, may modify the conclusions stated in this report.

November 2022 PAGE 2

# **Additional Services**

13. GZA recommends that we be retained to provide services during any future investigations, design, implementation activities, construction, and/or property development/ redevelopment at the Site. This will allow us the opportunity to: i) observe conditions and compliance with our design concepts and opinions; ii) allow for changes in the event that conditions are other than anticipated; iii) provide modifications to our design; and iv) assess the consequences of changes in technologies and/or regulations.

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# APPENDIX B ANALYTICAL TEST RESULTS



#### ANALYTICAL REPORT

Lab Number: L2249775

Client: GZA GeoEnvironmental of New York

300 Pearl Street

Suite 700

Buffalo, NY 14202

ATTN: Dan Troy

Phone: (716) 844-7050

Project Name: STEELWINDS ANNUAL/SEMIANNUAL

Project Number: 03.0033579.15

Report Date: 10/11/22

The original project report/data package is held by Alpha Analytical. This report/data package is paginated and should be reproduced only in its entirety. Alpha Analytical holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0574), IL (200077), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #P330-17-00196).

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

**Lab Number:** L2249775 **Report Date:** 10/11/22

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2249775-01	WT1-05-091322	WATER	LACKAWANNA NY	09/13/22 07:55	09/13/22
L2249775-02	MWN-01-091322	WATER	LACKAWANNA NY	09/13/22 09:05	09/13/22
L2249775-03	MWN-01B-091322	WATER	LACKAWANNA NY	09/13/22 10:15	09/13/22
L2249775-04	WT1-04-091322	WATER	LACKAWANNA NY	09/13/22 10:55	09/13/22
L2249775-05	BCP-ORC-1-091322	WATER	LACKAWANNA NY	09/13/22 13:25	09/13/22
L2249775-06	WT1-02-091322	WATER	LACKAWANNA NY	09/13/22 14:15	09/13/22
L2249775-07	TRIP BLANK-1	WATER	LACKAWANNA NY	09/13/22 00:00	09/13/22

Serial No:10112209:57

**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL Lab Number: L2249775

**Project Number:** 03.0033579.15 **Report Date:** 10/11/22

#### **Case Narrative**

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments, solids and tissues are reported on a dry weight basis unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

HOLD POLICY - For samples submitted on hold, Alpha's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Alpha Project Manager and made arrangements for Alpha to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.	



Serial\_No:10112209:57

Project Name: STEELWINDS ANNUAL/SEMIANNUAL Lab Number: L2249775

**Project Number:** 03.0033579.15 **Report Date:** 10/11/22

### **Case Narrative (continued)**

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

Title: Technical Director/Representative Date: 10/11/22

Custen Walker Cristin Walker

# **ORGANICS**



# **VOLATILES**



L2249775

10/11/22

**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Date Collected: 09/13/22 07:55

Lab Number:

Report Date:

Lab ID: L2249775-01 D Client ID: Date Received: 09/13/22 WT1-05-091322 Field Prep: Sample Location: Not Specified LACKAWANNA NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/24/22 20:33

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westbord	ough Lab					
Benzene	9.7		ug/l	1.0	0.32	2
Toluene	2.3	J	ug/l	5.0	1.4	2
Ethylbenzene	ND		ug/l	5.0	1.4	2
Methyl tert butyl ether	ND		ug/l	5.0	1.4	2
p/m-Xylene	5.4		ug/l	5.0	1.4	2
o-Xylene	4.0	J	ug/l	5.0	1.4	2
n-Butylbenzene	ND		ug/l	5.0	1.4	2
sec-Butylbenzene	ND		ug/l	5.0	1.4	2
tert-Butylbenzene	ND		ug/l	5.0	1.4	2
Isopropylbenzene	ND		ug/l	5.0	1.4	2
p-Isopropyltoluene	ND		ug/l	5.0	1.4	2
Naphthalene	220		ug/l	5.0	1.4	2
n-Propylbenzene	ND		ug/l	5.0	1.4	2
1,3,5-Trimethylbenzene	2.7	J	ug/l	5.0	1.4	2
1,2,4-Trimethylbenzene	2.7	J	ug/l	5.0	1.4	2

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	107	70-130	
Toluene-d8	94	70-130	
4-Bromofluorobenzene	99	70-130	
Dibromofluoromethane	101	70-130	



L2249775

**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

L2249775-02

MWN-01-091322

LACKAWANNA NY

D

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

10/11/22

Report Date:

Lab Number:

Date Collected: 09/13/22 09:05

Date Received: 09/13/22 Field Prep: Not Specified

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/24/22 20:12

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Wes	tborough Lab					
Benzene	12		ug/l	1.0	0.32	2
Toluene	2.8	J	ug/l	5.0	1.4	2
Ethylbenzene	ND		ug/l	5.0	1.4	2
Methyl tert butyl ether	ND		ug/l	5.0	1.4	2
p/m-Xylene	6.0		ug/l	5.0	1.4	2
o-Xylene	5.0		ug/l	5.0	1.4	2
n-Butylbenzene	ND		ug/l	5.0	1.4	2
sec-Butylbenzene	ND		ug/l	5.0	1.4	2
tert-Butylbenzene	ND		ug/l	5.0	1.4	2
Isopropylbenzene	ND		ug/l	5.0	1.4	2
p-Isopropyltoluene	ND		ug/l	5.0	1.4	2
Naphthalene	240		ug/l	5.0	1.4	2
n-Propylbenzene	ND		ug/l	5.0	1.4	2
1,3,5-Trimethylbenzene	2.8	J	ug/l	5.0	1.4	2
1,2,4-Trimethylbenzene	3.0	J	ug/l	5.0	1.4	2

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	119	70-130	
Toluene-d8	96	70-130	
4-Bromofluorobenzene	98	70-130	
Dibromofluoromethane	110	70-130	



L2249775

**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Lab Number:

Report Date: 10/11/22

Lab ID: L2249775-03 Date Collected: 09/13/22 10:15

Client ID: Date Received: 09/13/22 MWN-01B-091322 Field Prep: Sample Location: Not Specified LACKAWANNA NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/24/22 20:54

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westbo	rough Lab					
Benzene	55		ug/l	0.50	0.16	1
Toluene	20		ug/l	2.5	0.70	1
Ethylbenzene	0.95	J	ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	15		ug/l	2.5	0.70	1
o-Xylene	11		ug/l	2.5	0.70	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
Isopropylbenzene	1.4	J	ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	780	E	ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	5.2		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	7.4		ug/l	2.5	0.70	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	109	70-130	
Toluene-d8	96	70-130	
4-Bromofluorobenzene	98	70-130	
Dibromofluoromethane	99	70-130	



L2249775

**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

**Report Date:** 10/11/22

Lab Number:

Lab ID: D L2249775-03 Client ID: MWN-01B-091322 Sample Location: LACKAWANNA NY

Date Collected: 09/13/22 10:15 Date Received: 09/13/22 Field Prep: Not Specified

70-130

70-130

70-130

Sample Depth:

Toluene-d8

4-Bromofluorobenzene

Dibromofluoromethane

Matrix: Water 1,8260C Analytical Method: Analytical Date: 09/27/22 11:27

Analyst: MKS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Wes	stborough Lab					
Naphthalene	1500		ug/l	50	14.	20
Surrogate			% Recovery	Qualifier		eptance riteria
1,2-Dichloroethane-d4			102			70-130

96

93

113



L2249775

09/13/22 10:55

**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

10/11/22

Report Date:

Lab Number:

Date Collected:

Lab ID: L2249775-04

Client ID: WT1-04-091322 Sample Location: LACKAWANNA NY

Date Received: 09/13/22 Field Prep: Not Specified

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/24/22 19:51

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westbor	ough Lab					
Benzene	14		ug/l	0.50	0.16	1
Toluene	2.1	J	ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	3.5		ug/l	2.5	0.70	1
o-Xylene	2.6		ug/l	2.5	0.70	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	66		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	1.7	J	ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	1.4	J	ug/l	2.5	0.70	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	125	70-130	
Toluene-d8	96	70-130	
4-Bromofluorobenzene	93	70-130	
Dibromofluoromethane	113	70-130	



L2249775

10/11/22

**Project Name:** Lab Number: STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Date Collected: 09/13/22 13:25

Report Date:

Lab ID: D L2249775-05 Client ID:

Date Received: 09/13/22 BCP-ORC-1-091322 Field Prep: Sample Location: Not Specified LACKAWANNA NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/24/22 19:30

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Volatile Organics by GC/MS - Westbe	orough Lab						
Benzene	25		ug/l	2.0	0.64	4	
Toluene	3.2	J	ug/l	10	2.8	4	
Ethylbenzene	ND		ug/l	10	2.8	4	
Methyl tert butyl ether	ND		ug/l	10	2.8	4	
p/m-Xylene	3.4	J	ug/l	10	2.8	4	
o-Xylene	4.8	J	ug/l	10	2.8	4	
n-Butylbenzene	ND		ug/l	10	2.8	4	
sec-Butylbenzene	ND		ug/l	10	2.8	4	
tert-Butylbenzene	ND		ug/l	10	2.8	4	
Isopropylbenzene	ND		ug/l	10	2.8	4	
p-Isopropyltoluene	ND		ug/l	10	2.8	4	
Naphthalene	460		ug/l	10	2.8	4	
n-Propylbenzene	ND		ug/l	10	2.8	4	
1,3,5-Trimethylbenzene	ND		ug/l	10	2.8	4	
1,2,4-Trimethylbenzene	ND		ug/l	10	2.8	4	

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	121	70-130	
Toluene-d8	87	70-130	
4-Bromofluorobenzene	94	70-130	
Dibromofluoromethane	109	70-130	

L2249775

Project Name: STEELWINDS ANNUAL/SEMIANNUAL

L2249775-06

WT1-02-091322

LACKAWANNA NY

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

**Report Date:** 10/11/22

Lab Number:

Date Collected: 09/13/22 14:15
Date Received: 09/13/22
Field Prep: Not Specified

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Matrix: Water
Analytical Method: 1,8260C
Analytical Date: 09/24/22 19:09

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westboro	ugh Lab					
Benzene	8.7		ug/l	0.50	0.16	1
Toluene	1.7	J	ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	2.6		ug/l	2.5	0.70	1
o-Xylene	1.9	J	ug/l	2.5	0.70	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	27		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	1.2	J	ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	0.84	J	ug/l	2.5	0.70	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	114	70-130	
Toluene-d8	95	70-130	
4-Bromofluorobenzene	84	70-130	
Dibromofluoromethane	108	70-130	

L2249775

09/13/22 00:00

Not Specified

09/13/22

**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

**Report Date:** 10/11/22

Lab Number:

Date Collected:

Date Received:

Field Prep:

Lab ID: L2249775-07

Client ID: TRIP BLANK-1 Sample Location: LACKAWANNA NY

Sample Depth:

Matrix: Water
Analytical Method: 1,8260C
Analytical Date: 09/24/22 18:47

Parameter	Result	Qualifier Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - W	estborough Lab				
Benzene	ND	ug/l	0.50	0.16	1
Toluene	ND	ug/l	2.5	0.70	1
Ethylbenzene	ND	ug/l	2.5	0.70	1
Methyl tert butyl ether	ND	ug/l	2.5	0.70	1
p/m-Xylene	ND	ug/l	2.5	0.70	1
o-Xylene	ND	ug/l	2.5	0.70	1
n-Butylbenzene	ND	ug/l	2.5	0.70	1
sec-Butylbenzene	ND	ug/l	2.5	0.70	1
tert-Butylbenzene	ND	ug/l	2.5	0.70	1
Isopropylbenzene	ND	ug/l	2.5	0.70	1
p-Isopropyltoluene	ND	ug/l	2.5	0.70	1
Naphthalene	ND	ug/l	2.5	0.70	1
n-Propylbenzene	ND	ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND	ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	ND	ug/l	2.5	0.70	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	105	70-130	
Toluene-d8	97	70-130	
4-Bromofluorobenzene	97	70-130	
Dibromofluoromethane	114	70-130	



L2249775

Project Name: STEELWINDS ANNUAL/SEMIANNUAL Lab Number:

**Project Number:** 03.0033579.15 **Report Date:** 10/11/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/24/22 13:34

Analyst: TMS

Parameter	Result	Qualifier Units	RL	MDL	
olatile Organics by GC/MS	- Westborough Lab	for sample(s):	01-07 Batch:	WG1692517-5	
Benzene	ND	ug/l	0.50	0.16	
Toluene	ND	ug/l	2.5	0.70	
Ethylbenzene	ND	ug/l	2.5	0.70	
Methyl tert butyl ether	ND	ug/l	2.5	0.70	
p/m-Xylene	ND	ug/l	2.5	0.70	
o-Xylene	ND	ug/l	2.5	0.70	
n-Butylbenzene	ND	ug/l	2.5	0.70	
sec-Butylbenzene	ND	ug/l	2.5	0.70	
tert-Butylbenzene	ND	ug/l	2.5	0.70	
Isopropylbenzene	ND	ug/l	2.5	0.70	
p-Isopropyltoluene	ND	ug/l	2.5	0.70	
Naphthalene	ND	ug/l	2.5	0.70	
n-Propylbenzene	ND	ug/l	2.5	0.70	
1,3,5-Trimethylbenzene	ND	ug/l	2.5	0.70	
1,2,4-Trimethylbenzene	ND	ug/l	2.5	0.70	

	Acceptance					
Surrogate	%Recovery Qualifi	er Criteria				
1,2-Dichloroethane-d4	111	70-130				
Toluene-d8	101	70-130				
4-Bromofluorobenzene	100	70-130				
Dibromofluoromethane	104	70-130				



Project Name: STEELWINDS ANNUAL/SEMIANNUAL Lab Number: L2249775

**Project Number:** 03.0033579.15 **Report Date:** 10/11/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/27/22 11:02

Analyst: PD

Parameter	Result	Qualifier Units	RL	MDL
Volatile Organics by GC/MS - W	estborough Lab	for sample(s): 03	Batch:	WG1692957-5
Benzene	ND	ug/l	0.50	0.16
Toluene	ND	ug/l	2.5	0.70
Ethylbenzene	ND	ug/l	2.5	0.70
Methyl tert butyl ether	ND	ug/l	2.5	0.70
p/m-Xylene	ND	ug/l	2.5	0.70
o-Xylene	ND	ug/l	2.5	0.70
n-Butylbenzene	ND	ug/l	2.5	0.70
sec-Butylbenzene	ND	ug/l	2.5	0.70
tert-Butylbenzene	ND	ug/l	2.5	0.70
Isopropylbenzene	ND	ug/l	2.5	0.70
p-Isopropyltoluene	ND	ug/l	2.5	0.70
Naphthalene	ND	ug/l	2.5	0.70
n-Propylbenzene	ND	ug/l	2.5	0.70
1,3,5-Trimethylbenzene	ND	ug/l	2.5	0.70
1,2,4-Trimethylbenzene	ND	ug/l	2.5	0.70

	Acceptan					
Surrogate	%Recovery Qualifi	er Criteria				
1,2-Dichloroethane-d4	102	70-130				
Toluene-d8	97	70-130				
4-Bromofluorobenzene	94	70-130				
Dibromofluoromethane	113	70-130				



**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

Lab Number: L2249775

**Report Date:** 10/11/22

arameter	LCS %Recovery	Qual	LCSD %Recovery		%Recovery Limits	RPD	Qual	RPD Limits
	·					10. 2		
olatile Organics by GC/MS - Westborough L	ab Associated	sample(s):	01-07 Batch:	WG1692517-3	WG1692517-4			
Benzene	95		97		70-130	2		20
Toluene	100		100		70-130	0		20
Ethylbenzene	96		95		70-130	1		20
Methyl tert butyl ether	82		69		63-130	17		20
p/m-Xylene	100		100		70-130	0		20
o-Xylene	95		95		70-130	0		20
n-Butylbenzene	100		99		53-136	1		20
sec-Butylbenzene	97		93		70-130	4		20
tert-Butylbenzene	92		88		70-130	4		20
Isopropylbenzene	88		86		70-130	2		20
p-Isopropyltoluene	95		90		70-130	5		20
Naphthalene	72		74		70-130	3		20
n-Propylbenzene	95		93		69-130	2		20
1,3,5-Trimethylbenzene	96		92		64-130	4		20
1,2,4-Trimethylbenzene	93		88		70-130	6		20

Surrogate	LCS %Recovery Qua	LCSD I %Recovery Qual	Acceptance Criteria
1,2-Dichloroethane-d4	110	112	70-130
Toluene-d8	102	101	70-130
4-Bromofluorobenzene	93	94	70-130
Dibromofluoromethane	102	101	70-130



**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

Lab Number: L2249775

**Report Date:** 10/11/22

arameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
olatile Organics by GC/MS - Westborough L	ab Associated	sample(s): 03	Batch: WG1	692957-3	WG1692957-4			
Benzene	110		110		70-130	0		20
Toluene	98		100		70-130	2		20
Ethylbenzene	100		100		70-130	0		20
Methyl tert butyl ether	85		89		63-130	5		20
p/m-Xylene	100		105		70-130	5		20
o-Xylene	95		100		70-130	5		20
n-Butylbenzene	97		100		53-136	3		20
sec-Butylbenzene	96		100		70-130	4		20
tert-Butylbenzene	81		98		70-130	19		20
Isopropylbenzene	95		100		70-130	5		20
p-Isopropyltoluene	94		100		70-130	6		20
Naphthalene	84		90		70-130	7		20
n-Propylbenzene	99		100		69-130	1		20
1,3,5-Trimethylbenzene	92		97		64-130	5		20
1,2,4-Trimethylbenzene	92		97		70-130	5		20

	LCS	LCSD	Acceptance
Surrogate	%Recovery Qua	I %Recovery Qual	Criteria
1,2-Dichloroethane-d4	93	96	70-130
Toluene-d8	96	96	70-130
4-Bromofluorobenzene	94	94	70-130
Dibromofluoromethane	103	105	70-130



### **SEMIVOLATILES**



L2249775

10/11/22

Project Name: STEELWINDS ANNUAL/SEMIANNUAL Lab Number:

**Project Number:** 03.0033579.15

L2249775-01

**SAMPLE RESULTS** 

Date Collected: 09/13/22 07:55

**Report Date:** 

Client ID: WT1-05-091322 Date Received: 09/13/22
Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Lab ID:

Analyst:

Matrix: Water Extraction Method: EPA 3510C
Analytical Method: 1,8270D Extraction Date: 09/17/22 08:30

Analytical Date: 09/19/22 12:11

PS

Result Qualifier Units RL MDL **Dilution Factor Parameter** Semivolatile Organics by GC/MS - Mansfield Lab bis(2-Chloroethyl)ether ND 0.495 0.092 1 ug/l 1,3-Dichlorobenzene ND 0.495 0.078 ug/l 1,4-Dichlorobenzene ND ug/l 0.495 0.082 1 1,2-Dichlorobenzene ND ug/l 0.495 0.067 1 Benzyl alcohol ND ug/l 0.495 0.122 1 bis(2-chloroisopropyl)ether ND ug/l 0.495 0.107 1 0.561 J 0.990 0.205 Acetophenone ug/l 1 Hexachloroethane ND 0.495 0.101 1 ug/l ND Nitrobenzene ug/l 0.495 0.101 1 Isophorone ND 0.495 0.125 1 ug/l bis(2-Chloroethoxy)methane ND 0.495 0.085 1 ug/l ND 1,2,4-Trichlorobenzene 0.495 0.095 ug/l 1 Е 0.087 Naphthalene 62.1 0.495 1 ug/l 4-Chloroaniline ND ug/l 0.495 0.127 1 Hexachlorobutadiene ND 0.495 0.085 1 ug/l 27.0 0.495 0.090 2-Methylnaphthalene 1 ug/l ND 1,2,4,5-Tetrachlorobenzene ug/l 0.495 0.079 1 ND Hexachlorocyclopentadiene 0.495 0.151 1 ug/l Biphenyl 5.70 0.495 0.110 1 ug/l 2-Chloronaphthalene ND 0.089 1 ug/l 0.495 ND 2-Nitroaniline 0.495 0.137 1 ug/l Acenaphthylene 22.1 ug/l 0.495 0.111 1 ND Dimethylphthalate 0.495 0.116 1 ug/l 2,6-Dinitrotoluene ND 0.495 0.166 1 ug/l Acenaphthene 8.69 0.495 0.095 1 ug/l 3-Nitroaniline ND 0.110 1 ug/l 0.495 24.5 0.090 Dibenzofuran 0.495 1 ug/l ND 2,4-Dinitrotoluene ug/l 0.495 0.161 1



10/11/22

Project Name: STEELWINDS ANNUAL/SEMIANNUAL Lab Number: L2249775

**Project Number:** 03.0033579.15

L2249775-01

**SAMPLE RESULTS** 

Date Collected: 09/13/22 07:55

Report Date:

Client ID: WT1-05-091322 Date Received: 09/13/22 Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Lab ID:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS - Mansfield Lab								
Fluorene	34.7		ug/l	0.495	0.103	1		
Diethylphthalate	ND		ug/l	0.495	0.178	1		
4-Nitroaniline	ND		ug/l	0.495	0.111	1		
n-Nitrosodiphenylamine	ND		ug/l	0.495	0.071	1		
Hexachlorobenzene	ND		ug/l	0.495	0.121	1		
Phenanthrene	30.7		ug/l	0.495	0.110	1		
Anthracene	4.93		ug/l	0.495	0.136	1		
Carbazole	19.8		ug/l	0.495	0.142	1		
Di-n-butylphthalate	ND		ug/l	0.495	0.099	1		
Fluoranthene	3.38		ug/l	0.495	0.154	1		
Pyrene	2.59		ug/l	0.495	0.168	1		
Butylbenzylphthalate	ND		ug/l	0.495	0.084	1		
3,3'-Dichlorobenzidine	ND		ug/l	0.495	0.191	1		
Benz(a)anthracene	ND		ug/l	0.495	0.182	1		
Chrysene	ND		ug/l	0.495	0.140	1		
bis(2-Ethylhexyl)phthalate	ND		ug/l	0.495	0.080	1		
Di-n-octylphthalate	ND		ug/l	0.990	0.078	1		
Benzo(b)fluoranthene	ND		ug/l	0.495	0.065	1		
Benzo(k)fluoranthene	ND		ug/l	0.495	0.159	1		
Benzo(a)pyrene	ND		ug/l	0.495	0.060	1		
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.495	0.089	1		
Dibenz(a,h)anthracene	ND		ug/l	0.495	0.064	1		
Benzo(g,h,i)perylene	ND		ug/l	0.495	0.108	1		

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	60	15-115
Phenol-d5	42	15-115
Nitrobenzene-d5	94	30-130
2-Fluorobiphenyl	78	30-130
2,4,6-Tribromophenol	112	15-115
Terphenyl-d14	97	30-130



10/11/22

Report Date:

Project Name: STEELWINDS ANNUAL/SEMIANNUAL Lab Number: L2249775

**Project Number:** 03.0033579.15

09/20/22 12:24

SAMPLE RESULTS

Lab ID: L2249775-01 D Date Collected: 09/13/22 07:55

Client ID: WT1-05-091322 Date Received: 09/13/22 Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Analytical Date:

Matrix: Water Extraction Method: EPA 3510C
Analytical Method: 1,8270D Extraction Date: 09/17/22 08:30

Analyst: PS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Semivolatile Organics by GC/MS - Mansfield Lab							
Naphthalene	106		ug/l	2.48	0.434	5	

		Acceptance
Surrogate	% Recovery	Qualifier Criteria
2-Fluorophenol	61	15-115
Phenol-d5	41	15-115
Nitrobenzene-d5	88	30-130
2-Fluorobiphenyl	86	30-130
2,4,6-Tribromophenol	101	15-115
Terphenyl-d14	100	30-130



L2249775

10/11/22

Project Name: STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

L2249775-02

MWN-01-091322

LACKAWANNA NY

**SAMPLE RESULTS** 

Date Collected: 09/13/22 09:05

Date Received: 09/13/22

Lab Number:

Report Date:

Field Prep: Not Specified

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Matrix: Water
Analytical Method: 1,8270D
Analytical Date: 09/19/22 12:40

Analyst: PS

Extraction Method: EPA 3510C Extraction Date: 09/17/22 08:30

1,3-Dichlorobenzene         ND         ug/l         0.495         0.078         1           1,4-Dichlorobenzene         ND         ug/l         0.495         0.082         1           1,2-Dichlorobenzene         ND         ug/l         0.495         0.067         1           Benzyl alcohol         ND         ug/l         0.495         0.122         1           Isio(2-chloroisopropyl)ether         ND         ug/l         0.495         0.107         1           Acetophenone         0.570         J         ug/l         0.495         0.101         1           Hexachloroethane         ND         ug/l         0.495         0.101         1           Nitrobenzene         ND         ug/l         0.495         0.101         1           Isophorone         ND         ug/l         0.495         0.101         1           Isophorone         ND         ug/l         0.495         0.085         1           Isophorone         ND         ug/l         0.495         0.085         1           ND         ug/l         0.495         0.085         1           Isophithalene         59.4         E         ug/l         0.495         0.085	Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
1,3-Dichlorobenzene         ND         ug/l         0.495         0.078         1           1,4-Dichlorobenzene         ND         ug/l         0.495         0.082         1           1,2-Dichlorobenzene         ND         ug/l         0.495         0.067         1           Benzyl alcohol         ND         ug/l         0.495         0.122         1           Benzyl alcohol         ND         ug/l         0.495         0.102         1           Acetophenone         0.570         J         ug/l         0.495         0.101         1           Acetophenone         0.570         J         ug/l         0.495         0.101         1           Hexachloroethrane         ND         ug/l         0.495         0.101         1           Nitrobenzene         ND         ug/l         0.495         0.101         1           Isophorone         ND         ug/l         0.495         0.011         1           Isophorone         ND         ug/l         0.495         0.085         1           Isophitalene         ND         ug/l         0.495         0.085         1           I-2,4-Trichlorobenzene         ND         ug/l <td< td=""><td>Semivolatile Organics by GC/MS - Ma</td><td>nsfield Lab</td><td></td><td></td><td></td><td></td><td></td></td<>	Semivolatile Organics by GC/MS - Ma	nsfield Lab					
1,4-Dichlorobenzene   ND   ug/l   0,495   0,062   1	bis(2-Chloroethyl)ether	ND		ug/l	0.495	0.092	1
1,2-Dichlorobenzene   ND   ug/l   0,495   0,067   1	1,3-Dichlorobenzene	ND		ug/l	0.495	0.078	1
Benzyl alcohol         ND         ug/l         0.495         0.122         1           bis(2-chloroisopropyl)ether         ND         ug/l         0.495         0.107         1           Acetophenone         0.570         J         ug/l         0.990         0.205         1           Hexachloroethane         ND         ug/l         0.495         0.101         1           Nitrobenzene         ND         ug/l         0.495         0.101         1           Isophorone         ND         ug/l         0.495         0.101         1           Isophorone         ND         ug/l         0.495         0.102         1           Isophorone         ND         ug/l         0.495         0.125         1           Isophorone         ND         ug/l         0.495         0.085         1           Isophorone         ND         ug/l         0.495         0.085         1           L2.4-Trichlorobarzene         ND         ug/l         0.495         0.085         1           Hexachlorobutadiene         ND         ug/l         0.495         0.090         1           Hexachlorobutadiene         ND         ug/l         0.495	1,4-Dichlorobenzene	ND		ug/l	0.495	0.082	1
ND   Ug/l   0.495   0.107   1   1   1   1   1   1   1   1   1	1,2-Dichlorobenzene	ND		ug/l	0.495	0.067	1
Acetophenone 0.570 J ug/l 0.990 0.205 1 Hexachloroethane ND ug/l 0.495 0.101 1 Nitrobenzene ND ug/l 0.495 0.101 1 Isophorone ND ug/l 0.495 0.101 1 Isophorone ND ug/l 0.495 0.125 1 Isophorone ND ug/l 0.495 0.125 1 Isophorone ND ug/l 0.495 0.085 1 I.2.4-Trichlorobenzene ND ug/l 0.495 0.085 1 I.2.4-Trichlorobenzene ND ug/l 0.495 0.095 1 I.2.4-Trichlorobenzene ND ug/l 0.495 0.095 1 Isophorone ND ug/l 0.495 0.087 1 Isophorone ND ug/l 0.495 0.085 1 Isophorone ND ug/l 0.495 0.090 1 Isophorone ND ug/l 0.495 0.151 1 Isophorone ND ug/l 0.495 0.110 1 Isophorone ND ug/l 0.495 0.110 1 Isophorone ND ug/l 0.495 0.111 1 Isophorone ND ug/l 0.495 0.116 1 Isophorone ND u	Benzyl alcohol	ND		ug/l	0.495	0.122	1
ND	bis(2-chloroisopropyl)ether	ND		ug/l	0.495	0.107	1
Nitrobenzene ND ug/l 0.495 0.101 1 Isophorone ND ug/l 0.495 0.125 1 Isophorone ND ug/l 0.495 0.085 1 I.2.4-Trichlorobenzene ND ug/l 0.495 0.095 1 I.2.4-Trichlorobenzene ND ug/l 0.495 0.095 1 I.2.4-Trichlorobenzene ND ug/l 0.495 0.095 1 I.2.4-Trichlorobenzene ND ug/l 0.495 0.087 1 I.2.4-Trichlorobenzene ND ug/l 0.495 0.087 1 I.2.4-Trichlorobenzene ND ug/l 0.495 0.087 1 I.2.4-Trichlorobenzene ND ug/l 0.495 0.085 1 I.2.4-S-Tetrachlorobenzene ND ug/l 0.495 0.085 1 I.2.4-S-Tetrachlorobenzene ND ug/l 0.495 0.090 1 I.2.4-S-Tetrachlorobenzene ND ug/l 0.495 0.079 1 I.2.4-S-Tetrachlorobenzene ND ug/l 0.495 0.151 1 Isophenyl 6.48 ug/l 0.495 0.110 1 I.2.4-Chloronaphthalene ND ug/l 0.495 0.110 1 I.2.4-Chloronaphthalene ND ug/l 0.495 0.110 1 I.2.4-Chloronaphthalene ND ug/l 0.495 0.111 1 I.2.4-Chloronaphthalene ND ug/l 0.495 0.111 1 I.2.4-Chloronaphthalene ND ug/l 0.495 0.116 1 I.2.4-Chloronaphthalene ND ug/l 0.495 0.110 1 I.2.4-Chloronaphthal	Acetophenone	0.570	J	ug/l	0.990	0.205	1
ND	Hexachloroethane	ND		ug/l	0.495	0.101	1
ND	Nitrobenzene	ND		ug/l	0.495	0.101	1
1,2,4-Trichlorobenzene   ND	Isophorone	ND		ug/l	0.495	0.125	1
Naphthalene 59.4 E ug/l 0.495 0.087 1 4-Chloroaniline ND ug/l 0.495 0.127 1 Hexachlorobutadiene ND ug/l 0.495 0.085 1 2-Methylnaphthalene 27.8 ug/l 0.495 0.090 1 1,2,4,5-Tetrachlorobenzene ND ug/l 0.495 0.079 1 Hexachlorocyclopentadiene ND ug/l 0.495 0.151 1 Biphenyl 6.48 ug/l 0.495 0.151 1 Biphenyl 6.48 ug/l 0.495 0.110 1 2-Chloronaphthalene ND ug/l 0.495 0.110 1 2-Chloronaphthalene ND ug/l 0.495 0.110 1 2-Chloronaphthalene ND ug/l 0.495 0.137 1 Chloronaphthalene ND ug/l 0.495 0.111 1 Dimethylphthalate ND ug/l 0.495 0.111 1 Dimethylphthalate ND ug/l 0.495 0.116 1 2,6-Dinitrotoluene ND ug/l 0.495 0.166 1 Acenaphthene 10.1 ug/l 0.495 0.095 1 3-Nitroaniline ND ug/l 0.495 0.095 1 3-Nitroaniline ND ug/l 0.495 0.110 1 Dibenzofuran 29.7 ug/l 0.495 0.110 1	bis(2-Chloroethoxy)methane	ND		ug/l	0.495	0.085	1
A-Chloroaniline ND ug/l 0.495 0.127 1  Hexachlorobutadiene ND ug/l 0.495 0.085 1  2-Methylnaphthalene 27.8 ug/l 0.495 0.090 1  1,2,4,5-Tetrachlorobenzene ND ug/l 0.495 0.079 1  Hexachlorocyclopentadiene ND ug/l 0.495 0.151 1  Biphenyl 6.48 ug/l 0.495 0.110 1  2-Chloronaphthalene ND ug/l 0.495 0.110 1  2-Nitroaniline ND ug/l 0.495 0.110 1  Acenaphthylene 23.5 ug/l 0.495 0.117 1  Dimethylphthalate ND ug/l 0.495 0.111 1  Dimethylphthalate ND ug/l 0.495 0.111 1  2,6-Dinitrotoluene ND ug/l 0.495 0.116 1  2,6-Dinitrotoluene ND ug/l 0.495 0.116 1  3-Nitroaniline ND ug/l 0.495 0.166 1  Acenaphthene 10.1 ug/l 0.495 0.095 1  3-Nitroaniline ND ug/l 0.495 0.095 1  3-Nitroaniline ND ug/l 0.495 0.090 1	1,2,4-Trichlorobenzene	ND		ug/l	0.495	0.095	1
Hexachlorobutadiene   ND	Naphthalene	59.4	E	ug/l	0.495	0.087	1
2-Methylnaphthalene 27.8 ug/l 0.495 0.090 1 1,2,4,5-Tetrachlorobenzene ND ug/l 0.495 0.079 1 Hexachlorocyclopentadiene ND ug/l 0.495 0.151 1 Biphenyl 6.48 ug/l 0.495 0.110 1 2-Chloronaphthalene ND ug/l 0.495 0.089 1 2-Nitroaniline ND ug/l 0.495 0.137 1 Acenaphthylene 23.5 ug/l 0.495 0.111 1 Dimethylphthalate ND ug/l 0.495 0.111 1 2,6-Dinitrotoluene ND ug/l 0.495 0.116 1 Acenaphthene 10.1 ug/l 0.495 0.166 1 Acenaphthene 10.1 ug/l 0.495 0.095 1 3-Nitroaniline ND ug/l 0.495 0.095 1 3-Nitroaniline ND ug/l 0.495 0.110 1	4-Chloroaniline	ND		ug/l	0.495	0.127	1
1,2,4,5-Tetrachlorobenzene ND ug/l 0.495 0.079 1  Hexachlorocyclopentadiene ND ug/l 0.495 0.151 1  Biphenyl 6.48 ug/l 0.495 0.110 1  2-Chloronaphthalene ND ug/l 0.495 0.089 1  2-Nitroaniline ND ug/l 0.495 0.137 1  Acenaphthylene 23.5 ug/l 0.495 0.111 1  Dimethylphthalate ND ug/l 0.495 0.116 1  2,6-Dinitrotoluene ND ug/l 0.495 0.116 1  Acenaphthene 10.1 ug/l 0.495 0.166 1  Acenaphthene 10.1 ug/l 0.495 0.095 1  3-Nitroaniline ND ug/l 0.495 0.095 1  3-Nitroaniline ND ug/l 0.495 0.095 1	Hexachlorobutadiene	ND		ug/l	0.495	0.085	1
Hexachlorocyclopentadiene ND ug/l 0.495 0.151 1  Biphenyl 6.48 ug/l 0.495 0.110 1  2-Chloronaphthalene ND ug/l 0.495 0.089 1  2-Nitroaniline ND ug/l 0.495 0.137 1  Acenaphthylene 23.5 ug/l 0.495 0.111 1  Dimethylphthalate ND ug/l 0.495 0.116 1  2,6-Dinitrotoluene ND ug/l 0.495 0.166 1  Acenaphthene 10.1 ug/l 0.495 0.095 1  3-Nitroaniline ND ug/l 0.495 0.095 1  3-Nitroaniline ND ug/l 0.495 0.095 1  3-Nitroaniline ND ug/l 0.495 0.110 1  Dibenzofuran 29.7 ug/l 0.495 0.090 1	2-Methylnaphthalene	27.8		ug/l	0.495	0.090	1
Biphenyl   6.48   ug/l   0.495   0.110   1	1,2,4,5-Tetrachlorobenzene	ND		ug/l	0.495	0.079	1
2-Chloronaphthalene ND ug/l 0.495 0.089 1 2-Nitroaniline ND ug/l 0.495 0.137 1 Acenaphthylene 23.5 ug/l 0.495 0.111 1 Dimethylphthalate ND ug/l 0.495 0.116 1 2,6-Dinitrotoluene ND ug/l 0.495 0.166 1 Acenaphthene 10.1 ug/l 0.495 0.095 1 3-Nitroaniline ND ug/l 0.495 0.110 1 Dibenzofuran 29.7 ug/l 0.495 0.090 1	Hexachlorocyclopentadiene	ND		ug/l	0.495	0.151	1
2-Nitroaniline ND ug/l 0.495 0.137 1 Acenaphthylene 23.5 ug/l 0.495 0.111 1 Dimethylphthalate ND ug/l 0.495 0.116 1 2,6-Dinitrotoluene ND ug/l 0.495 0.166 1 Acenaphthene 10.1 ug/l 0.495 0.095 1 3-Nitroaniline ND ug/l 0.495 0.110 1 Dibenzofuran 29.7 ug/l 0.495 0.090 1	Biphenyl	6.48		ug/l	0.495	0.110	1
Acenaphthylene 23.5 ug/l 0.495 0.111 1 Dimethylphthalate ND ug/l 0.495 0.116 1 2,6-Dinitrotoluene ND ug/l 0.495 0.166 1 Acenaphthene 10.1 ug/l 0.495 0.095 1 3-Nitroaniline ND ug/l 0.495 0.110 1 Dibenzofuran 29.7 ug/l 0.495 0.090 1	2-Chloronaphthalene	ND		ug/l	0.495	0.089	1
Dimethylphthalate ND ug/l 0.495 0.116 1 2,6-Dinitrotoluene ND ug/l 0.495 0.166 1 Acenaphthene 10.1 ug/l 0.495 0.095 1 3-Nitroaniline ND ug/l 0.495 0.110 1 Dibenzofuran 29.7 ug/l 0.495 0.090 1	2-Nitroaniline	ND		ug/l	0.495	0.137	1
2,6-Dinitrotoluene ND ug/l 0.495 0.166 1  Acenaphthene 10.1 ug/l 0.495 0.095 1  3-Nitroaniline ND ug/l 0.495 0.110 1  Dibenzofuran 29.7 ug/l 0.495 0.090 1	Acenaphthylene	23.5		ug/l	0.495	0.111	1
Acenaphthene         10.1         ug/l         0.495         0.095         1           3-Nitroaniline         ND         ug/l         0.495         0.110         1           Dibenzofuran         29.7         ug/l         0.495         0.090         1	Dimethylphthalate	ND		ug/l	0.495	0.116	1
3-Nitroaniline ND ug/l 0.495 0.110 1 Dibenzofuran 29.7 ug/l 0.495 0.090 1	2,6-Dinitrotoluene	ND		ug/l	0.495	0.166	1
Dibenzofuran 29.7 ug/l 0.495 0.090 1	Acenaphthene	10.1		ug/l	0.495	0.095	1
	3-Nitroaniline	ND		ug/l	0.495	0.110	1
2,4-Dinitrotoluene ND ug/l 0.495 0.161 1	Dibenzofuran	29.7		ug/l	0.495	0.090	1
<del>-</del>	2,4-Dinitrotoluene	ND		ug/l	0.495	0.161	1



10/11/22

**Project Name:** Lab Number: STEELWINDS ANNUAL/SEMIANNUAL L2249775

**Project Number:** 03.0033579.15

L2249775-02

**SAMPLE RESULTS** 

Date Collected: 09/13/22 09:05

Report Date:

Date Received: Client ID: 09/13/22 MWN-01-091322 Sample Location: Field Prep: LACKAWANNA NY Not Specified

Sample Depth:

Lab ID:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor				
Semivolatile Organics by GC/MS -	Semivolatile Organics by GC/MS - Mansfield Lab									
Fluorene	44.4		ug/l	0.495	0.103	1				
Diethylphthalate	ND		ug/l	0.495	0.178	1				
4-Nitroaniline	ND		ug/l	0.495	0.111	1				
n-Nitrosodiphenylamine	ND		ug/l	0.495	0.071	1				
Hexachlorobenzene	ND		ug/l	0.495	0.121	1				
Phenanthrene	57.0	E	ug/l	0.495	0.110	1				
Anthracene	12.2		ug/l	0.495	0.136	1				
Carbazole	19.7		ug/l	0.495	0.142	1				
Di-n-butylphthalate	ND		ug/l	0.495	0.099	1				
Fluoranthene	12.3		ug/l	0.495	0.154	1				
Pyrene	6.81		ug/l	0.495	0.168	1				
Butylbenzylphthalate	ND		ug/l	0.495	0.084	1				
3,3'-Dichlorobenzidine	ND		ug/l	0.495	0.191	1				
Benz(a)anthracene	0.380	J	ug/l	0.495	0.182	1				
Chrysene	0.214	J	ug/l	0.495	0.140	1				
bis(2-Ethylhexyl)phthalate	ND		ug/l	0.495	0.080	1				
Di-n-octylphthalate	ND		ug/l	0.990	0.078	1				
Benzo(b)fluoranthene	0.079	J	ug/l	0.495	0.065	1				
Benzo(k)fluoranthene	ND		ug/l	0.495	0.159	1				
Benzo(a)pyrene	ND		ug/l	0.495	0.060	1				
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.495	0.089	1				
Dibenz(a,h)anthracene	ND		ug/l	0.495	0.064	1				
Benzo(g,h,i)perylene	ND		ug/l	0.495	0.108	1				

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	63	15-115
Phenol-d5	42	15-115
Nitrobenzene-d5	92	30-130
2-Fluorobiphenyl	78	30-130
2,4,6-Tribromophenol	102	15-115
Terphenyl-d14	92	30-130



10/11/22

Report Date:

Project Name: STEELWINDS ANNUAL/SEMIANNUAL Lab Number: L2249775

**Project Number:** 03.0033579.15

09/20/22 12:54

SAMPLE RESULTS

Lab ID: L2249775-02 D Date Collected: 09/13/22 09:05

Client ID: MWN-01-091322 Date Received: 09/13/22 Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Analytical Date:

Matrix: Water Extraction Method: EPA 3510C
Analytical Method: 1,8270D Extraction Date: 09/17/22 08:30

Analyst: PS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Semivolatile Organics by GC/MS - Mansfield Lab							
Naphthalene	91.9		ug/l	2.48	0.434	5	
Phenanthrene	69.6		ug/l	2.48	0.550	5	

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	62	15-115
Phenol-d5	40	15-115
Nitrobenzene-d5	89	30-130
2-Fluorobiphenyl	85	30-130
2,4,6-Tribromophenol	95	15-115
Terphenyl-d14	91	30-130



10/11/22

**Project Name:** Lab Number: STEELWINDS ANNUAL/SEMIANNUAL L2249775

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Date Collected: 09/13/22 10:15

Report Date:

Lab ID: D L2249775-03 Date Received: Client ID: 09/13/22 MWN-01B-091322 Sample Location: Field Prep: LACKAWANNA NY Not Specified

Sample Depth:

Extraction Method: EPA 3510C Matrix: Water **Extraction Date:** 09/17/22 08:30 Analytical Method: 1,8270D

Analytical Date: 09/20/22 13:24 Analyst: PS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS - I	Mansfield Lab					
bis(2-Chloroethyl)ether	ND		ug/l	9.80	1.82	20
1,3-Dichlorobenzene	ND		ug/l	9.80	1.54	20
1,4-Dichlorobenzene	ND		ug/l	9.80	1.62	20
1,2-Dichlorobenzene	ND		ug/l	9.80	1.33	20
Benzyl alcohol	ND		ug/l	9.80	2.41	20
bis(2-chloroisopropyl)ether	ND		ug/l	9.80	2.12	20
Acetophenone	ND		ug/l	19.6	4.06	20
Hexachloroethane	ND		ug/l	9.80	2.00	20
Nitrobenzene	ND		ug/l	9.80	2.00	20
Isophorone	ND		ug/l	9.80	2.47	20
bis(2-Chloroethoxy)methane	ND		ug/l	9.80	1.67	20
1,2,4-Trichlorobenzene	ND		ug/l	9.80	1.88	20
Naphthalene	742		ug/l	9.80	1.72	20
4-Chloroaniline	ND		ug/l	9.80	2.51	20
Hexachlorobutadiene	ND		ug/l	9.80	1.68	20
2-Methylnaphthalene	52.4		ug/l	9.80	1.79	20
1,2,4,5-Tetrachlorobenzene	ND		ug/l	9.80	1.56	20
Hexachlorocyclopentadiene	ND		ug/l	9.80	3.00	20
Biphenyl	7.84	J	ug/l	9.80	2.18	20
2-Chloronaphthalene	ND		ug/l	9.80	1.76	20
2-Nitroaniline	ND		ug/l	9.80	2.70	20
Acenaphthylene	54.3		ug/l	9.80	2.20	20
Dimethylphthalate	ND		ug/l	9.80	2.29	20
2,6-Dinitrotoluene	ND		ug/l	9.80	3.29	20
Acenaphthene	11.8		ug/l	9.80	1.87	20
3-Nitroaniline	ND		ug/l	9.80	2.18	20
Dibenzofuran	30.6		ug/l	9.80	1.78	20
2,4-Dinitrotoluene	ND		ug/l	9.80	3.20	20



10/11/22

**Project Name:** Lab Number: STEELWINDS ANNUAL/SEMIANNUAL L2249775

**Project Number:** 03.0033579.15

L2249775-03

**SAMPLE RESULTS** 

Date Collected: 09/13/22 10:15

Report Date:

Date Received: Client ID: 09/13/22 MWN-01B-091322 Sample Location: LACKAWANNA NY

D

Field Prep: Not Specified

Sample Depth:

Lab ID:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor			
Semivolatile Organics by GC/MS - Mansfield Lab									
Fluorene	42.3		ug/l	9.80	2.04	20			
Diethylphthalate	ND		ug/l	9.80	3.53	20			
4-Nitroaniline	ND		ug/l	9.80	2.20	20			
n-Nitrosodiphenylamine	ND		ug/l	9.80	1.41	20			
Hexachlorobenzene	ND		ug/l	9.80	2.39	20			
Phenanthrene	69.5		ug/l	9.80	2.18	20			
Anthracene	11.8		ug/l	9.80	2.69	20			
Carbazole	61.3		ug/l	9.80	2.80	20			
Di-n-butylphthalate	ND		ug/l	9.80	1.95	20			
Fluoranthene	10.8		ug/l	9.80	3.06	20			
Pyrene	5.57	J	ug/l	9.80	3.33	20			
Butylbenzylphthalate	ND		ug/l	9.80	1.66	20			
3,3'-Dichlorobenzidine	ND		ug/l	9.80	3.78	20			
Benz(a)anthracene	ND		ug/l	9.80	3.61	20			
Chrysene	ND		ug/l	9.80	2.78	20			
bis(2-Ethylhexyl)phthalate	ND		ug/l	9.80	1.59	20			
Di-n-octylphthalate	ND		ug/l	19.6	1.54	20			
Benzo(b)fluoranthene	ND		ug/l	9.80	1.28	20			
Benzo(k)fluoranthene	ND		ug/l	9.80	3.16	20			
Benzo(a)pyrene	ND		ug/l	9.80	1.18	20			
Indeno(1,2,3-cd)pyrene	ND		ug/l	9.80	1.76	20			
Dibenz(a,h)anthracene	ND		ug/l	9.80	1.26	20			
Benzo(g,h,i)perylene	ND		ug/l	9.80	2.14	20			

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	58	15-115
Phenol-d5	37	15-115
Nitrobenzene-d5	85	30-130
2-Fluorobiphenyl	84	30-130
2,4,6-Tribromophenol	83	15-115
Terphenyl-d14	89	30-130



L2249775

10/11/22

Project Name: STEELWINDS ANNUAL/SEMIANNUAL Lab Number:

**Project Number:** 03.0033579.15

L2249775-04

**SAMPLE RESULTS** 

Date Collected: 09/13/22 10:55

Report Date:

Client ID: WT1-04-091322 Date Received: 09/13/22 Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Lab ID:

Matrix: Water Extraction Method: EPA 3510C
Analytical Method: 1,8270D Extraction Date: 09/17/22 08:30

Analytical Date: 09/19/22 13:39

Analyst: PS

Semivolatile Organics by GC/MS - Mansfield bis(2-Chloroethyl)ether 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene Benzyl alcohol	ND ND ND ND ND		ug/l ug/l	0.495	0.092	4
1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene	ND ND				0.092	4
1,4-Dichlorobenzene 1,2-Dichlorobenzene	ND		ug/l			1
1,2-Dichlorobenzene				0.495	0.078	1
·	ND		ug/l	0.495	0.082	1
Ponzul alaahal			ug/l	0.495	0.067	1
Delizyi alconol	ND		ug/l	0.495	0.122	1
bis(2-chloroisopropyl)ether	ND		ug/l	0.495	0.107	1
Acetophenone	0.413	J	ug/l	0.990	0.205	1
Hexachloroethane	ND		ug/l	0.495	0.101	1
Nitrobenzene	ND		ug/l	0.495	0.101	1
Isophorone	ND		ug/l	0.495	0.125	1
bis(2-Chloroethoxy)methane	ND		ug/l	0.495	0.085	1
1,2,4-Trichlorobenzene	ND		ug/l	0.495	0.095	1
Naphthalene	32.6		ug/l	0.495	0.087	1
4-Chloroaniline	ND		ug/l	0.495	0.127	1
Hexachlorobutadiene	ND		ug/l	0.495	0.085	1
2-Methylnaphthalene	8.39		ug/l	0.495	0.090	1
1,2,4,5-Tetrachlorobenzene	ND		ug/l	0.495	0.079	1
Hexachlorocyclopentadiene	ND		ug/l	0.495	0.151	1
Biphenyl	1.96		ug/l	0.495	0.110	1
2-Chloronaphthalene	ND		ug/l	0.495	0.089	1
2-Nitroaniline	ND		ug/l	0.495	0.137	1
Acenaphthylene	3.24		ug/l	0.495	0.111	1
Dimethylphthalate	ND		ug/l	0.495	0.116	1
2,6-Dinitrotoluene	ND		ug/l	0.495	0.166	1
Acenaphthene	3.42		ug/l	0.495	0.095	1
3-Nitroaniline	ND		ug/l	0.495	0.110	1
Dibenzofuran	10.1		ug/l	0.495	0.090	1
2,4-Dinitrotoluene	ND		ug/l	0.495	0.161	1



10/11/22

**Project Name:** Lab Number: STEELWINDS ANNUAL/SEMIANNUAL L2249775

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Date Collected: 09/13/22 10:55

Report Date:

L2249775-04 Date Received: Client ID: 09/13/22 WT1-04-091322 Sample Location: LACKAWANNA NY

Field Prep: Not Specified

Sample Depth:

Lab ID:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS -	Mansfield Lab					
Fluorene	15.2		ug/l	0.495	0.103	1
Diethylphthalate	ND		ug/l	0.495	0.178	1
4-Nitroaniline	ND		ug/l	0.495	0.111	1
n-Nitrosodiphenylamine	ND		ug/l	0.495	0.071	1
Hexachlorobenzene	ND		ug/l	0.495	0.121	1
Phenanthrene	36.3		ug/l	0.495	0.110	1
Anthracene	7.70		ug/l	0.495	0.136	1
Carbazole	8.48		ug/l	0.495	0.142	1
Di-n-butylphthalate	ND		ug/l	0.495	0.099	1
Fluoranthene	10.9		ug/l	0.495	0.154	1
Pyrene	6.39		ug/l	0.495	0.168	1
Butylbenzylphthalate	ND		ug/l	0.495	0.084	1
3,3'-Dichlorobenzidine	ND		ug/l	0.495	0.191	1
Benz(a)anthracene	0.342	J	ug/l	0.495	0.182	1
Chrysene	0.250	J	ug/l	0.495	0.140	1
bis(2-Ethylhexyl)phthalate	ND		ug/l	0.495	0.080	1
Di-n-octylphthalate	ND		ug/l	0.990	0.078	1
Benzo(b)fluoranthene	0.093	J	ug/l	0.495	0.065	1
Benzo(k)fluoranthene	ND		ug/l	0.495	0.159	1
Benzo(a)pyrene	ND		ug/l	0.495	0.060	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.495	0.089	1
Dibenz(a,h)anthracene	ND		ug/l	0.495	0.064	1
Benzo(g,h,i)perylene	ND		ug/l	0.495	0.108	1

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	62	15-115
Phenol-d5	43	15-115
Nitrobenzene-d5	88	30-130
2-Fluorobiphenyl	79	30-130
2,4,6-Tribromophenol	106	15-115
Terphenyl-d14	97	30-130



L2249775

10/11/22

Project Name: STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

L2249775-05

BCP-ORC-1-091322

LACKAWANNA NY

**SAMPLE RESULTS** 

Date Collected: 09/13/22 13:25

Date Received: 09/13/22

Lab Number:

Report Date:

Field Prep: Not Specified

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Matrix: Water
Analytical Method: 1,8270D
Analytical Date: 09/19/22 14:09

Analyst: PS

Extraction Method: EPA 3510C Extraction Date: 09/17/22 08:30

1,3-Dichlorobenzene         ND         ug/l         0.490         0.077         1           1,4-Dichlorobenzene         ND         ug/l         0.490         0.081         1           1,2-Dichlorobenzene         ND         ug/l         0.490         0.067         1           Benzyl alcohol         ND         ug/l         0.490         0.120         1           Benzyl alcohol         ND         ug/l         0.490         0.120         1           Acetophenone         0.492         J         ug/l         0.490         0.100         1           Hexachloroethane         ND         ug/l         0.490         0.100         1           Nitrobenzene         ND         ug/l         0.490         0.100         1           Isophorone         ND         ug/l         0.490         0.100         1           Isophorone         ND         ug/l         0.490         0.084         1           ND         ug/l         0.490         0.084         1           1,2,4-Trichlorobenzene         ND         ug/l         0.490         0.086         1           4-Chloroshiline         ND         ug/l         0.490         0.084         1 </th <th>Parameter</th> <th>Result</th> <th>Qualifier</th> <th>Units</th> <th>RL</th> <th>MDL</th> <th>Dilution Factor</th>	Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
1,3-Dichlorobenzene   ND   ug/l   0,490   0,077   1   1   1,4-Dichlorobenzene   ND   ug/l   0,490   0,081   1   1   1,2-Dichlorobenzene   ND   ug/l   0,490   0,067   1   1   1,2-Dichlorobenzene   ND   ug/l   0,490   0,067   1   1   1   1   1   1   1   1   1	Semivolatile Organics by GC/MS - Ma	ansfield Lab					
1,4-Dichlorobenzene   ND   ug/l   0,490   0,081   1   1,2-Dichlorobenzene   ND   ug/l   0,490   0,067   1   1   1,2-Dichlorobenzene   ND   ug/l   0,490   0,120   1   1   1   1   1   1   1   1   1	bis(2-Chloroethyl)ether	ND		ug/l	0.490	0.091	1
1,2-Dichlorobenzene   ND   ug/l   0.490   0.067   1	1,3-Dichlorobenzene	ND		ug/l	0.490	0.077	1
Benzyl alcohol         ND         ug/l         0.490         0.120         1           bis(2-chloroisopropyl)ether         ND         ug/l         0.490         0.106         1           Acetophenone         0.492         J         ug/l         0.980         0.203         1           Hexachloroethane         ND         ug/l         0.490         0.100         1           Nitrobenzene         ND         ug/l         0.490         0.100         1           Isisphorone         ND         ug/l         0.490         0.100         1           Isisphorone         ND         ug/l         0.490         0.084         1           Isisphorone         ND         ug/l         0.490         0.086         1           Isisphorone         ND         ug/l         0.490         0.125         1           Isisphorone         ND         ug/l         0.490         0.086 <t< td=""><td>1,4-Dichlorobenzene</td><td>ND</td><td></td><td>ug/l</td><td>0.490</td><td>0.081</td><td>1</td></t<>	1,4-Dichlorobenzene	ND		ug/l	0.490	0.081	1
Dis(2-chloroispropy))ether   ND	1,2-Dichlorobenzene	ND		ug/l	0.490	0.067	1
Acetophenone 0.492 J ug/l 0.980 0.203 1 Hexachloroethane ND ug/l 0.490 0.100 1 Isophorone ND ug/l 0.490 0.100 1 Isophorone ND ug/l 0.490 0.100 1 Isophorone ND ug/l 0.490 0.124 1 Isophorone ND ug/l 0.490 0.084 1 I.2.4-Trichlorobenzene ND ug/l 0.490 0.084 1 I.2.4-Trichlorobenzene ND ug/l 0.490 0.094 1 I.2.4-Trichlorobenzene ND ug/l 0.490 0.086 1 I.2.4-Trichlorobenzene ND ug/l 0.490 0.084 1 I.2.4-Trichlorobenzene ND ug/l 0.490 0.084 1 I.2.4-Trichlorobenzene ND ug/l 0.490 0.089 1 I.2.4-S-Tetrachlorobenzene ND ug/l 0.490 0.089 1 I.2.4-S-Tetrachlorobenzene ND ug/l 0.490 0.078 1 I.2.4-S-Tetrachlorobenzene ND ug/l 0.490 0.150 1 III Dibenoropy 0.150 1 II Dibenoropy 0.150 1 II Dimethylphthalate ND ug/l 0.490 0.195 1 II Dimethylphthalate ND ug/l 0.490 0.115 1 II Dimethylphthalate ND ug/l 0.490 0.109 1	Benzyl alcohol	ND		ug/l	0.490	0.120	1
ND	bis(2-chloroisopropyl)ether	ND		ug/l	0.490	0.106	1
Nitrobenzene ND ug/l 0.490 0.100 1 Isophorone ND ug/l 0.490 0.100 1 Isophorone ND ug/l 0.490 0.124 1 Ibis(2-Chloroethoxy)methane ND ug/l 0.490 0.084 1 I.2,4-Trichlorobenzene ND ug/l 0.490 0.094 1 Naphthalene 90.4 E ug/l 0.490 0.086 1 I-Chloroaniline ND ug/l 0.490 0.125 1 I-Chloroaniline ND ug/l 0.490 0.086 1 I-Chloroaniline ND ug/l 0.490 0.086 1 I-Chloroaniline ND ug/l 0.490 0.086 1 I-Chloroaniline ND ug/l 0.490 0.084 1 I-Chloroaniline ND ug/l 0.490 0.084 1 I-Chloroaniline ND ug/l 0.490 0.089 1 I-Chlorophyloaniline ND ug/l 0.490 0.089 1 I-Chloronaphthalene ND ug/l 0.490 0.078 1 I-Chloronaphthalene ND ug/l 0.490 0.150 1 I-Chloronaphthalene ND ug/l 0.490 0.150 1 I-Chloronaphthalene ND ug/l 0.490 0.109 1 I-Chloronaphthalene ND ug/l 0.490 0.135 1 I-Chloronaphthalene ND ug/l 0.490 0.115 1 I-Chlorophylone ND ug/l 0.490 0.190 1 I-Chlorophylone ND ug/l 0.490 0.190 1 I-Chlorophylone ND ug/l 0.490 0.094 1 I-Chlorophylone ND ug/l 0.490 0.099 1	Acetophenone	0.492	J	ug/l	0.980	0.203	1
ND	Hexachloroethane	ND		ug/l	0.490	0.100	1
bis(2-Chloroethoxy)methane         ND         ug/l         0.490         0.084         1           1,2,4-Trichlorobenzene         ND         ug/l         0.490         0.094         1           Naphthalene         90.4         E         ug/l         0.490         0.086         1           4-Chloroaniline         ND         ug/l         0.490         0.125         1           Hexachlorobutadiene         ND         ug/l         0.490         0.084         1           2-Methylnaphthalene         23.2         ug/l         0.490         0.089         1           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.490         0.078         1           Hexachlorocyclopentadiene         ND         ug/l         0.490         0.150         1           Biphenyl         3.06         ug/l         0.490         0.109         1           2-Chloronaphthalene         ND         ug/l         0.490         0.135         1           2-Nitroaniline         ND         ug/l         0.490         0.135         1           Acenaphthylene         17.0         ug/l         0.490         0.115         1           2,6-Dinitrotoluene         ND	Nitrobenzene	ND		ug/l	0.490	0.100	1
1,2,4-Trichlorobenzene   ND   ug/l   0.490   0.094   1	Isophorone	ND		ug/l	0.490	0.124	1
Naphthalene         90.4         E         ug/l         0.490         0.086         1           4-Chloroaniline         ND         ug/l         0.490         0.125         1           Hexachlorobutadiene         ND         ug/l         0.490         0.084         1           2-Methylnaphthalene         23.2         ug/l         0.490         0.089         1           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.490         0.078         1           Hexachlorocyclopentadiene         ND         ug/l         0.490         0.150         1           Biphenyl         3.06         ug/l         0.490         0.109         1           2-Chloronaphthalene         ND         ug/l         0.490         0.088         1           2-Nittroaniline         ND         ug/l         0.490         0.135         1           Acenaphthylene         17.0         ug/l         0.490         0.110         1           Dimethylphthalate         ND         ug/l         0.490         0.115         1           2,6-Dinitrotoluene         ND         ug/l         0.490         0.165         1           Acenaphthene         5.68         ug/l	bis(2-Chloroethoxy)methane	ND		ug/l	0.490	0.084	1
A-Chloroaniline ND ug/l 0.490 0.125 1  Hexachlorobutadiene ND ug/l 0.490 0.084 1  2-Methylnaphthalene 23.2 ug/l 0.490 0.089 1  1,2,4,5-Tetrachlorobenzene ND ug/l 0.490 0.078 1  Hexachlorocyclopentadiene ND ug/l 0.490 0.150 1  Biphenyl 3.06 ug/l 0.490 0.150 1  Biphenyl 3.06 ug/l 0.490 0.109 1  2-Chloronaphthalene ND ug/l 0.490 0.109 1  2-Chloronaphthalene ND ug/l 0.490 0.135 1  Acenaphthylene 17.0 ug/l 0.490 0.110 1  Dimethylphthalate ND ug/l 0.490 0.115 1  2,6-Dinitrotoluene ND ug/l 0.490 0.115 1  2,6-Dinitrotoluene ND ug/l 0.490 0.115 1  3-Nitroaniline ND ug/l 0.490 0.165 1  Acenaphthene 5.68 ug/l 0.490 0.094 1  3-Nitroaniline ND ug/l 0.490 0.099 1	1,2,4-Trichlorobenzene	ND		ug/l	0.490	0.094	1
Hexachlorobutadiene   ND	Naphthalene	90.4	E	ug/l	0.490	0.086	1
2-Methylnaphthalene       23.2       ug/l       0.490       0.089       1         1,2,4,5-Tetrachlorobenzene       ND       ug/l       0.490       0.078       1         Hexachlorocyclopentadiene       ND       ug/l       0.490       0.150       1         Biphenyl       3.06       ug/l       0.490       0.109       1         2-Chloronaphthalene       ND       ug/l       0.490       0.088       1         2-Nitroaniline       ND       ug/l       0.490       0.135       1         Acenaphthylene       17.0       ug/l       0.490       0.110       1         Dimethylphthalate       ND       ug/l       0.490       0.115       1         2,6-Dinitrotoluene       ND       ug/l       0.490       0.165       1         Acenaphthene       5.68       ug/l       0.490       0.094       1         3-Nitroaniline       ND       ug/l       0.490       0.109       1         Dibenzofuran       13.8       ug/l       0.490       0.089       1	4-Chloroaniline	ND		ug/l	0.490	0.125	1
1,2,4,5-Tetrachlorobenzene       ND       ug/l       0.490       0.078       1         Hexachlorocyclopentadiene       ND       ug/l       0.490       0.150       1         Biphenyl       3.06       ug/l       0.490       0.109       1         2-Chloronaphthalene       ND       ug/l       0.490       0.088       1         2-Nitroaniline       ND       ug/l       0.490       0.135       1         Acenaphthylene       17.0       ug/l       0.490       0.110       1         Dimethylphthalate       ND       ug/l       0.490       0.115       1         2,6-Dinitrotoluene       ND       ug/l       0.490       0.165       1         Acenaphthene       5.68       ug/l       0.490       0.094       1         3-Nitroaniline       ND       ug/l       0.490       0.109       1         Dibenzofuran       13.8       ug/l       0.490       0.089       1	Hexachlorobutadiene	ND		ug/l	0.490	0.084	1
Hexachlorocyclopentadiene   ND   ug/l   0.490   0.150   1	2-Methylnaphthalene	23.2		ug/l	0.490	0.089	1
Biphenyl   3.06   ug/l   0.490   0.109   1	1,2,4,5-Tetrachlorobenzene	ND		ug/l	0.490	0.078	1
2-Chloronaphthalene ND ug/l 0.490 0.088 1 2-Nitroaniline ND ug/l 0.490 0.135 1 Acenaphthylene 17.0 ug/l 0.490 0.110 1 Dimethylphthalate ND ug/l 0.490 0.115 1 2,6-Dinitrotoluene ND ug/l 0.490 0.165 1 Acenaphthene 5.68 ug/l 0.490 0.094 1 3-Nitroaniline ND ug/l 0.490 0.109 1 Dibenzofuran 13.8 ug/l 0.490 0.089 1	Hexachlorocyclopentadiene	ND		ug/l	0.490	0.150	1
2-Nitroaniline ND ug/l 0.490 0.135 1 Acenaphthylene 17.0 ug/l 0.490 0.110 1 Dimethylphthalate ND ug/l 0.490 0.115 1 2,6-Dinitrotoluene ND ug/l 0.490 0.165 1 Acenaphthene 5.68 ug/l 0.490 0.094 1 3-Nitroaniline ND ug/l 0.490 0.109 1 Dibenzofuran 13.8 ug/l 0.490 0.089 1	Biphenyl	3.06		ug/l	0.490	0.109	1
Acenaphthylene 17.0 ug/l 0.490 0.110 1  Dimethylphthalate ND ug/l 0.490 0.115 1  2,6-Dinitrotoluene ND ug/l 0.490 0.165 1  Acenaphthene 5.68 ug/l 0.490 0.094 1  3-Nitroaniline ND ug/l 0.490 0.109 1  Dibenzofuran 13.8 ug/l 0.490 0.089 1	2-Chloronaphthalene	ND		ug/l	0.490	0.088	1
Dimethylphthalate	2-Nitroaniline	ND		ug/l	0.490	0.135	1
2,6-Dinitrotoluene ND ug/l 0.490 0.165 1  Acenaphthene 5.68 ug/l 0.490 0.094 1  3-Nitroaniline ND ug/l 0.490 0.109 1  Dibenzofuran 13.8 ug/l 0.490 0.089 1	Acenaphthylene	17.0		ug/l	0.490	0.110	1
Acenaphthene         5.68         ug/l         0.490         0.094         1           3-Nitroaniline         ND         ug/l         0.490         0.109         1           Dibenzofuran         13.8         ug/l         0.490         0.089         1	Dimethylphthalate	ND		ug/l	0.490	0.115	1
3-Nitroaniline ND ug/l 0.490 0.109 1 Dibenzofuran 13.8 ug/l 0.490 0.089 1	2,6-Dinitrotoluene	ND		ug/l	0.490	0.165	1
Dibenzofuran 13.8 ug/l 0.490 0.089 1	Acenaphthene	5.68		ug/l	0.490	0.094	1
-9.	3-Nitroaniline	ND		ug/l	0.490	0.109	1
2,4-Dinitrotoluene ND ug/l 0.490 0.160 1	Dibenzofuran	13.8		ug/l	0.490	0.089	1
	2,4-Dinitrotoluene	ND		ug/l	0.490	0.160	1



10/11/22

**Project Name:** Lab Number: STEELWINDS ANNUAL/SEMIANNUAL L2249775

**Project Number:** 03.0033579.15

L2249775-05

**SAMPLE RESULTS** 

Date Collected: 09/13/22 13:25

Report Date:

BCP-ORC-1-091322 Date Received: Client ID: 09/13/22 Sample Location: Field Prep: LACKAWANNA NY Not Specified

Sample Depth:

Lab ID:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS	- Mansfield Lab					
Fluorene	21.4		ug/l	0.490	0.102	1
Diethylphthalate	ND		ug/l	0.490	0.176	1
4-Nitroaniline	ND		ug/l	0.490	0.110	1
n-Nitrosodiphenylamine	ND		ug/l	0.490	0.071	1
Hexachlorobenzene	ND		ug/l	0.490	0.120	1
Phenanthrene	30.0		ug/l	0.490	0.109	1
Anthracene	3.76		ug/l	0.490	0.134	1
Carbazole	26.2		ug/l	0.490	0.140	1
Di-n-butylphthalate	ND		ug/l	0.490	0.098	1
Fluoranthene	6.32		ug/l	0.490	0.153	1
Pyrene	4.06		ug/l	0.490	0.167	1
Butylbenzylphthalate	ND		ug/l	0.490	0.083	1
3,3'-Dichlorobenzidine	ND		ug/l	0.490	0.189	1
Benz(a)anthracene	0.214	J	ug/l	0.490	0.180	1
Chrysene	0.145	J	ug/l	0.490	0.139	1
bis(2-Ethylhexyl)phthalate	ND		ug/l	0.490	0.079	1
Di-n-octylphthalate	ND		ug/l	0.980	0.077	1
Benzo(b)fluoranthene	ND		ug/l	0.490	0.064	1
Benzo(k)fluoranthene	ND		ug/l	0.490	0.158	1
Benzo(a)pyrene	ND		ug/l	0.490	0.059	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.490	0.088	1
Dibenz(a,h)anthracene	ND		ug/l	0.490	0.063	1
Benzo(g,h,i)perylene	ND		ug/l	0.490	0.107	1

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	38	15-115
Phenol-d5	24	15-115
Nitrobenzene-d5	83	30-130
2-Fluorobiphenyl	67	30-130
2,4,6-Tribromophenol	88	15-115
Terphenyl-d14	84	30-130



10/11/22

Report Date:

Project Name: STEELWINDS ANNUAL/SEMIANNUAL Lab Number: L2249775

**Project Number:** 03.0033579.15

09/20/22 13:54

SAMPLE RESULTS

Lab ID: L2249775-05 D Date Collected: 09/13/22 13:25

Client ID: BCP-ORC-1-091322 Date Received: 09/13/22 Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Analytical Date:

Matrix: Water Extraction Method: EPA 3510C
Analytical Method: 1,8270D Extraction Date: 09/17/22 08:30

Analyst: PS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS - Mansfield	d Lab					
Naphthalene	198		ug/l	4.90	0.859	10

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	37	15-115
Phenol-d5	22	15-115
Nitrobenzene-d5	75	30-130
2-Fluorobiphenyl	75	30-130
2,4,6-Tribromophenol	78	15-115
Terphenyl-d14	86	30-130



10/11/22

Project Name: STEELWINDS ANNUAL/SEMIANNUAL Lab Number: L2249775

**Project Number:** 03.0033579.15

L2249775-06

09/19/22 14:38

SAMPLE RESULTS

Date Collected: 09/13/22 14:15

Report Date:

Client ID: WT1-02-091322 Date Received: 09/13/22 Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Analytical Date:

Lab ID:

Matrix: Water Extraction Method: EPA 3510C
Analytical Method: 1,8270D Extraction Date: 09/17/22 08:30

Analyst: PS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Semivolatile Organics by GC/MS -	Mansfield Lab						
bis(2-Chloroethyl)ether	ND		ug/l	0.495	0.092	1	
1,3-Dichlorobenzene	ND		ug/l	0.495	0.078	1	
1,4-Dichlorobenzene	ND		ug/l	0.495	0.082	1	
1,2-Dichlorobenzene	ND		ug/l	0.495	0.067	1	
Benzyl alcohol	ND		ug/l	0.495	0.122	1	
bis(2-chloroisopropyl)ether	ND		ug/l	0.495	0.107	1	
Acetophenone	0.317	J	ug/l	0.990	0.205	1	
Hexachloroethane	ND		ug/l	0.495	0.101	1	
Nitrobenzene	ND		ug/l	0.495	0.101	1	
Isophorone	ND		ug/l	0.495	0.125	1	
bis(2-Chloroethoxy)methane	ND		ug/l	0.495	0.085	1	
1,2,4-Trichlorobenzene	ND		ug/l	0.495	0.095	1	
Naphthalene	17.2		ug/l	0.495	0.087	1	
4-Chloroaniline	ND		ug/l	0.495	0.127	1	
Hexachlorobutadiene	ND		ug/l	0.495	0.085	1	
2-Methylnaphthalene	4.62		ug/l	0.495	0.090	1	
1,2,4,5-Tetrachlorobenzene	ND		ug/l	0.495	0.079	1	
Hexachlorocyclopentadiene	ND		ug/l	0.495	0.151	1	
Biphenyl	1.13		ug/l	0.495	0.110	1	
2-Chloronaphthalene	ND		ug/l	0.495	0.089	1	
2-Nitroaniline	ND		ug/l	0.495	0.137	1	
Acenaphthylene	1.16		ug/l	0.495	0.111	1	
Dimethylphthalate	ND		ug/l	0.495	0.116	1	
2,6-Dinitrotoluene	ND		ug/l	0.495	0.166	1	
Acenaphthene	1.47		ug/l	0.495	0.095	1	
3-Nitroaniline	ND		ug/l	0.495	0.110	1	
Dibenzofuran	4.92		ug/l	0.495	0.090	1	
2,4-Dinitrotoluene	ND		ug/l	0.495	0.161	1	



10/11/22

Project Name: STEELWINDS ANNUAL/SEMIANNUAL Lab Number: L2249775

**Project Number:** 03.0033579.15

L2249775-06

**SAMPLE RESULTS** 

Date Collected: 09/13/22 14:15

Report Date:

Client ID: WT1-02-091322 Date Received: 09/13/22 Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Lab ID:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS - Mansfie	eld Lab					
Fluorene	7.48		ug/l	0.495	0.103	1
Diethylphthalate	ND		ug/l	0.495	0.178	1
4-Nitroaniline	ND		ug/l	0.495	0.111	1
n-Nitrosodiphenylamine	ND		ug/l	0.495	0.071	1
Hexachlorobenzene	ND		ug/l	0.495	0.121	1
Phenanthrene	13.7		ug/l	0.495	0.110	1
Anthracene	2.74		ug/l	0.495	0.136	1
Carbazole	6.02		ug/l	0.495	0.142	1
Di-n-butylphthalate	ND		ug/l	0.495	0.099	1
Fluoranthene	4.61		ug/l	0.495	0.154	1
Pyrene	2.93		ug/l	0.495	0.168	1
Butylbenzylphthalate	ND		ug/l	0.495	0.084	1
3,3'-Dichlorobenzidine	ND		ug/l	0.495	0.191	1
Benz(a)anthracene	ND		ug/l	0.495	0.182	1
Chrysene	ND		ug/l	0.495	0.140	1
bis(2-Ethylhexyl)phthalate	ND		ug/l	0.495	0.080	1
Di-n-octylphthalate	ND		ug/l	0.990	0.078	1
Benzo(b)fluoranthene	ND		ug/l	0.495	0.065	1
Benzo(k)fluoranthene	ND		ug/l	0.495	0.159	1
Benzo(a)pyrene	ND		ug/l	0.495	0.060	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.495	0.089	1
Dibenz(a,h)anthracene	ND		ug/l	0.495	0.064	1
Benzo(g,h,i)perylene	ND		ug/l	0.495	0.108	1

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	41	15-115
Phenol-d5	27	15-115
Nitrobenzene-d5	81	30-130
2-Fluorobiphenyl	75	30-130
2,4,6-Tribromophenol	103	15-115
Terphenyl-d14	92	30-130



L2249775

Lab Number:

Project Name: STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15 **Report Date:** 10/11/22

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8270D Extraction Method: EPA 3510C Analytical Date: 09/19/22 10:43 Extraction Date: 09/17/22 08:30

Analyst: PS

Parameter	Result	Qualifier Ur	nits		RL	MDL	
Semivolatile Organics by GC/MS	s - Mansfield La	ab for sample(s	s):	01-06	Batch:	WG1688692-1	
bis(2-Chloroethyl)ether	ND	ι	ıg/l	(	0.500	0.093	
1,3-Dichlorobenzene	ND	ι	ıg/l	(	0.500	0.078	
1,4-Dichlorobenzene	ND	ι	ıg/l	(	0.500	0.083	
1,2-Dichlorobenzene	ND	ι	ıg/l	(	0.500	0.068	
Benzyl alcohol	ND	ι	ıg/l	(	0.500	0.123	
bis(2-chloroisopropyl)ether	ND	ι	ıg/l	(	0.500	0.108	
Acetophenone	ND	ι	ıg/l		1.00	0.207	
Hexachloroethane	ND	ι	ıg/l	(	0.500	0.102	
Nitrobenzene	ND	ι	ıg/l		0.500	0.102	
Isophorone	ND	ι	ıg/l		0.500	0.126	
bis(2-Chloroethoxy)methane	ND	ι	ıg/l		0.500	0.085	
1,2,4-Trichlorobenzene	ND	ι	ıg/l		0.500	0.096	
Naphthalene	ND	ι	ıg/l		0.500	0.088	
4-Chloroaniline	ND	ι	ıg/l	(	0.500	0.128	
Hexachlorobutadiene	ND	ι	ıg/l		0.500	0.086	
2-Methylnaphthalene	ND	ι	ıg/l		0.500	0.091	
1,2,4,5-Tetrachlorobenzene	ND	ι	ıg/l		0.500	0.080	
Hexachlorocyclopentadiene	ND	ι	ıg/l	(	0.500	0.153	
Biphenyl	ND	ι	ıg/l	(	0.500	0.111	
2-Chloronaphthalene	ND	ι	ıg/l	(	0.500	0.090	
2-Nitroaniline	ND	ι	ıg/l	(	0.500	0.138	
Acenaphthylene	ND	ι	ıg/l	(	0.500	0.112	
Dimethylphthalate	ND	ι	ıg/l	(	0.500	0.117	
2,6-Dinitrotoluene	ND	ι	ıg/l		0.500	0.168	
Acenaphthene	ND	ι	ıg/l	(	0.500	0.096	
3-Nitroaniline	ND	ι	ıg/l	(	0.500	0.111	
Dibenzofuran	ND	ι	ıg/l	(	0.500	0.091	
2,4-Dinitrotoluene	ND	ι	ıg/l	(	0.500	0.163	
Fluorene	ND	U	ıg/l		0.500	0.104	



L2249775

Project Name: STEELWINDS ANNUAL/SEMIANNUAL Lab Number:

**Project Number:** 03.0033579.15 **Report Date:** 10/11/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270D Extraction Method: EPA 3510C
Analytical Date: 09/19/22 10:43 Extraction Date: 09/17/22 08:30

Analyst: PS

Parameter	Result	Qualifier	Units		RL	MDL	
Semivolatile Organics by GC/MS - I	Mansfield La	b for samp	ole(s):	01-06	Batch:	WG1688692-1	
Diethylphthalate	ND		ug/l		0.500	0.180	
4-Nitroaniline	ND		ug/l		0.500	0.112	
n-Nitrosodiphenylamine	ND		ug/l		0.500	0.072	
Hexachlorobenzene	ND		ug/l		0.500	0.122	
Phenanthrene	ND		ug/l		0.500	0.111	
Anthracene	ND		ug/l		0.500	0.137	
Carbazole	ND		ug/l		0.500	0.143	
Di-n-butylphthalate	ND		ug/l		0.500	0.100	
Fluoranthene	ND		ug/l		0.500	0.156	
Pyrene	ND		ug/l		0.500	0.170	
Butylbenzylphthalate	ND		ug/l		0.500	0.085	
3,3'-Dichlorobenzidine	ND		ug/l		0.500	0.193	
Benz(a)anthracene	ND		ug/l		0.500	0.184	
Chrysene	ND		ug/l		0.500	0.142	
bis(2-Ethylhexyl)phthalate	ND		ug/l		0.500	0.081	
Di-n-octylphthalate	ND		ug/l		1.00	0.079	
Benzo(b)fluoranthene	ND		ug/l		0.500	0.066	
Benzo(k)fluoranthene	ND		ug/l		0.500	0.161	
Benzo(a)pyrene	ND		ug/l		0.500	0.060	
Indeno(1,2,3-cd)pyrene	ND		ug/l		0.500	0.090	
Dibenz(a,h)anthracene	ND		ug/l		0.500	0.064	
Benzo(g,h,i)perylene	ND		ug/l		0.500	0.109	



L2249775

Lab Number:

**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

**Project Number: Report Date:** 03.0033579.15 10/11/22

**Method Blank Analysis** 

**Batch Quality Control** 

Analytical Method: 1,8270D Extraction Method: EPA 3510C Analytical Date: 09/19/22 10:43 09/17/22 08:30 **Extraction Date:** 

Analyst: PS

> Result Qualifier Units RLMDL **Parameter**

Semivolatile Organics by GC/MS - Mansfield Lab for sample(s): 01-06 Batch: WG1688692-1

Surrogate	%Recovery Qualifi	Acceptance er Criteria
	•	
2-Fluorophenol	64	15-115
Phenol-d5	42	15-115
Nitrobenzene-d5	93	30-130
2-Fluorobiphenyl	90	30-130
2,4,6-Tribromophenol	95	15-115
Terphenyl-d14	106	30-130



**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

Lab Number: L2249775

**Report Date:** 10/11/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Semivolatile Organics by GC/MS - Mansfield	Lab Associated	d sample(s):	01-06 Batch:	WG1688692-2	WG1688692-3			
bis(2-Chloroethyl)ether	84		88		40-140	5		20
1,3-Dichlorobenzene	44		53		40-140	19		20
1,4-Dichlorobenzene	46		54		40-140	16		20
1,2-Dichlorobenzene	48		56		40-140	15		20
bis(2-chloroisopropyl)ether	75		80		40-140	6		20
Acetophenone	82		85		40-140	4		20
Hexachloroethane	40		49		10-97	20		20
Nitrobenzene	83		88		40-140	6		20
Isophorone	82		86		40-140	5		20
bis(2-Chloroethoxy)methane	89		93		40-140	4		20
1,2,4-Trichlorobenzene	50		57		40-140	13		20
Naphthalene	65		71		40-140	9		20
4-Chloroaniline	93		97		40-140	4		20
Hexachlorobutadiene	41		47		40-140	14		20
2-Methylnaphthalene	64		69		40-140	8		20
1,2,4,5-Tetrachlorobenzene	56		61		40-140	9		20
Hexachlorocyclopentadiene	40		48		10-109	18		20
Biphenyl	75		74		40-140	1		20
2-Chloronaphthalene	68		73		40-140	7		20
2-Nitroaniline	93		98		40-140	5		20
Acenaphthylene	77		81		40-140	5		20
Dimethylphthalate	84		80		40-140	5		20
2,6-Dinitrotoluene	101		106		40-140	5		20



**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15

Lab Number: L2249775

**Report Date:** 10/11/22

Parameter	LCS %Recovery	Qual		CSD covery	Qual	%Recovery Limits	RPD	Qual	RPD Limits	
Semivolatile Organics by GC/MS - Mansfield	Lab Associate	d sample(s):	01-06	Batch:	WG1688692-2	2 WG1688692-3				
Acenaphthene	76			80		40-140	5		20	
3-Nitroaniline	100			106		40-140	6		20	
Dibenzofuran	84			87		40-140	4		20	
2,4-Dinitrotoluene	106			110		40-140	4		20	
Fluorene	86			89		40-140	3		20	
Diethylphthalate	92			92		40-140	0		20	
4-Nitroaniline	108			113		40-140	5		20	
n-Nitrosodiphenylamine	99			103		40-140	4		20	
Hexachlorobenzene	95			98		40-140	3		20	
Phenanthrene	93			96		40-140	3		20	
Anthracene	96			100		40-140	4		20	
Carbazole	100			106		40-140	6		20	
Di-n-butylphthalate	101			104		40-140	3		20	
Fluoranthene	98			103		40-140	5		20	
Pyrene	94			99		40-140	5		20	
Butylbenzylphthalate	97			101		40-140	4		20	
3,3'-Dichlorobenzidine	92			99		40-140	7		20	
Benz(a)anthracene	91			97		40-140	6		20	
Chrysene	99			105		40-140	6		20	
bis(2-Ethylhexyl)phthalate	99			105		40-140	6		20	
Di-n-octylphthalate	103			108		40-140	5		20	
Benzo(b)fluoranthene	89			93		40-140	4		20	
Benzo(k)fluoranthene	105			111		40-140	6		20	



**Project Name:** STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 

03.0033579.15

Lab Number:

L2249775

Report Date:

10/11/22

<u>Parameter</u>	LCS %Recovery	Qual	_	CSD ecovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits	
Semivolatile Organics by GC/MS - Mansfield	Lab Associated	d sample(s):	01-06	Batch:	WG1688692-2	WG1688692-3				
Benzo(a)pyrene	97			101		40-140	4		20	
Indeno(1,2,3-cd)pyrene	102			109		40-140	7		20	
Dibenz(a,h)anthracene	101			105		40-140	4		20	
Benzo(g,h,i)perylene	91			95		40-140	4		20	

	LCS	LCSD	Acceptance
Surrogate	%Recovery Qu	al %Recovery Qua	al Criteria
2-Fluorophenol	58	61	15-115
Phenol-d5	42	43	15-115
Nitrobenzene-d5	86	89	30-130
2-Fluorobiphenyl	80	81	30-130
2,4,6-Tribromophenol	99	102	15-115
Terphenyl-d14	95	100	30-130

Serial\_No:10112209:57 **Lab Number:** L2249775

Project Name: STEELWINDS ANNUAL/SEMIANNUAL

**Project Number:** 03.0033579.15 **Report Date:** 10/11/22

### Sample Receipt and Container Information

YES Were project specific reporting limits specified?

**Cooler Information** 

**Custody Seal** Cooler

Α Absent

Container Information			Initial	Final	Temp			Frozen		
Container ID	Container Type	Cooler	рН	pН	deg C	Pres	Seal	Date/Time	Analysis(*)	
L2249775-01A	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-01B	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-01C	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-01D	Amber 1000ml unpreserved	Α	7	7	3.1	Υ	Absent		A2-SVOC-8270(7)	
L2249775-01E	Amber 1000ml unpreserved	Α	7	7	3.1	Υ	Absent		A2-SVOC-8270(7)	
L2249775-02A	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-02B	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-02C	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-02D	Amber 1000ml unpreserved	Α	7	7	3.1	Υ	Absent		A2-SVOC-8270(7)	
L2249775-02E	Amber 1000ml unpreserved	Α	7	7	3.1	Υ	Absent		A2-SVOC-8270(7)	
L2249775-03A	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-03B	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-03C	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-03D	Amber 1000ml unpreserved	Α	7	7	3.1	Υ	Absent		A2-SVOC-8270(7)	
L2249775-03E	Amber 1000ml unpreserved	Α	7	7	3.1	Υ	Absent		A2-SVOC-8270(7)	
L2249775-04A	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-04B	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-04C	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-04D	Amber 1000ml unpreserved	Α	7	7	3.1	Υ	Absent		A2-SVOC-8270(7)	
L2249775-04E	Amber 1000ml unpreserved	Α	7	7	3.1	Υ	Absent		A2-SVOC-8270(7)	
L2249775-05A	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-05B	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	
L2249775-05C	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)	



**Lab Number:** L2249775

Report Date: 10/11/22

**Project Number:** 03.0033579.15

STEELWINDS ANNUAL/SEMIANNUAL

Project Name:

Container Information			Initial	Final	Temp			Frozen			
Container ID	Container Type	Cooler	рН	pН	deg C	Pres	Seal	Date/Time	Analysis(*)		
L2249775-05D	Amber 1000ml unpreserved	Α	7	7	3.1	Υ	Absent		A2-SVOC-8270(7)		
L2249775-05E	Amber 1000ml unpreserved	Α	7	7	3.1	Υ	Absent		A2-SVOC-8270(7)		
L2249775-06A	Vial HCI preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)		
L2249775-06B	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)		
L2249775-06C	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)		
L2249775-06D	Amber 1000ml unpreserved	Α	7	7	3.1	Υ	Absent		A2-SVOC-8270(7)		
L2249775-06E	Amber 1000ml unpreserved	Α	7	7	3.1	Υ	Absent		A2-SVOC-8270(7)		
L2249775-07A	Vial HCl preserved	Α	NA		3.1	Υ	Absent		NYCP51-8260(14)		

**Project Name:** Lab Number: STEELWINDS ANNUAL/SEMIANNUAL L2249775 **Report Date: Project Number:** 03.0033579.15 10/11/22

### **GLOSSARY**

#### **Acronyms**

**EDL** 

LOQ

MS

DL - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

- Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis of PAHs using Solid-Phase Microextraction (SPME).

**EMPC** - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case

estimate of the concentration. **EPA** 

LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of

analytes or a material containing known and verified amounts of analytes.

LCSD Laboratory Control Sample Duplicate: Refer to LCS.

Environmental Protection Agency.

LFB - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LOD - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content,

where applicable. (DoD report formats only.) - Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The

LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats

Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats

MDI - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

> - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated

using the native concentration, including estimated values.

MSD - Matrix Spike Sample Duplicate: Refer to MS.

NA - Not Applicable.

NC - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's

reporting unit.

NDPA/DPA - N-Nitrosodiphenylamine/Diphenylamine.

NI - Not Ignitable.

NP - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.

- No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile NR

Organic TIC only requests.

RL - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL

includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the

values; although the RPD value will be provided in the report.

SRM - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the

associated field samples.

STLP - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.

TEF - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.

TEO - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF

and then summing the resulting values.

TIC - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.

Report Format: DU Report with 'J' Qualifiers



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#### **Footnotes**

1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

#### **Terms**

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Chlordane: The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA,this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

Difference: With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

Final pH: As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

Frozen Date/Time: With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

Gasoline Range Organics (GRO): Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

Initial pH: As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

PAH Total: With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benza(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

PFAS Total: With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

Total: With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

# Data Qualifiers

- A -Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentrations of the analyte at less than ten times (10x) the concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G The concentration may be biased high due to matrix interferences (i.e, co-elution) with non-target compound(s). The result should be considered estimated.
- H The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I The lower value for the two columns has been reported due to obvious interference.
- J Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively

Report Format: DU Report with 'J' Qualifiers



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#### **Data Qualifiers**

Identified Compounds (TICs).

- $\label{eq:main_main_model} \textbf{M} \qquad \text{-Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.}$
- ND Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.
- **NJ** Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- ${f P}$  The RPD between the results for the two columns exceeds the method-specified criteria.
- Q The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- RE Analytical results are from sample re-extraction.
- S Analytical results are from modified screening analysis.
- V The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits.
   (Applicable to MassDEP DW Compliance samples only.)

Report Format: DU Report with 'J' Qualifiers



Project Name:STEELWINDS ANNUAL/SEMIANNUALLab Number:L2249775Project Number:03.0033579.15Report Date:10/11/22

### REFERENCES

Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

# **LIMITATION OF LIABILITIES**

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Alpha Analytical, Inc. Facility: Company-wide

Department: Quality Assurance

Title: Certificate/Approval Program Summary

ID No.:17873 Revision 19

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# Certification Information

### The following analytes are not included in our Primary NELAP Scope of Accreditation:

Westborough Facility

EPA 624/624.1: m/p-xylene, o-xylene, Naphthalene

EPA 625/625.1: alpha-Terpineol

EPA 8260C/8260D: NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene;

EPA 8270D/8270E: NPW: Dimethylnaphthalene,1,4-Diphenylhydrazine, alpha-Terpineol; SCM: Dimethylnaphthalene,1,4-Diphenylhydrazine.

SM4500: NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO2, NO3.

**Mansfield Facility** 

**SM 2540D:** TSS

EPA 8082A: NPW: PCB: 1, 5, 31, 87,101, 110, 141, 151, 153, 180, 183, 187.

EPA TO-15: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

Biological Tissue Matrix: EPA 3050B

### The following analytes are included in our Massachusetts DEP Scope of Accreditation

### Westborough Facility:

### **Drinking Water**

EPA 300.0: Chloride, Nitrate-N, Fluoride, Sulfate; EPA 353.2: Nitrate-N, Nitrite-N; SM4500NO3-F: Nitrate-N, Nitrite-N; SM4500F-C, SM4500CN-CE,

EPA 180.1, SM2130B, SM4500CI-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B

EPA 332: Perchlorate; EPA 524.2: THMs and VOCs; EPA 504.1: EDB, DBCP.

Microbiology: SM9215B; SM9223-P/A, SM9223B-Colilert-QT,SM9222D.

#### Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, EPA 350.1: Ammonia-N, LACHAT 10-107-06-1-B: Ammonia-N, EPA 351.1, SM4500NO3-F, EPA 353.2: Nitrate-N, SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300: Chloride, Sulfate, Nitrate. EPA 624.1: Volatile Halocarbons & Aromatics,

EPA 608.3: Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan II, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

EPA 625.1: SVOC (Acid/Base/Neutral Extractables), EPA 600/4-81-045: PCB-Oil.

Microbiology: SM9223B-Colilert-QT; Enterolert-QT, SM9221E, EPA 1600, EPA 1603, SM9222D.

## Mansfield Facility:

### **Drinking Water**

EPA 200.7: Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. EPA 200.8: Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. EPA 245.1 Hg. EPA 522, EPA 537.1.

#### Non-Potable Water

EPA 200.7: Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.

EPA 200.8: Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.

EPA 245.1 Hg

SM2340B

For a complete listing of analytes and methods, please contact your Alpha Project Manager.

Pre-Qualtrax Document ID: 08-113 Document Type: Form

Westborough, MA 01581 8 Walkup Dr. TEL: 508-898-9220 FAX: 508-898-9193	NEW YORK CHAIN OF CUSTODY Mansfield, MA 02048 320 Forbes Blvd TEL: 508-822-9300 FAX: 508-822-3288	Service Centers Mahwah, NJ 07430: 35 Whitney Albany, NY 12205: 14 Walker W Tonawanda, NY 14150: 275 Coo  Project Information  Project Name: STEE  Project Location: LAC	L WINDS	A NY	Page / of	1	Delive	Date Rec'd in Lab erables ASP-A EQuIS (1 F	q/		ASP-B EQuIS		ALPHA Job # 75  Billing Information  Same as Client Info
Client Information	100 10 27 10 10	Project # 03.003	3579.15				_			AT B	*	7-11-	
Client: GZA		(Use Project name as Project name)	oject#)				DOM:	latory Requ	iremer				Disposal Site Information
Address: 300 Pea	rl St. Suite 700	Project Manager: DA	NIEL +	ROY				NY TOGS			NY Part		Please identify below location of
BUTTALO, NY	14202	ALPHAQuote #:						AWQ Standa	ards	-	VY CP-	51	applicable disposal facilities.
Phone: (716) 517	-5708	Turn-Around Time		ST BOOK	200			NY Restricte	d Use		Other		Disposal Facility:
Fax:		Standard		Due Date:				NY Unrestric					NJ NY
Email: DANIE LITE	204 @ GZA. COM	Rush (only if pre approved	)	# of Days:				NYC Sewer	Dischar	ge			Other:
These samples have be	een previously analyze	ed by Alpha					ANA	LYSIS					Sample Filtration
Other project specific  Please specify Metals		ents:					STARS	РАН					Done Lab to do  Preservation Lab to do  (Please Specify below)
ALPHA Lab ID		0000000	Coll	ection	Sample	Sampler's	8260	270		ΙI			t
(Lab Use Only)	Sa	mple ID	Date	Time	Matrix	Initials	82	8		1 1			Sample Specific Comments e
99775.01	WT1-05-	091322	9-13-22	0755	GW	PSN	×	X					
.02	MWN-01-0	A CALLES COMMENT OF THE STATE O	i	0905	1	1	×	×					
603	MWN-01B-			1015			×	X					
Pa	WT1-04-			1055			×	×					
-07	BEP-ORC-1			1325			X	V		$\Box$			
,06	WT4-02-0			1415	1		×	X					
.07	TRIP BLAN		1	×	w	9	X						
THE RESERVE OF THE PARTY OF THE							$\top$						
									1				
B = HCI C = HNO <sub>3</sub> D = H <sub>2</sub> SO <sub>4</sub>	Container Code P = Plastic A = Amber Glass V = Vial G = Glass B = Bacteria Cup	Westboro: Certification N Mansfield: Certification N				tainer Type	-						Please print clearly, legibly and completely. Samples can not be logged in and turnaround time clock will not start until any ambiguities are
F = MeOH	C = Cube	Relinquished	By:	Date/	Time		Recei	ved By:			Date/	Time	resolved. BY EXECUTING
0 - 14011001	O = Other E = Encore	Pin	2	9-13-22		Elle		SAL		9-13	22	1441	THIS COC, THE CLIENT HAS READ AND AGREES
H = 14020203	D = BOD Bottle	TELES AA	۷	9-13-22	1445	1	_		٠	9/14	1/22	030	TO BE BOUND BY ALPHA'S TERMS & CONDITIONS.
Form No: 01-25 HC (rev. 30	0-Sept-2013)												(See reverse side.)



## ANALYTICAL REPORT

Lab Number: L2250076

Client: GZA GeoEnvironmental of New York

300 Pearl Street

Suite 700

Buffalo, NY 14202

ATTN: Dan Troy

Phone: (716) 844-7050

Project Name: STEELWINDS ANNUAL/ SEMIANNUAL

Project Number: 03.0033579.15

Report Date: 10/12/22

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Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0574), IL (200077), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #P330-17-00196).

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



**Project Name:** STEELWINDS ANNUAL/ SEMIANNUAL

**Project Number:** 03.0033579.15

 Lab Number:
 L2250076

 Report Date:
 10/12/22

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2250076-01	MWN-02-091422	WATER	LACKAWANNA NY	09/14/22 08:00	09/14/22
L2250076-02	MWN-02B-091422	WATER	LACKAWANNA NY	09/14/22 08:45	09/14/22
L2250076-03	TRIP BLANK-2	WATER	LACKAWANNA NY	09/14/22 00:00	09/14/22



Project Name: STEELWINDS ANNUAL/ SEMIANNUAL Lab Number: L2250076

**Project Number:** 03.0033579.15 **Report Date:** 10/12/22

### **Case Narrative**

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments, solids and tissues are reported on a dry weight basis unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

HOLD POLICY - For samples submitted on hold, Alpha's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Alpha Project Manager and made arrangements for Alpha to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.	



Project Name:STEELWINDS ANNUAL/ SEMIANNUALLab Number:L2250076Project Number:03.0033579.15Report Date:10/12/22

# **Case Narrative (continued)**

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

Sample Receipt

L2250076-03: A sample identified as "TRIP BLANK-2" was listed on the Chain of Custody, but not received. This was verified by the client.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Jufani Morrissey-Tiffani Morrissey

Authorized Signature:

Title: Technical Director/Representative Date: 10/12/22

# **ORGANICS**



# **VOLATILES**



Project Name: STEELWINDS ANNUAL/ SEMIANNUAL

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Lab Number: L2250076

**Report Date:** 10/12/22

Lab ID: L2250076-01 Date Collected: 09/14/22 08:00

Client ID: MWN-02-091422 Date Received: 09/14/22 Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Matrix: Water
Analytical Method: 1,8260C
Analytical Date: 09/24/22 22:08

Analyst: PD

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westbor	ough Lab					
Benzene	1.5		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	ND		ug/l	2.5	0.70	1
o-Xylene	ND		ug/l	2.5	0.70	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	4.2		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	ND		ug/l	2.5	0.70	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	102	70-130	
Toluene-d8	100	70-130	
4-Bromofluorobenzene	93	70-130	
Dibromofluoromethane	108	70-130	



L2250076

09/14/22 08:45

**Project Name:** STEELWINDS ANNUAL/ SEMIANNUAL

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Lab Number:

Date Collected:

Report Date: 10/12/22

Lab ID: D L2250076-02

Client ID: MWN-02B-091422 Sample Location: LACKAWANNA NY Date Received: 09/14/22 Field Prep: Not Specified

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/24/22 22:31

Analyst: PD

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor					
Volatile Organics by GC/MS - Westborough Lab											
Benzene	62		ug/l	1.2	0.40	2.5					
Toluene	10		ug/l	6.2	1.8	2.5					
Ethylbenzene	ND		ug/l	6.2	1.8	2.5					
Methyl tert butyl ether	ND		ug/l	6.2	1.8	2.5					
p/m-Xylene	7.2		ug/l	6.2	1.8	2.5					
o-Xylene	10		ug/l	6.2	1.8	2.5					
n-Butylbenzene	ND		ug/l	6.2	1.8	2.5					
sec-Butylbenzene	ND		ug/l	6.2	1.8	2.5					
tert-Butylbenzene	ND		ug/l	6.2	1.8	2.5					
Isopropylbenzene	ND		ug/l	6.2	1.8	2.5					
p-Isopropyltoluene	ND		ug/l	6.2	1.8	2.5					
Naphthalene	320		ug/l	6.2	1.8	2.5					
n-Propylbenzene	ND		ug/l	6.2	1.8	2.5					
1,3,5-Trimethylbenzene	ND		ug/l	6.2	1.8	2.5					
1,2,4-Trimethylbenzene	1.9	J	ug/l	6.2	1.8	2.5					

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	101	70-130	
Toluene-d8	102	70-130	
4-Bromofluorobenzene	91	70-130	
Dibromofluoromethane	106	70-130	



Project Name: STEELWINDS ANNUAL/ SEMIANNUAL Lab Number: L2250076

**Project Number:** 03.0033579.15 **Report Date:** 10/12/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/24/22 17:52

Analyst: LAC

Parameter	Result	Qualifier Units	s RL	MDL	
Volatile Organics by GC/MS - We	stborough Lab	for sample(s):	01-02 Batch:	WG1692585-5	
Benzene	ND	ug/l	0.50	0.16	
Toluene	ND	ug/l	2.5	0.70	
Ethylbenzene	ND	ug/l	2.5	0.70	
Methyl tert butyl ether	ND	ug/l	2.5	0.70	
p/m-Xylene	ND	ug/l	2.5	0.70	
o-Xylene	ND	ug/l	2.5	0.70	
n-Butylbenzene	ND	ug/l	2.5	0.70	
sec-Butylbenzene	ND	ug/l	2.5	0.70	
tert-Butylbenzene	ND	ug/l	2.5	0.70	
Isopropylbenzene	ND	ug/l	2.5	0.70	
p-Isopropyltoluene	ND	ug/l	2.5	0.70	
Naphthalene	ND	ug/l	2.5	0.70	
n-Propylbenzene	ND	ug/l	2.5	0.70	
1,3,5-Trimethylbenzene	ND	ug/l	2.5	0.70	
1,2,4-Trimethylbenzene	ND	ug/l	2.5	0.70	

		Acceptance	
Surrogate	%Recovery Qualif	er Criteria	
1,2-Dichloroethane-d4	102	70-130	
Toluene-d8	100	70-130	
4-Bromofluorobenzene	93	70-130	
Dibromofluoromethane	107	70-130	



**Project Name:** STEELWINDS ANNUAL/ SEMIANNUAL

**Project Number:** 03.0033579.15

Lab Number: L2250076

**Report Date:** 10/12/22

arameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
olatile Organics by GC/MS - Westborough L	ab Associated	sample(s):	01-02 Batch: W	G1692585-3	WG1692585-4			
Benzene	100		100		70-130	0		20
Toluene	100		100		70-130	0		20
Ethylbenzene	98		100		70-130	2		20
Methyl tert butyl ether	94		100		63-130	6		20
p/m-Xylene	100		100		70-130	0		20
o-Xylene	95		100		70-130	5		20
n-Butylbenzene	99		100		53-136	1		20
sec-Butylbenzene	99		100		70-130	1		20
tert-Butylbenzene	97		97		70-130	0		20
Isopropylbenzene	96		99		70-130	3		20
p-Isopropyltoluene	98		99		70-130	1		20
Naphthalene	89		100		70-130	12		20
n-Propylbenzene	98		100		69-130	2		20
1,3,5-Trimethylbenzene	94		96		64-130	2		20
1,2,4-Trimethylbenzene	95		98		70-130	3		20

Surrogate	LCS %Recovery Qual	LCSD %Recovery Qual	Acceptance Criteria
1,2-Dichloroethane-d4	104	107	70-130
Toluene-d8	103	103	70-130
4-Bromofluorobenzene	94	94	70-130
Dibromofluoromethane	107	106	70-130



# **SEMIVOLATILES**



L2250076

10/12/22

09/17/22 08:30

Project Name: STEELWINDS ANNUAL/ SEMIANNUAL Lab Number:

**Project Number:** 03.0033579.15

L2250076-01

MWN-02-091422

LACKAWANNA NY

**SAMPLE RESULTS** 

Date Collected: 09/14/22 08:00

Date Received: 09/14/22

Extraction Method: EPA 3510C

Field Prep: Not Specified

**Extraction Date:** 

Report Date:

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Matrix: Water
Analytical Method: 1,8270D

Analytical Date: 09/19/22 15:08

Analyst: PS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS - M	ansfield Lab					
bis(2-Chloroethyl)ether	ND		ug/l	0.495	0.092	1
1,3-Dichlorobenzene	ND		ug/l	0.495	0.078	1
1,4-Dichlorobenzene	ND		ug/l	0.495	0.082	1
1,2-Dichlorobenzene	ND		ug/l	0.495	0.067	1
Benzyl alcohol	ND		ug/l	0.495	0.122	1
bis(2-chloroisopropyl)ether	ND		ug/l	0.495	0.107	1
Acetophenone	0.246	J	ug/l	0.990	0.205	1
Hexachloroethane	ND		ug/l	0.495	0.101	1
Nitrobenzene	ND		ug/l	0.495	0.101	1
Isophorone	ND		ug/l	0.495	0.125	1
bis(2-Chloroethoxy)methane	ND		ug/l	0.495	0.085	1
1,2,4-Trichlorobenzene	ND		ug/l	0.495	0.095	1
Naphthalene	3.44		ug/l	0.495	0.087	1
4-Chloroaniline	ND		ug/l	0.495	0.127	1
Hexachlorobutadiene	ND		ug/l	0.495	0.085	1
2-Methylnaphthalene	1.01		ug/l	0.495	0.090	1
1,2,4,5-Tetrachlorobenzene	ND		ug/l	0.495	0.079	1
Hexachlorocyclopentadiene	ND		ug/l	0.495	0.151	1
Biphenyl	0.332	J	ug/l	0.495	0.110	1
2-Chloronaphthalene	ND		ug/l	0.495	0.089	1
2-Nitroaniline	ND		ug/l	0.495	0.137	1
Acenaphthylene	1.03		ug/l	0.495	0.111	1
Dimethylphthalate	ND		ug/l	0.495	0.116	1
2,6-Dinitrotoluene	ND		ug/l	0.495	0.166	1
Acenaphthene	0.603		ug/l	0.495	0.095	1
3-Nitroaniline	ND		ug/l	0.495	0.110	1
Dibenzofuran	0.967		ug/l	0.495	0.090	1
2,4-Dinitrotoluene	ND		ug/l	0.495	0.161	1



10/12/22

**Project Name:** Lab Number: STEELWINDS ANNUAL/ SEMIANNUAL L2250076

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Date Collected: 09/14/22 08:00

Report Date:

L2250076-01 Date Received: 09/14/22 MWN-02-091422 Field Prep: LACKAWANNA NY Not Specified

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Diethylphthalate   ND	Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Diethylphthalate   ND	Semivolatile Organics by GC/M	S - Mansfield Lab					
Diethylphthalate   ND   ug/l   0.495   0.178   1   1   1   1   1   1   1   1   1	Fluorene	2.26		ua/l	0.495	0.103	1
A-Nitroaniline   ND   ug/l   0.495   0.111   1   1   1   1   1   1   1   1   1	Diethylphthalate	ND			0.495	0.178	1
ND	4-Nitroaniline	ND			0.495	0.111	1
Phenanthrene 1.76 ug/l 0.495 0.110 1 Anthracene 0.588 ug/l 0.495 0.136 1 Carbazole 1.28 ug/l 0.495 0.142 1 Di-n-butylphthalate ND ug/l 0.495 0.099 1 Fluoranthene 0.971 ug/l 0.495 0.154 1 Pyrene 1.70 ug/l 0.495 0.168 1 Betylbenzylphthalate ND ug/l 0.495 0.084 1 Benzo(a)anthracene ND ug/l 0.495 0.191 1 Benzo(b)fluoranthene ND ug/l 0.495 0.182 1 Chrysene ND ug/l 0.495 0.182 1 Chrysene ND ug/l 0.495 0.182 1 Chrysene ND ug/l 0.495 0.180 1 Di-n-octylphthalate ND ug/l 0.495 0.080 1 Di-n-octylphthalate ND ug/l 0.495 0.080 1 Di-n-octylphthalate ND ug/l 0.495 0.065 1 Benzo(b)fluoranthene ND ug/l 0.495 0.065 1 Benzo(b)fluoranthene ND ug/l 0.495 0.060 1 Benzo(b)fluoranthene ND ug/l 0.495 0.060 1 Benzo(a)pyrene ND ug/l 0.495 0.060 1 Benzo(a)pyrene ND ug/l 0.495 0.060 1 Indeno(1,2,3-cd)pyrene ND ug/l 0.495 0.089 1 Dibenz(a,h)anthracene ND ug/l 0.495 0.089 1	n-Nitrosodiphenylamine	ND		ug/l	0.495	0.071	1
Anthracene	Hexachlorobenzene	ND		ug/l	0.495	0.121	1
1.28	Phenanthrene	1.76		ug/l	0.495	0.110	1
Di-n-butylphthalate	Anthracene	0.588		ug/l	0.495	0.136	1
Pyrene   1.70   ug/l   0.495   0.154   1	Carbazole	1.28		ug/l	0.495	0.142	1
Pyrene         1.70         ug/l         0.495         0.168         1           Butylbenzylphthalate         ND         ug/l         0.495         0.084         1           3,3'-Dichlorobenzidine         ND         ug/l         0.495         0.191         1           Benz(a)anthracene         ND         ug/l         0.495         0.182         1           Chrysene         ND         ug/l         0.495         0.140         1           Dis(2-Ethylhexyl)phthalate         ND         ug/l         0.495         0.080         1           Di-n-octylphthalate         ND         ug/l         0.990         0.078         1           Benzo(b)fluoranthene         ND         ug/l         0.495         0.065         1           Benzo(k)fluoranthene         ND         ug/l         0.495         0.159         1           Benzo(a)pyrene         ND         ug/l         0.495         0.060         1           Indeno(1,2,3-cd)pyrene         ND         ug/l         0.495         0.089         1           Dibenz(a,h)anthracene         ND         ug/l         0.495         0.064         1	Di-n-butylphthalate	ND		ug/l	0.495	0.099	1
Butylbenzylphthalate ND ug/l 0.495 0.084 1 3,3'-Dichlorobenzidine ND ug/l 0.495 0.191 1 Benz(a)anthracene ND ug/l 0.495 0.182 1 Chrysene ND ug/l 0.495 0.182 1 Chrysene ND ug/l 0.495 0.080 1 Di-n-octylphthalate ND ug/l 0.495 0.080 1 Di-n-octylphthalate ND ug/l 0.495 0.080 1 Di-n-octylphthalate ND ug/l 0.495 0.065 1 Benzo(b)fluoranthene ND ug/l 0.495 0.065 1 Benzo(k)fluoranthene ND ug/l 0.495 0.159 1 Benzo(a)pyrene ND ug/l 0.495 0.060 1 Indeno(1,2,3-cd)pyrene ND ug/l 0.495 0.089 1 Dibenz(a,h)anthracene ND ug/l 0.495 0.089 1	Fluoranthene	0.971		ug/l	0.495	0.154	1
ND	Pyrene	1.70		ug/l	0.495	0.168	1
ND	Butylbenzylphthalate	ND		ug/l	0.495	0.084	1
Chrysene         ND         ug/l         0.495         0.140         1           bis(2-Ethylhexyl)phthalate         ND         ug/l         0.495         0.080         1           Di-n-octylphthalate         ND         ug/l         0.990         0.078         1           Benzo(b)fluoranthene         ND         ug/l         0.495         0.065         1           Benzo(k)fluoranthene         ND         ug/l         0.495         0.159         1           Benzo(a)pyrene         ND         ug/l         0.495         0.060         1           Indeno(1,2,3-cd)pyrene         ND         ug/l         0.495         0.089         1           Dibenz(a,h)anthracene         ND         ug/l         0.495         0.064         1	3,3'-Dichlorobenzidine	ND		ug/l	0.495	0.191	1
ND   ug/l   0.495   0.080   1     Di-n-octylphthalate   ND   ug/l   0.990   0.078   1     Di-n-octylphthalate   ND   ug/l   0.495   0.065   1     Di-n-octylphthalate   ND   ug/l   0.495   0.065   1     Di-n-octylphthalate   ND   ug/l   0.495   0.159   1     Di-n-octylphthalate   ND   ug/l   0.495   0.159   1     Di-n-octylphthalate   ND   ug/l   0.495   0.060   1     Di-n-octylphthalate   ND   ug/l   0.495   0.060   1     Di-n-octylphthalate   ND   ug/l   0.495   0.089   1   Di-n-octylphthalate   ND   ug/l   0.495   0.064   1   Di-n-octylphthalate   ND   ug/l   0.495   0.065   1   Di-n-octylphthalate   Di-n-octylphthalate	Benz(a)anthracene	ND		ug/l	0.495	0.182	1
Di-n-octylphthalate	Chrysene	ND		ug/l	0.495	0.140	1
Benzo(b)fluoranthene         ND         ug/l         0.495         0.065         1           Benzo(k)fluoranthene         ND         ug/l         0.495         0.159         1           Benzo(a)pyrene         ND         ug/l         0.495         0.060         1           Indeno(1,2,3-cd)pyrene         ND         ug/l         0.495         0.089         1           Dibenz(a,h)anthracene         ND         ug/l         0.495         0.064         1	bis(2-Ethylhexyl)phthalate	ND		ug/l	0.495	0.080	1
ND   ug/l   0.495   0.159   1	Di-n-octylphthalate	ND		ug/l	0.990	0.078	1
ND   ug/l   0.495   0.060   1	Benzo(b)fluoranthene	ND		ug/l	0.495	0.065	1
Indeno(1,2,3-cd)pyrene         ND         ug/l         0.495         0.089         1           Dibenz(a,h)anthracene         ND         ug/l         0.495         0.064         1	Benzo(k)fluoranthene	ND		ug/l	0.495	0.159	1
Dibenz(a,h)anthracene ND ug/l 0.495 0.064 1	Benzo(a)pyrene	ND		ug/l	0.495	0.060	1
<b>49.</b>	Indeno(1,2,3-cd)pyrene	ND		ug/l	0.495	0.089	1
Benzo(g,h,i)perylene ND ug/l 0.495 0.108 1	Dibenz(a,h)anthracene	ND		ug/l	0.495	0.064	1
	Benzo(g,h,i)perylene	ND		ug/l	0.495	0.108	1

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	43	15-115
Phenol-d5	27	15-115
Nitrobenzene-d5	86	30-130
2-Fluorobiphenyl	78	30-130
2,4,6-Tribromophenol	107	15-115
Terphenyl-d14	96	30-130



L2250076

10/12/22

09/17/22 08:30

**Project Name:** STEELWINDS ANNUAL/ SEMIANNUAL

L2250076-02

MWN-02B-091422

LACKAWANNA NY

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Date Collected: 09/14/22 08:45

Date Received: 09/14/22

Extraction Method: EPA 3510C

Lab Number:

Report Date:

**Extraction Date:** 

Field Prep: Not Specified

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Matrix: Water Analytical Method: 1,8270D

Analytical Date: 09/19/22 15:38

Analyst: PS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Semivolatile Organics by GC/MS - M	lansfield Lab						
bis(2-Chloroethyl)ether	ND		ug/l	0.490	0.091	1	
1,3-Dichlorobenzene	ND		ug/l	0.490	0.077	1	
1,4-Dichlorobenzene	ND		ug/l	0.490	0.081	1	
1,2-Dichlorobenzene	0.200	J	ug/l	0.490	0.067	1	
Benzyl alcohol	ND		ug/l	0.490	0.120	1	
bis(2-chloroisopropyl)ether	ND		ug/l	0.490	0.106	1	
Acetophenone	0.770	J	ug/l	0.980	0.203	1	
Hexachloroethane	ND		ug/l	0.490	0.100	1	
Nitrobenzene	ND		ug/l	0.490	0.100	1	
Isophorone	ND		ug/l	0.490	0.124	1	
bis(2-Chloroethoxy)methane	ND		ug/l	0.490	0.084	1	
1,2,4-Trichlorobenzene	ND		ug/l	0.490	0.094	1	
Naphthalene	83.8	Е	ug/l	0.490	0.086	1	
4-Chloroaniline	ND		ug/l	0.490	0.125	1	
Hexachlorobutadiene	ND		ug/l	0.490	0.084	1	
2-Methylnaphthalene	8.48		ug/l	0.490	0.089	1	
1,2,4,5-Tetrachlorobenzene	ND		ug/l	0.490	0.078	1	
Hexachlorocyclopentadiene	ND		ug/l	0.490	0.150	1	
Biphenyl	1.11		ug/l	0.490	0.109	1	
2-Chloronaphthalene	ND		ug/l	0.490	0.088	1	
2-Nitroaniline	ND		ug/l	0.490	0.135	1	
Acenaphthylene	2.83		ug/l	0.490	0.110	1	
Dimethylphthalate	ND		ug/l	0.490	0.115	1	
2,6-Dinitrotoluene	ND		ug/l	0.490	0.165	1	
Acenaphthene	6.20		ug/l	0.490	0.094	1	
3-Nitroaniline	ND		ug/l	0.490	0.109	1	
Dibenzofuran	4.50		ug/l	0.490	0.089	1	
2,4-Dinitrotoluene	ND		ug/l	0.490	0.160	1	



10/12/22

Project Name: STEELWINDS ANNUAL/ SEMIANNUAL Lab Number: L2250076

**Project Number:** 03.0033579.15

L2250076-02

**SAMPLE RESULTS** 

Date Collected: 09/14/22 08:45

Report Date:

Client ID: MWN-02B-091422 Date Received: 09/14/22 Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Lab ID:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS	- Mansfield Lab					
Fluorene	7.72		ug/l	0.490	0.102	1
Diethylphthalate	ND		ug/l	0.490	0.176	1
4-Nitroaniline	ND		ug/l	0.490	0.110	1
n-Nitrosodiphenylamine	ND		ug/l	0.490	0.071	1
Hexachlorobenzene	ND		ug/l	0.490	0.120	1
Phenanthrene	13.7		ug/l	0.490	0.109	1
Anthracene	1.88		ug/l	0.490	0.134	1
Carbazole	21.2		ug/l	0.490	0.140	1
Di-n-butylphthalate	ND		ug/l	0.490	0.098	1
Fluoranthene	3.51		ug/l	0.490	0.153	1
Pyrene	2.00		ug/l	0.490	0.167	1
Butylbenzylphthalate	ND		ug/l	0.490	0.083	1
3,3'-Dichlorobenzidine	ND		ug/l	0.490	0.189	1
Benz(a)anthracene	ND		ug/l	0.490	0.180	1
Chrysene	ND		ug/l	0.490	0.139	1
bis(2-Ethylhexyl)phthalate	ND		ug/l	0.490	0.079	1
Di-n-octylphthalate	ND		ug/l	0.980	0.077	1
Benzo(b)fluoranthene	ND		ug/l	0.490	0.064	1
Benzo(k)fluoranthene	ND		ug/l	0.490	0.158	1
Benzo(a)pyrene	ND		ug/l	0.490	0.059	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.490	0.088	1
Dibenz(a,h)anthracene	ND		ug/l	0.490	0.063	1
Benzo(g,h,i)perylene	ND		ug/l	0.490	0.107	1

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	39	15-115
Phenol-d5	27	15-115
Nitrobenzene-d5	92	30-130
2-Fluorobiphenyl	69	30-130
2,4,6-Tribromophenol	104	15-115
Terphenyl-d14	92	30-130



10/12/22

Report Date:

**Project Name:** Lab Number: STEELWINDS ANNUAL/ SEMIANNUAL L2250076

**Project Number:** 03.0033579.15

09/20/22 14:24

**SAMPLE RESULTS** 

Lab ID: D Date Collected: 09/14/22 08:45 L2250076-02

Date Received: Client ID: 09/14/22 MWN-02B-091422 Sample Location: Field Prep: LACKAWANNA NY Not Specified

Sample Depth:

Analytical Date:

Extraction Method: EPA 3510C Matrix: Water **Extraction Date:** 09/17/22 08:30 Analytical Method: 1,8270D

Analyst: PS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS - Mansfield	Lab					
Naphthalene	146		ug/l	2.45	0.429	5

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	39	15-115
Phenol-d5	26	15-115
Nitrobenzene-d5	84	30-130
2-Fluorobiphenyl	80	30-130
2,4,6-Tribromophenol	99	15-115
Terphenyl-d14	96	30-130



L2250076

Project Name: STEELWINDS ANNUAL/ SEMIANNUAL Lab Number:

**Project Number:** 03.0033579.15 **Report Date:** 10/12/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270D Extraction Method: EPA 3510C
Analytical Date: 09/19/22 10:43 Extraction Date: 09/17/22 08:30

Analyst: PS

Parameter	Result	Qualifier	Units		RL	MDL	
Semivolatile Organics by GC/MS	- Mansfield La	ab for samp	le(s):	01-02	Batch:	WG1688692-1	
bis(2-Chloroethyl)ether	ND		ug/l		0.500	0.093	
1,3-Dichlorobenzene	ND		ug/l		0.500	0.078	
1,4-Dichlorobenzene	ND		ug/l		0.500	0.083	
1,2-Dichlorobenzene	ND		ug/l		0.500	0.068	
Benzyl alcohol	ND		ug/l		0.500	0.123	
bis(2-chloroisopropyl)ether	ND		ug/l		0.500	0.108	
Acetophenone	ND		ug/l		1.00	0.207	
Hexachloroethane	ND		ug/l		0.500	0.102	
Nitrobenzene	ND		ug/l		0.500	0.102	
Isophorone	ND		ug/l		0.500	0.126	
bis(2-Chloroethoxy)methane	ND		ug/l		0.500	0.085	
1,2,4-Trichlorobenzene	ND		ug/l		0.500	0.096	
Naphthalene	ND		ug/l		0.500	0.088	
4-Chloroaniline	ND		ug/l		0.500	0.128	
Hexachlorobutadiene	ND		ug/l		0.500	0.086	
2-Methylnaphthalene	ND		ug/l		0.500	0.091	
1,2,4,5-Tetrachlorobenzene	ND		ug/l		0.500	0.080	
Hexachlorocyclopentadiene	ND		ug/l		0.500	0.153	
Biphenyl	ND		ug/l		0.500	0.111	
2-Chloronaphthalene	ND		ug/l		0.500	0.090	
2-Nitroaniline	ND		ug/l		0.500	0.138	
Acenaphthylene	ND		ug/l		0.500	0.112	
Dimethylphthalate	ND		ug/l		0.500	0.117	
2,6-Dinitrotoluene	ND		ug/l		0.500	0.168	
Acenaphthene	ND		ug/l		0.500	0.096	
3-Nitroaniline	ND		ug/l		0.500	0.111	
Dibenzofuran	ND		ug/l		0.500	0.091	
2,4-Dinitrotoluene	ND		ug/l		0.500	0.163	
Fluorene	ND		ug/l		0.500	0.104	



L2250076

Project Name: STEELWINDS ANNUAL/ SEMIANNUAL Lab Number:

**Project Number:** 03.0033579.15 **Report Date:** 10/12/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270D Extraction Method: EPA 3510C
Analytical Date: 09/19/22 10:43 Extraction Date: 09/17/22 08:30

Analyst: PS

Parameter	Result	Qualifier	Units		RL	MDL	
Semivolatile Organics by GC/MS -	Mansfield La	ab for samp	ole(s):	01-02	Batch:	WG1688692-1	
Diethylphthalate	ND		ug/l		0.500	0.180	
4-Nitroaniline	ND		ug/l		0.500	0.112	
n-Nitrosodiphenylamine	ND		ug/l		0.500	0.072	
Hexachlorobenzene	ND		ug/l		0.500	0.122	
Phenanthrene	ND		ug/l		0.500	0.111	
Anthracene	ND		ug/l		0.500	0.137	
Carbazole	ND		ug/l		0.500	0.143	
Di-n-butylphthalate	ND		ug/l		0.500	0.100	
Fluoranthene	ND		ug/l		0.500	0.156	
Pyrene	ND		ug/l		0.500	0.170	
Butylbenzylphthalate	ND		ug/l		0.500	0.085	
3,3'-Dichlorobenzidine	ND		ug/l		0.500	0.193	
Benz(a)anthracene	ND		ug/l		0.500	0.184	
Chrysene	ND		ug/l		0.500	0.142	
bis(2-Ethylhexyl)phthalate	ND		ug/l		0.500	0.081	
Di-n-octylphthalate	ND		ug/l		1.00	0.079	
Benzo(b)fluoranthene	ND		ug/l		0.500	0.066	
Benzo(k)fluoranthene	ND		ug/l		0.500	0.161	
Benzo(a)pyrene	ND		ug/l		0.500	0.060	
Indeno(1,2,3-cd)pyrene	ND		ug/l		0.500	0.090	
Dibenz(a,h)anthracene	ND		ug/l		0.500	0.064	
Benzo(g,h,i)perylene	ND		ug/l		0.500	0.109	



Project Name: STEELWINDS ANNUAL/ SEMIANNUAL Lab Number: L2250076

**Project Number:** 03.0033579.15 **Report Date:** 10/12/22

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8270D Extraction Method: EPA 3510C
Analytical Date: 09/19/22 10:43 Extraction Date: 09/17/22 08:30

Analyst: PS

Parameter Result Qualifier Units RL MDL

Semivolatile Organics by GC/MS - Mansfield Lab for sample(s): 01-02 Batch: WG1688692-1

Surrogate	%Recovery Qualit	Acceptance fier Criteria
2-Fluorophenol	64	15-115
Phenol-d5	42	15-115
Nitrobenzene-d5	93	30-130
2-Fluorobiphenyl	90	30-130
2,4,6-Tribromophenol	95	15-115
Terphenyl-d14	106	30-130



**Project Name:** STEELWINDS ANNUAL/ SEMIANNUAL

**Project Number:** 03.0033579.15

Lab Number: L2250076

**Report Date:** 10/12/22

Parameter	LCS %Recovery	Qual		SD covery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Semivolatile Organics by GC/MS - Mansfield	Lab Associate	d sample(s):	01-02	Batch:	WG1688692-2	2 WG1688692-3			
bis(2-Chloroethyl)ether	84			88		40-140	5		20
1,3-Dichlorobenzene	44			53		40-140	19		20
1,4-Dichlorobenzene	46			54		40-140	16		20
1,2-Dichlorobenzene	48			56		40-140	15		20
bis(2-chloroisopropyl)ether	75			80		40-140	6		20
Acetophenone	82			85		40-140	4		20
Hexachloroethane	40			49		10-97	20		20
Nitrobenzene	83			88		40-140	6		20
Isophorone	82			86		40-140	5		20
bis(2-Chloroethoxy)methane	89			93		40-140	4		20
1,2,4-Trichlorobenzene	50			57		40-140	13		20
Naphthalene	65			71		40-140	9		20
4-Chloroaniline	93			97		40-140	4		20
Hexachlorobutadiene	41			47		40-140	14		20
2-Methylnaphthalene	64			69		40-140	8		20
1,2,4,5-Tetrachlorobenzene	56			61		40-140	9		20
Hexachlorocyclopentadiene	40			48		10-109	18		20
Biphenyl	75			74		40-140	1		20
2-Chloronaphthalene	68			73		40-140	7		20
2-Nitroaniline	93			98		40-140	5		20
Acenaphthylene	77			81		40-140	5		20
Dimethylphthalate	84			80		40-140	5		20
2,6-Dinitrotoluene	101		1	106		40-140	5		20



**Project Name:** STEELWINDS ANNUAL/ SEMIANNUAL

**Project Number:** 03.0033579.15

Lab Number: L2250076

**Report Date:** 10/12/22

Parameter	LCS %Recovery	Qual		SD covery	Qual	%Recovery Limits	RPD	Qual	RPD Limits	
Semivolatile Organics by GC/MS - Mansfield	Lab Associated	d sample(s):	01-02	Batch:	WG1688692-2	2 WG1688692-3				
Acenaphthene	76			80		40-140	5		20	
3-Nitroaniline	100		•	106		40-140	6		20	
Dibenzofuran	84			87		40-140	4		20	
2,4-Dinitrotoluene	106		•	110		40-140	4		20	
Fluorene	86			89		40-140	3		20	
Diethylphthalate	92			92		40-140	0		20	
4-Nitroaniline	108		•	113		40-140	5		20	
n-Nitrosodiphenylamine	99		•	103		40-140	4		20	
Hexachlorobenzene	95			98		40-140	3		20	
Phenanthrene	93			96		40-140	3		20	
Anthracene	96		•	100		40-140	4		20	
Carbazole	100		•	106		40-140	6		20	
Di-n-butylphthalate	101		•	104		40-140	3		20	
Fluoranthene	98		•	103		40-140	5		20	
Pyrene	94			99		40-140	5		20	
Butylbenzylphthalate	97		•	101		40-140	4		20	
3,3'-Dichlorobenzidine	92			99		40-140	7		20	
Benz(a)anthracene	91			97		40-140	6		20	
Chrysene	99		,	105		40-140	6		20	
bis(2-Ethylhexyl)phthalate	99		•	105		40-140	6		20	
Di-n-octylphthalate	103		,	108		40-140	5		20	
Benzo(b)fluoranthene	89			93		40-140	4		20	
Benzo(k)fluoranthene	105			111		40-140	6		20	

**Project Name:** STEELWINDS ANNUAL/ SEMIANNUAL

Lab Number: L2250076

**Project Number:** 03.0033579.15 Report Date: 10/12/22

<u>Parameter</u>	LCS %Recovery	Qual		SD covery	Qual	%Recovery Limits	RPD	Qual	RPD Limits	
Semivolatile Organics by GC/MS - Mansfiel	d Lab Associated	sample(s):	01-02	Batch:	WG1688692-2	2 WG1688692-3				
Benzo(a)pyrene	97		1	101		40-140	4		20	
Indeno(1,2,3-cd)pyrene	102		1	109		40-140	7		20	
Dibenz(a,h)anthracene	101		1	105		40-140	4		20	
Benzo(g,h,i)perylene	91			95		40-140	4		20	

Surveyede	LCS	LCSD	Acceptance Griteria
Surrogate	%Recovery Qu	al %Recovery Qua	ar Ciliena
2-Fluorophenol	58	61	15-115
Phenol-d5	42	43	15-115
Nitrobenzene-d5	86	89	30-130
2-Fluorobiphenyl	80	81	30-130
2,4,6-Tribromophenol	99	102	15-115
Terphenyl-d14	95	100	30-130



# **METALS**



Project Name: STEELWINDS ANNUAL/ SEMIANNUAL Lab Number: L2250076

**Project Number:** 03.0033579.15 **Report Date:** 10/12/22

**SAMPLE RESULTS** 

 Lab ID:
 L2250076-02
 Date Collected:
 09/14/22 08:45

 Client ID:
 MWN-02B-091422
 Date Received:
 09/14/22

Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Matrix: Water

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	Date Prepared	Date Analyzed	Prep Method	Analytical Method	Analyst
Total Metals - Ma	ansfield Lab										
Arsenic, Total	0.0379		mg/l	0.0050	0.0019	1	09/17/22 07:0	2 10/10/22 07:54	4 EPA 3005A	1,6010D	EW



L2250076

**Project Name:** STEELWINDS ANNUAL/ SEMIANNUAL

Lab Number:

**Project Number: Report Date:** 03.0033579.15 10/12/22

> **Method Blank Analysis Batch Quality Control**

**Dilution Date Date** Analytical Method Analyst **Parameter Result Qualifier** Units RL**Factor Prepared** Analyzed MDL Total Metals - Mansfield Lab for sample(s): 02 Batch: WG1688660-1 Arsenic, Total ND mg/l 0.0050 0.0019 10/06/22 12:46 1,6010D ΝB 1 09/17/22 07:02

**Prep Information** 

Digestion Method: EPA 3005A



**Project Name:** STEELWINDS ANNUAL/ SEMIANNUAL

Lab Number: L2250076

**Project Number:** 03.0033579.15

Report Date:

10/12/22

Parameter	LCS %Recovery Qu	LCSD al %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits	
Total Metals - Mansfield Lab Associated sa	ample(s): 02 Batch: WG1	688660-2						
Arsenic. Total	105	-		80-120	_			



# Matrix Spike Analysis Batch Quality Control

Project Name: STEELWINDS ANNUAL/ SEMIANNUAL

**Project Number:** 03.0033579.15

Lab Number:

L2250076

Report Date:

10/12/22

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	Qua	MSD   Found	MSD %Recovery	Qual	Recovery Limits	RPD	Qual	RPD Limits
Total Metals - Mansfield La	b Associated san	nple(s): 02	QC Batch	ID: WG168866	0-3	QC Sample	: L2249023-01	Clier	nt ID: MS Sa	ample		
Arsenic, Total	ND	0.12	0.131	109		-	-		75-125	-		20



Serial\_No:10122213:28 *Lab Number:* L2250076

Project Name: STEELWINDS ANNUAL/ SEMIANNUAL

**Project Number:** 03.0033579.15 **Report Date:** 10/12/22

# Sample Receipt and Container Information

Were project specific reporting limits specified?

**Cooler Information** 

Cooler Custody Seal

A Absent

Container Information		Initial	Final	Temp			Frozen		
Container ID	Container Type	Cooler	рН	рН	deg C	Pres	Seal	Date/Time	Analysis(*)
L2250076-01A	Vial HCl preserved	Α	NA		2.4	Υ	Absent		NYCP51-8260(14)
L2250076-01B	Vial HCI preserved	Α	NA		2.4	Υ	Absent		NYCP51-8260(14)
L2250076-01C	Vial HCI preserved	Α	NA		2.4	Υ	Absent		NYCP51-8260(14)
L2250076-01D	Amber 1000ml unpreserved	Α	7	7	2.4	Υ	Absent		A2-SVOC-8270(7)
L2250076-01E	Amber 1000ml unpreserved	Α	7	7	2.4	Υ	Absent		A2-SVOC-8270(7)
L2250076-02A	Vial HCI preserved	Α	NA		2.4	Υ	Absent		NYCP51-8260(14)
L2250076-02B	Vial HCI preserved	Α	NA		2.4	Υ	Absent		NYCP51-8260(14)
L2250076-02C	Vial HCl preserved	Α	NA		2.4	Υ	Absent		NYCP51-8260(14)
L2250076-02D	Plastic 250ml HNO3 preserved	Α	<2	<2	2.4	Υ	Absent		AS-TI(180)
L2250076-02E	Amber 1000ml unpreserved	Α	7	7	2.4	Υ	Absent		A2-SVOC-8270(7)
L2250076-02F	Amber 1000ml unpreserved	Α	7	7	2.4	Υ	Absent		A2-SVOC-8270(7)



**Project Name:** Lab Number: STEELWINDS ANNUAL/ SEMIANNUAL L2250076 03.0033579.15 **Report Date: Project Number:** 10/12/22

### **GLOSSARY**

#### **Acronyms**

**EDL** 

LOQ

MS

RPD

DL - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

- Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis

of PAHs using Solid-Phase Microextraction (SPME).

**EMPC** - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case estimate of the concentration.

**EPA** Environmental Protection Agency.

LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LCSD Laboratory Control Sample Duplicate: Refer to LCS.

LFB - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LOD - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

- Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats

Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats

MDI - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

- Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.

MSD - Matrix Spike Sample Duplicate: Refer to MS.

NA - Not Applicable.

NC - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.

NDPA/DPA - N-Nitrosodiphenylamine/Diphenylamine.

NI - Not Ignitable.

NP - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.

- No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile NR Organic TIC only requests.

RL - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

- Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.

SRM - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples.

STLP - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.

TEF - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.

TEO - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF and then summing the resulting values.

TIC - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.

Report Format: DU Report with 'J' Qualifiers



Project Name:STEELWINDS ANNUAL/ SEMIANNUALLab Number:L2250076Project Number:03.0033579.15Report Date:10/12/22

#### **Footnotes**

1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

#### **Terms**

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Chlordane: The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA,this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

Difference: With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

Final pH: As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

Frozen Date/Time: With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

Gasoline Range Organics (GRO): Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

Initial pH: As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

PAH Total: With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benz(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

PFAS Total: With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

Total: With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

# Data Qualifiers

- A Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentrations of the analyte at less than ten times (10x) the concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G The concentration may be biased high due to matrix interferences (i.e, co-elution) with non-target compound(s). The result should be considered estimated.
- H The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I The lower value for the two columns has been reported due to obvious interference.
- Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit
   (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively

Report Format: DU Report with 'J' Qualifiers



Project Name:STEELWINDS ANNUAL/ SEMIANNUALLab Number:L2250076Project Number:03.0033579.15Report Date:10/12/22

#### **Data Qualifiers**

Identified Compounds (TICs).

- $\label{eq:main_main_model} \textbf{M} \qquad \text{-Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.}$
- ND Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.
- NJ Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- ${f P}$  The RPD between the results for the two columns exceeds the method-specified criteria.
- Q The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- **RE** Analytical results are from sample re-extraction.
- S Analytical results are from modified screening analysis.
- The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits.
   (Applicable to MassDEP DW Compliance samples only.)

Report Format: DU Report with 'J' Qualifiers



Project Name:STEELWINDS ANNUAL/ SEMIANNUALLab Number:L2250076Project Number:03.0033579.15Report Date:10/12/22

#### REFERENCES

Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

#### **LIMITATION OF LIABILITIES**

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Alpha Analytical, Inc.
Facility: Company-wide

Department: Quality Assurance

Title: Certificate/Approval Program Summary

\_ ID No.:**17873** 

Page 1 of 1

Revision 19 Published Date: 4/2/2021 1:14:23 PM

#### Certification Information

#### The following analytes are not included in our Primary NELAP Scope of Accreditation:

#### Westborough Facility

EPA 624/624.1: m/p-xylene, o-xylene, Naphthalene

EPA 625/625.1: alpha-Terpineol

**EPA 8260C/8260D:** NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; SCM: lodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene

4-Ethyltoluene.

EPA 8270D/8270E: NPW: Dimethylnaphthalene,1,4-Diphenylhydrazine, alpha-Terpineol; SCM: Dimethylnaphthalene,1,4-Diphenylhydrazine.

SM4500: NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO2, NO3.

#### Mansfield Facility

SM 2540D: TSS

EPA 8082A: NPW: PCB: 1, 5, 31, 87,101, 110, 141, 151, 153, 180, 183, 187.

**EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

Biological Tissue Matrix: EPA 3050B

#### The following analytes are included in our Massachusetts DEP Scope of Accreditation

#### Westborough Facility:

#### **Drinking Water**

EPA 300.0: Chloride, Nitrate-N, Fluoride, Sulfate; EPA 353.2: Nitrate-N, Nitrite-N; SM4500NO3-F: Nitrate-N, Nitrite-N; SM4500F-C, SM4500CN-CE,

EPA 180.1, SM2130B, SM4500CI-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B

EPA 332: Perchlorate; EPA 524.2: THMs and VOCs; EPA 504.1: EDB, DBCP.

Microbiology: SM9215B; SM9223-P/A, SM9223B-Colilert-QT,SM9222D.

#### Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, EPA 350.1: Ammonia-N, LACHAT 10-107-06-1-B: Ammonia-N, EPA 351.1, SM4500NO3-F, EPA 353.2: Nitrate-N, SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300: Chloride, Sulfate, Nitrate. EPA 624.1: Volatile Halocarbons & Aromatics.

EPA 608.3: Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan II, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

EPA 625.1: SVOC (Acid/Base/Neutral Extractables), EPA 600/4-81-045: PCB-Oil.

Microbiology: SM9223B-Colilert-QT; Enterolert-QT, SM9221E, EPA 1600, EPA 1603, SM9222D.

#### Mansfield Facility:

#### **Drinking Water**

**EPA 200.7:** Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. **EPA 200.8:** Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. **EPA 245.1** Hg. **EPA 522, EPA 537.1.** 

#### Non-Potable Water

EPA 200.7: Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.

EPA 200.8: Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.

EPA 245.1 Hg.

SM2340B

For a complete listing of analytes and methods, please contact your Alpha Project Manager.

Document Type: Form Pre-Qualtrax Document ID: 08-113

Westborough, MA 01581 8 Walkup Dr. TEL: 508-898-9220	NEW YORK CHAIN OF CUSTODY Mansfield, MA 02048 320 Forbes Blvd TEL: 508-822-9300	Service Centers Mahwah, NJ 07430: 35 Whitne Albany, NY 12205: 14 Walker Tonawanda, NY 14150: 275 Co Project Information Project Name: STEEL	Way coper Ave, Suite 1			of j	Deliv	Date R in L erables	ab	9	15 ASP-			ALPHA Job # L 2 2 5 00 7 4  Billing Information  Same as Client Info
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H = Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub>	E = Encore D = BOD Bottle	1230		9-14-22 / 1	250	force			(eni		14/22 5/2	12:	50	THIS COC, THE CLIENT HAS READ AND AGREES TO BE BOUND BY ALPHA'S TERMS & CONDITIONS.
Form No: 01-25 HC (rev. 30	-Sept-2013)													(See reverse side.)



#### ANALYTICAL REPORT

Lab Number: L2250575

Client: GZA GeoEnvironmental of New York

300 Pearl Street

Suite 700

Buffalo, NY 14202

ATTN: Dan Troy

Phone: (716) 844-7050

Project Name: STEELWINDS ANNUAL/SEMI-ANNUAL

Project Number: 03.0033579.15

Report Date: 10/13/22

The original project report/data package is held by Alpha Analytical. This report/data package is paginated and should be reproduced only in its entirety. Alpha Analytical holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0574), IL (200077), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #P330-17-00196).

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

**Project Number:** 03.0033579.15

**Lab Number:** L2250575 **Report Date:** 10/13/22

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2250575-01	MWN-03D-091522	WATER	LACKAWANNA NY	09/15/22 08:30	09/15/22
L2250575-02	MWN-03B-091522	WATER	LACKAWANNA NY	09/15/22 09:45	09/15/22
L2250575-03	MWN-03-091522	WATER	LACKAWANNA NY	09/15/22 10:30	09/15/22
L2250575-04	MWN-02D-091522	WATER	LACKAWANNA NY	09/15/22 13:45	09/15/22
L2250575-05	MWN-04-091522	WATER	LACKAWANNA NY	09/15/22 14:50	09/15/22
L2250575-06	TRIP BLANK	WATER	LACKAWANNA NY	09/15/22 00:00	09/15/22



L2250575

Project Name: STEELWINDS ANNUAL/SEMI-ANNUAL Lab Number:

**Project Number:** 03.0033579.15 **Report Date:** 10/13/22

#### **Case Narrative**

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments, solids and tissues are reported on a dry weight basis unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

HOLD POLICY - For samples submitted on hold, Alpha's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Alpha Project Manager and made arrangements for Alpha to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.	



Project Name: STEELWINDS ANNUAL/SEMI-ANNUAL Lab Number: L2250575

**Project Number:** 03.0033579.15 **Report Date:** 10/13/22

**Case Narrative (continued)** 

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

Title: Technical Director/Representative Date: 10/13/22

600, Sharow Kelly Stenstrom

### **ORGANICS**



### **VOLATILES**



L2250575

**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

MWN-03D-091522

LACKAWANNA NY

L2250575-01

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Date Collected: 09/15/22 08:30

Lab Number:

Report Date:

10/13/22

Date Received: 09/15/22 Field Prep: Not Specified

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/26/22 15:59

Analyst: MKS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - West	borough Lab					
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	ND		ug/l	2.5	0.70	1
o-Xylene	ND		ug/l	2.5	0.70	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	ND		ug/l	2.5	0.70	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	101	70-130	
Toluene-d8	100	70-130	
4-Bromofluorobenzene	95	70-130	
Dibromofluoromethane	119	70-130	



L2250575

10/13/22

**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

L2250575-03

MWN-03-091522

LACKAWANNA NY

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Date Collected: 09/15/22 10:30

Lab Number:

Report Date:

Date Received: 09/15/22
Field Prep: Not Specified

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Matrix: Water
Analytical Method: 1,8260C
Analytical Date: 09/26/22 16:24

Analyst: MKS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westbe	orough Lab					
Benzene	11		ug/l	0.50	0.16	1
Toluene	2.4	J	ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	1.6	J	ug/l	2.5	0.70	1
o-Xylene	1.7	J	ug/l	2.5	0.70	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	25		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	0.97	J	ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	ND		ug/l	2.5	0.70	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	101	70-130	
Toluene-d8	98	70-130	
4-Bromofluorobenzene	94	70-130	
Dibromofluoromethane	115	70-130	



**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

L2250575

Lab Number:

Report Date: 10/13/22

Lab ID: L2250575-05 Date Collected: 09/15/22 14:50

Client ID: Date Received: 09/15/22 MWN-04-091522 Field Prep: Sample Location: Not Specified LACKAWANNA NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/26/22 16:48

Analyst: MKS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westl	oorough Lab					
Benzene	0.51		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	ND		ug/l	2.5	0.70	1
o-Xylene	ND		ug/l	2.5	0.70	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	16		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	ND		ug/l	2.5	0.70	1

Surrogate	% Recovery	Qualifier	Acceptance Criteria	
1,2-Dichloroethane-d4	104		70-130	
Toluene-d8	97		70-130	
4-Bromofluorobenzene	91		70-130	
Dibromofluoromethane	123		70-130	



L2250575

10/13/22

**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Lab Number:

Report Date:

Lab ID: L2250575-06 Date Collected: 09/15/22 00:00

Client ID: TRIP BLANK Date Received: 09/15/22 Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Matrix: Water
Analytical Method: 1,8260C
Analytical Date: 10/05/22 11:02

Analyst: PD

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - West	oorough Lab					
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	ND		ug/l	2.5	0.70	1
o-Xylene	ND		ug/l	2.5	0.70	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	ND		ug/l	2.5	0.70	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	96	70-130	
Toluene-d8	96	70-130	
4-Bromofluorobenzene	98	70-130	
Dibromofluoromethane	98	70-130	



Project Name: STEELWINDS ANNUAL/SEMI-ANNUAL Lab Number: L2250575

**Project Number:** 03.0033579.15 **Report Date:** 10/13/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/26/22 11:03

Analyst: PD

Parameter	Result	Qualifier Uni	ts R	L	MDL
Volatile Organics by GC/MS - Wes	tborough Lab	for sample(s):	01,03,05	Batch:	WG1692636-5
Benzene	ND	ug	/I 0.5	50	0.16
Toluene	ND	ug	/l 2.	5	0.70
Ethylbenzene	ND	ug	/l 2.	5	0.70
Methyl tert butyl ether	ND	ug	/l 2.	5	0.70
p/m-Xylene	ND	ug	/l 2.	5	0.70
o-Xylene	ND	ug	/I 2.	5	0.70
n-Butylbenzene	ND	ug	/I 2.	5	0.70
sec-Butylbenzene	ND	ug	/I 2.	5	0.70
tert-Butylbenzene	ND	ug	/l 2.	5	0.70
Isopropylbenzene	ND	ug	/l 2.	5	0.70
p-Isopropyltoluene	ND	ug	/l 2.	5	0.70
Naphthalene	ND	ug	/l 2.	5	0.70
n-Propylbenzene	ND	ug	/l 2.	5	0.70
1,3,5-Trimethylbenzene	ND	ug	/l 2.	5	0.70
1,2,4-Trimethylbenzene	ND	ug	/l 2.	5	0.70

		Acceptance	
Surrogate	%Recovery Qualifi	ier Criteria	
1,2-Dichloroethane-d4	101	70-130	
Toluene-d8	99	70-130	
4-Bromofluorobenzene	95	70-130	
Dibromofluoromethane	116	70-130	



Project Name: STEELWINDS ANNUAL/SEMI-ANNUAL Lab Number: L2250575

**Project Number:** 03.0033579.15 **Report Date:** 10/13/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 10/05/22 10:39

Analyst: PD

Parameter	Result	Qualifier Units	s RL	MDL	
Volatile Organics by GC/MS	- Westborough Lab	for sample(s):	06 Batch:	WG1696331-5	
Benzene	ND	ug/l	0.50	0.16	
Toluene	ND	ug/l	2.5	0.70	
Ethylbenzene	ND	ug/l	2.5	0.70	
Methyl tert butyl ether	ND	ug/l	2.5	0.70	
p/m-Xylene	ND	ug/l	2.5	0.70	
o-Xylene	ND	ug/l	2.5	0.70	
n-Butylbenzene	ND	ug/l	2.5	0.70	
sec-Butylbenzene	ND	ug/l	2.5	0.70	
tert-Butylbenzene	ND	ug/l	2.5	0.70	
Isopropylbenzene	ND	ug/l	2.5	0.70	
p-Isopropyltoluene	ND	ug/l	2.5	0.70	
Naphthalene	ND	ug/l	2.5	0.70	
n-Propylbenzene	ND	ug/l	2.5	0.70	
1,3,5-Trimethylbenzene	ND	ug/l	2.5	0.70	
1,2,4-Trimethylbenzene	ND	ug/l	2.5	0.70	

		Acceptance	
Surrogate	%Recovery Quality	fier Criteria	
1,2-Dichloroethane-d4	96	70-130	
Toluene-d8	97	70-130	
4-Bromofluorobenzene	98	70-130	
Dibromofluoromethane	97	70-130	



**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

**Project Number:** 03.0033579.15

Lab Number: L2250575

**Report Date:** 10/13/22

	LCS	0	LCSD	%Recovery	555	RPD	
arameter	%Recovery	Qual	%Recovery	Qual Limits	RPD	Qual Limits	
platile Organics by GC/MS - Westborough L	ab Associated	sample(s):	01,03,05 Batch:	WG1692636-3 WG1692636-	4		
Benzene	110		120	70-130	9	20	
Toluene	110		110	70-130	0	20	
Ethylbenzene	110		110	70-130	0	20	
Methyl tert butyl ether	85		89	63-130	5	20	
p/m-Xylene	105		110	70-130	5	20	
o-Xylene	100		110	70-130	10	20	
n-Butylbenzene	110		120	53-136	9	20	
sec-Butylbenzene	110		110	70-130	0	20	
tert-Butylbenzene	87		100	70-130	14	20	
Isopropylbenzene	100		110	70-130	10	20	
p-Isopropyltoluene	100		110	70-130	10	20	
Naphthalene	85		93	70-130	9	20	
n-Propylbenzene	110		120	69-130	9	20	
1,3,5-Trimethylbenzene	100		100	64-130	0	20	
1,2,4-Trimethylbenzene	100		100	70-130	0	20	

Surrogate	LCS	LCSD	Acceptance
	%Recovery Qual	%Recovery Qual	Criteria
1,2-Dichloroethane-d4	97	99	70-130
Toluene-d8	100	100	70-130
4-Bromofluorobenzene	97	95	70-130
Dibromofluoromethane	100	105	70-130



**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

**Project Number:** 03.0033579.15

Lab Number: L2250575

**Report Date:** 10/13/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
olatile Organics by GC/MS - Westborough L	ab Associated	sample(s): 06	Batch: WG1	1696331-3	WG1696331-4			
Benzene	110		110		70-130	0		20
Toluene	100		100		70-130	0		20
Ethylbenzene	99		99		70-130	0		20
Methyl tert butyl ether	99		100		63-130	1		20
p/m-Xylene	100		100		70-130	0		20
o-Xylene	100		100		70-130	0		20
n-Butylbenzene	98		99		53-136	1		20
sec-Butylbenzene	98		100		70-130	2		20
tert-Butylbenzene	98		99		70-130	1		20
Isopropylbenzene	99		99		70-130	0		20
p-Isopropyltoluene	97		100		70-130	3		20
Naphthalene	83		92		70-130	10		20
n-Propylbenzene	99		99		69-130	0		20
1,3,5-Trimethylbenzene	96		97		64-130	1		20
1,2,4-Trimethylbenzene	97		98		70-130	1		20

	LCS	LCSD	Acceptance	
Surrogate	%Recovery Qual	%Recovery Qual	Criteria	
1,2-Dichloroethane-d4	94	95	70-130	
Toluene-d8	96	96	70-130	
4-Bromofluorobenzene	99	99	70-130	
Dibromofluoromethane	101	100	70-130	



### **SEMIVOLATILES**



L2250575

10/13/22

**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

L2250575-01

MWN-03D-091522

LACKAWANNA NY

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Date Collected: 09/15/22 08:30

Date Received: 09/15/22

Lab Number:

Report Date:

Field Prep: Not Specified

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Matrix: Water Analytical Method: 1,8270D Analytical Date: 09/19/22 16:08

Analyst: PS

Extraction M	lethod: El	PA 3510C
Extraction D	ate: 09	9/17/22 08:30

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS - Ma	nsfield Lab					
bis(2-Chloroethyl)ether	ND		ug/l	0.495	0.092	1
1,3-Dichlorobenzene	ND		ug/l	0.495	0.078	1
1,4-Dichlorobenzene	ND		ug/l	0.495	0.082	1
1,2-Dichlorobenzene	ND		ug/l	0.495	0.067	1
Benzyl alcohol	ND		ug/l	0.495	0.122	1
bis(2-chloroisopropyl)ether	ND		ug/l	0.495	0.107	1
Acetophenone	ND		ug/l	0.990	0.205	1
Hexachloroethane	ND		ug/l	0.495	0.101	1
Nitrobenzene	ND		ug/l	0.495	0.101	1
Isophorone	ND		ug/l	0.495	0.125	1
bis(2-Chloroethoxy)methane	ND		ug/l	0.495	0.085	1
1,2,4-Trichlorobenzene	ND		ug/l	0.495	0.095	1
Naphthalene	ND		ug/l	0.495	0.087	1
4-Chloroaniline	ND		ug/l	0.495	0.127	1
Hexachlorobutadiene	ND		ug/l	0.495	0.085	1
2-Methylnaphthalene	ND		ug/l	0.495	0.090	1
1,2,4,5-Tetrachlorobenzene	ND		ug/l	0.495	0.079	1
Hexachlorocyclopentadiene	ND		ug/l	0.495	0.151	1
Biphenyl	ND		ug/l	0.495	0.110	1
2-Chloronaphthalene	ND		ug/l	0.495	0.089	1
2-Nitroaniline	ND		ug/l	0.495	0.137	1
Acenaphthylene	ND		ug/l	0.495	0.111	1
Dimethylphthalate	ND		ug/l	0.495	0.116	1
2,6-Dinitrotoluene	ND		ug/l	0.495	0.166	1
Acenaphthene	0.536		ug/l	0.495	0.095	1
3-Nitroaniline	ND		ug/l	0.495	0.110	1
Dibenzofuran	ND		ug/l	0.495	0.090	1
2,4-Dinitrotoluene	ND		ug/l	0.495	0.161	1



10/13/22

**Project Name:** Lab Number: STEELWINDS ANNUAL/SEMI-ANNUAL L2250575

**Project Number:** 03.0033579.15

L2250575-01

**SAMPLE RESULTS** 

Date Collected: 09/15/22 08:30

Report Date:

Date Received: 09/15/22 Client ID: MWN-03D-091522 Sample Location: Field Prep: LACKAWANNA NY Not Specified

Sample Depth:

Lab ID:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS	- Mansfield Lab					
Fluorene	0.187	J	ug/l	0.495	0.103	1
Diethylphthalate	ND		ug/l	0.495	0.178	1
4-Nitroaniline	ND		ug/l	0.495	0.111	1
n-Nitrosodiphenylamine	ND		ug/l	0.495	0.071	1
Hexachlorobenzene	ND		ug/l	0.495	0.121	1
Phenanthrene	0.434	J	ug/l	0.495	0.110	1
Anthracene	ND		ug/l	0.495	0.136	1
Carbazole	ND		ug/l	0.495	0.142	1
Di-n-butylphthalate	ND		ug/l	0.495	0.099	1
Fluoranthene	ND		ug/l	0.495	0.154	1
Pyrene	ND		ug/l	0.495	0.168	1
Butylbenzylphthalate	ND		ug/l	0.495	0.084	1
3,3'-Dichlorobenzidine	ND		ug/l	0.495	0.191	1
Benz(a)anthracene	ND		ug/l	0.495	0.182	1
Chrysene	ND		ug/l	0.495	0.140	1
bis(2-Ethylhexyl)phthalate	0.376	J	ug/l	0.495	0.080	1
Di-n-octylphthalate	ND		ug/l	0.990	0.078	1
Benzo(b)fluoranthene	ND		ug/l	0.495	0.065	1
Benzo(k)fluoranthene	ND		ug/l	0.495	0.159	1
Benzo(a)pyrene	ND		ug/l	0.495	0.060	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.495	0.089	1
Dibenz(a,h)anthracene	ND		ug/l	0.495	0.064	1
Benzo(g,h,i)perylene	ND		ug/l	0.495	0.108	1

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	36	15-115
Phenol-d5	26	15-115
Nitrobenzene-d5	64	30-130
2-Fluorobiphenyl	61	30-130
2,4,6-Tribromophenol	82	15-115
Terphenyl-d14	66	30-130



L2250575

10/13/22

09/17/22 08:30

Project Name: STEELWINDS ANNUAL/SEMI-ANNUAL

L2250575-03

MWN-03-091522

LACKAWANNA NY

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Date Collected: 09/15/22 10:30

Date Received: 09/15/22

Extraction Method: EPA 3510C

Lab Number:

Report Date:

**Extraction Date:** 

Field Prep: Not Specified

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Matrix: Water
Analytical Method: 1,8270D

Analytical Date: 09/19/22 16:38

Analyst: PS

Semivolatile Organics by GC/MS - Mansfield Lab           bis (2-Chloroethyl)ether         ND         ug/l         0.490         0.091         1           1,3-Dichlorobenzene         ND         ug/l         0.490         0.077         1           1,4-Dichlorobenzene         ND         ug/l         0.490         0.081         1           1,2-Dichlorobenzene         0.115         J         ug/l         0.490         0.067         1           Benzyl alcohol         ND         ug/l         0.490         0.102         1           Benzyl alcohol         ND         ug/l         0.490         0.106         1           Acetophenone         0.308         J         ug/l         0.490         0.106         1           Hexachloroethane         ND         ug/l         0.490         0.100         1           Nitrobenzene         ND         ug/l         0.490         0.100         1           Sleige-Chloroethoxylmethane         ND         ug/l         0.490         0.084         1           1,2-4-Trichlorobenzene         ND         ug/l         0.490         0.084         1           1,2-4-Trichlorobenzene         ND         ug/l         0.490         0.084 <th>Parameter</th> <th>Result</th> <th>Qualifier</th> <th>Units</th> <th>RL</th> <th>MDL</th> <th>Dilution Factor</th>	Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
1.3-Dichlorobenzene ND ug/l 0.490 0.077 1 1.4-Dichlorobenzene ND ug/l 0.490 0.081 1 1.2-Dichlorobenzene ND ug/l 0.490 0.081 1 1.2-Dichlorobenzene 0.115 J ug/l 0.490 0.067 1 1.2-Dichlorobenzene 0.115 J ug/l 0.490 0.067 1 1.2-Dichlorobenzene 0.115 J ug/l 0.490 0.067 1 1.2-Dichlorobenzene ND ug/l 0.490 0.120 1 1.2-Dichlorobenzene ND ug/l 0.490 0.106 1 1.2-Dichlorobenzene ND ug/l 0.490 0.106 1 1.2-Dichlorobenzene ND ug/l 0.490 0.100 1 1.2-Dichlorobenzene ND ug/l 0.490 0.084 1 1.2-L-Trichlorobenzene ND ug/l 0.490 0.084 1 1.2-L-Trichlorobenzene ND ug/l 0.490 0.084 1 1.2-L-Trichlorobenzene ND ug/l 0.490 0.086 1 1-Dichlorobenzene ND ug/l 0.490 0.086 1 1-Dichlorobenzene ND ug/l 0.490 0.084 1 1.2-L-Trichlorobenzene ND ug/l 0.490 0.084 1 1.2-L-Trichlorobenzene ND ug/l 0.490 0.084 1 1.2-L-Trichlorobenzene ND ug/l 0.490 0.086 1 1-Dichlorobenzene ND ug/l 0.490 0.089 1 1-Dichlorobenzene ND ug/l 0.490 0.089 1 1-Dichlorobenzene ND ug/l 0.490 0.089 1 1-Dichlorobenzene ND ug/l 0.490 0.078 1 1-Dichlorobenzene ND ug/l 0.490 0.078 1 1-Dichlorobenzene ND ug/l 0.490 0.088 1 1-Dichlorobenzene ND ug/l 0.490 0.0100	Semivolatile Organics by GC/MS - Ma	ansfield Lab					
1.3-Dichlorobenzene         ND         ug/l         0.490         0.077         1           1.4-Dichlorobenzene         ND         ug/l         0.490         0.081         1           1.2-Dichlorobenzene         0.115         J         ug/l         0.490         0.067         1           Benzyl alcohol         ND         ug/l         0.490         0.120         1           Sie(2-chlorospropylether         ND         ug/l         0.490         0.106         1           Acetophenone         0.308         J         ug/l         0.490         0.100         1           Hexachloroethane         ND         ug/l         0.490         0.100         1           Nitrobenzene         ND         ug/l         0.490         0.100         1           Isophorone         ND         ug/l         0.490         0.100         1           Isophorone         ND         ug/l         0.490         0.084         1           Isophorone         ND         ug/l         0.490         0.084         1           Isophorone         ND         ug/l         0.490         0.084         1           Isophorone         ND         ug/l         0.490<	bis(2-Chloroethyl)ether	ND		ug/l	0.490	0.091	1
1,4-Dichlorobenzene   ND	1,3-Dichlorobenzene	ND			0.490	0.077	1
Benzyl alcohol         ND         ug/l         0.490         0.120         1           bis(2-chloroisopropyl)ether         ND         ug/l         0.490         0.106         1           Acetophenone         0.308         J         ug/l         0.490         0.106         1           Hexachloroethane         ND         ug/l         0.490         0.100         1           Nitrobenzene         ND         ug/l         0.490         0.100         1           Isophorone         ND         ug/l         0.490         0.100         1           Isophorone         ND         ug/l         0.490         0.124         1           Isophorone         ND         ug/l         0.490         0.084         1           Lg-4-Trichlorobarzene         ND         ug/l         0.490         0.086         1           Hexachlorobutadiene         ND         ug/l         0.490         0.075 <td>1,4-Dichlorobenzene</td> <td>ND</td> <td></td> <td></td> <td>0.490</td> <td>0.081</td> <td>1</td>	1,4-Dichlorobenzene	ND			0.490	0.081	1
Dis(2-chloroispropyl)ether   ND	1,2-Dichlorobenzene	0.115	J	ug/l	0.490	0.067	1
Acetophenone         0.308         J         ug/l         0.980         0.203         1           Hexachloroethane         ND         ug/l         0.490         0.100         1           Nitrobenzene         ND         ug/l         0.490         0.100         1           Isophorone         ND         ug/l         0.490         0.124         1           bis(2-Chloroethoxy)methane         ND         ug/l         0.490         0.084         1           1,2.4-Trichlorobenzene         ND         ug/l         0.490         0.094         1           Naphthalene         15.0         ug/l         0.490         0.096         1           4-Chloroaniline         ND         ug/l         0.490         0.086         1           Hexachlorobutadiene         ND         ug/l         0.490         0.084         1           Hexachlorobutadiene         ND         ug/l         0.490         0.084         1           1,2.4,5-Tetrachlorobenzene         ND         ug/l         0.490         0.078         1           Hexachlorocyclopentadiene         ND         ug/l         0.490         0.150         1           Biphenyl         0.715         ug/l	Benzyl alcohol	ND		ug/l	0.490	0.120	1
Hexachloroethane         ND         ug/l         0.490         0.100         1           Nitrobenzene         ND         ug/l         0.490         0.100         1           Isophorone         ND         ug/l         0.490         0.124         1           bis(2-Chloroethoxy)methane         ND         ug/l         0.490         0.084         1           1,2,4-Trichlorobenzene         ND         ug/l         0.490         0.094         1           Naphthalene         15.0         ug/l         0.490         0.086         1           4-Chloroaniline         ND         ug/l         0.490         0.086         1           Hexachlorobutadiene         ND         ug/l         0.490         0.086         1           4-Chloroaniline         ND         ug/l         0.490         0.084         1           2-Methylnaphthalene         3.03         ug/l         0.490         0.089         1           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.490         0.078         1           Hexachlorocyclopentadiene         ND         ug/l         0.490         0.150         1           Biphenyl         0.715         ug/l         0.49	bis(2-chloroisopropyl)ether	ND		ug/l	0.490	0.106	1
NItrobenzene         ND         ug/l         0.490         0.100         1           Isophorone         ND         ug/l         0.490         0.124         1           bis(2-Chloroethoxy)methane         ND         ug/l         0.490         0.084         1           1,2,4-Trichlorobenzene         ND         ug/l         0.490         0.094         1           Naphthalene         15.0         ug/l         0.490         0.086         1           4-Chloroaniline         ND         ug/l         0.490         0.086         1           4-Chloroaniline         ND         ug/l         0.490         0.086         1           Hexachlorobutadiene         ND         ug/l         0.490         0.086         1           4-Chloroaniline         ND         ug/l         0.490         0.084         1           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.490         0.078         1           Hexachlorocyclopentadiene         ND         ug/l         0.490         0.150         1           Biphenyl         0.715         ug/l         0.490         0.109         1           2-Chloronaphthalene         ND         ug/l         0.490 </td <td>Acetophenone</td> <td>0.308</td> <td>J</td> <td>ug/l</td> <td>0.980</td> <td>0.203</td> <td>1</td>	Acetophenone	0.308	J	ug/l	0.980	0.203	1
Sophorone   ND   ug/l   0.490   0.124   1	Hexachloroethane	ND		ug/l	0.490	0.100	1
Discretion   ND   Ug/l   0.490   0.084   1	Nitrobenzene	ND		ug/l	0.490	0.100	1
1,2,4-Trichlorobenzene         ND         ug/l         0.490         0.094         1           Naphthalene         15.0         ug/l         0.490         0.086         1           4-Chloroaniline         ND         ug/l         0.490         0.125         1           Hexachlorobutadiene         ND         ug/l         0.490         0.084         1           2-Methylnaphthalene         3.03         ug/l         0.490         0.089         1           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.490         0.078         1           Hexachlorocyclopentadiene         ND         ug/l         0.490         0.150         1           Biphenyl         0.715         ug/l         0.490         0.109         1           2-Chloronaphthalene         ND         ug/l         0.490         0.109         1           2-Nitroaniline         ND         ug/l         0.490         0.135         1           Acenaphthylpene         2.70         ug/l         0.490         0.115         1           2,6-Dinitrotoluene         ND         ug/l         0.490         0.165         1           Acenaphthene         1.54         ug/l         0	Isophorone	ND		ug/l	0.490	0.124	1
Naphthalene         15.0         ug/l         0.490         0.086         1           4-Chloroaniline         ND         ug/l         0.490         0.125         1           Hexachlorobutadiene         ND         ug/l         0.490         0.084         1           2-Methylnaphthalene         3.03         ug/l         0.490         0.089         1           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.490         0.078         1           Hexachlorocyclopentadiene         ND         ug/l         0.490         0.150         1           Biphenyl         0.715         ug/l         0.490         0.109         1           2-Chloronaphthalene         ND         ug/l         0.490         0.109         1           2-Nitroaniline         ND         ug/l         0.490         0.135         1           Acenaphthylene         2.70         ug/l         0.490         0.115         1           2,6-Dinitrotoluene         ND         ug/l         0.490         0.165         1           Acenaphthene         1.54         ug/l         0.490         0.094         1           3-Nitroaniline         ND         ug/l         0.490	bis(2-Chloroethoxy)methane	ND		ug/l	0.490	0.084	1
4-Chloroaniline         ND         ug/l         0.490         0.125         1           Hexachlorobutadiene         ND         ug/l         0.490         0.084         1           2-Methylnaphthalene         3.03         ug/l         0.490         0.089         1           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.490         0.078         1           Hexachlorocyclopentadiene         ND         ug/l         0.490         0.150         1           Biphenyl         0.715         ug/l         0.490         0.109         1           2-Chloronaphthalene         ND         ug/l         0.490         0.109         1           2-Nitroaniline         ND         ug/l         0.490         0.135         1           Acenaphthylene         2.70         ug/l         0.490         0.110         1           Dimethylphthalate         ND         ug/l         0.490         0.115         1           2,6-Dinitrotoluene         ND         ug/l         0.490         0.165         1           Acenaphthene         1.54         ug/l         0.490         0.094         1           3-Nitroaniline         ND         ug/l         0.490<	1,2,4-Trichlorobenzene	ND		ug/l	0.490	0.094	1
Hexachlorobutadiene   ND   ug/l   0.490   0.084   1	Naphthalene	15.0		ug/l	0.490	0.086	1
2-Methylnaphthalene       3.03       ug/l       0.490       0.089       1         1,2,4,5-Tetrachlorobenzene       ND       ug/l       0.490       0.078       1         Hexachlorocyclopentadiene       ND       ug/l       0.490       0.150       1         Biphenyl       0.715       ug/l       0.490       0.109       1         2-Chloronaphthalene       ND       ug/l       0.490       0.088       1         2-Nitroaniline       ND       ug/l       0.490       0.135       1         Acenaphthylene       2.70       ug/l       0.490       0.110       1         Dimethylphthalate       ND       ug/l       0.490       0.115       1         2,6-Dinitrotoluene       ND       ug/l       0.490       0.165       1         Acenaphthene       1.54       ug/l       0.490       0.094       1         3-Nitroaniline       ND       ug/l       0.490       0.109       1         Dibenzofuran       2.92       ug/l       0.490       0.089       1	4-Chloroaniline	ND		ug/l	0.490	0.125	1
1,2,4,5-Tetrachlorobenzene       ND       ug/l       0.490       0.078       1         Hexachlorocyclopentadiene       ND       ug/l       0.490       0.150       1         Biphenyl       0.715       ug/l       0.490       0.109       1         2-Chloronaphthalene       ND       ug/l       0.490       0.088       1         2-Nitroaniline       ND       ug/l       0.490       0.135       1         Acenaphthylene       2.70       ug/l       0.490       0.110       1         Dimethylphthalate       ND       ug/l       0.490       0.115       1         2,6-Dinitrotoluene       ND       ug/l       0.490       0.165       1         Acenaphthene       1.54       ug/l       0.490       0.094       1         3-Nitroaniline       ND       ug/l       0.490       0.109       1         Dibenzofuran       2.92       ug/l       0.490       0.089       1	Hexachlorobutadiene	ND		ug/l	0.490	0.084	1
Hexachlorocyclopentadiene   ND   ug/l   0.490   0.150   1	2-Methylnaphthalene	3.03		ug/l	0.490	0.089	1
Biphenyl         0.715         ug/l         0.490         0.109         1           2-Chloronaphthalene         ND         ug/l         0.490         0.088         1           2-Nitroaniline         ND         ug/l         0.490         0.135         1           Acenaphthylene         2.70         ug/l         0.490         0.110         1           Dimethylphthalate         ND         ug/l         0.490         0.115         1           2,6-Dinitrotoluene         ND         ug/l         0.490         0.165         1           Acenaphthene         1.54         ug/l         0.490         0.094         1           3-Nitroaniline         ND         ug/l         0.490         0.109         1           Dibenzofuran         2.92         ug/l         0.490         0.089         1	1,2,4,5-Tetrachlorobenzene	ND		ug/l	0.490	0.078	1
2-Chloronaphthalene ND ug/l 0.490 0.088 1 2-Nitroaniline ND ug/l 0.490 0.135 1 Acenaphthylene 2.70 ug/l 0.490 0.110 1 Dimethylphthalate ND ug/l 0.490 0.115 1 2,6-Dinitrotoluene ND ug/l 0.490 0.165 1 Acenaphthene 1.54 ug/l 0.490 0.094 1 3-Nitroaniline ND ug/l 0.490 0.109 1 Dibenzofuran 2.92 ug/l 0.490 0.089 1	Hexachlorocyclopentadiene	ND		ug/l	0.490	0.150	1
2-Nitroaniline ND ug/l 0.490 0.135 1 Acenaphthylene 2.70 ug/l 0.490 0.110 1 Dimethylphthalate ND ug/l 0.490 0.115 1 2,6-Dinitrotoluene ND ug/l 0.490 0.165 1 Acenaphthene 1.54 ug/l 0.490 0.094 1 3-Nitroaniline ND ug/l 0.490 0.109 1 Dibenzofuran 2.92 ug/l 0.490 0.089 1	Biphenyl	0.715		ug/l	0.490	0.109	1
Acenaphthylene         2.70         ug/l         0.490         0.110         1           Dimethylphthalate         ND         ug/l         0.490         0.115         1           2,6-Dinitrotoluene         ND         ug/l         0.490         0.165         1           Acenaphthene         1.54         ug/l         0.490         0.094         1           3-Nitroaniline         ND         ug/l         0.490         0.109         1           Dibenzofuran         2.92         ug/l         0.490         0.089         1	2-Chloronaphthalene	ND		ug/l	0.490	0.088	1
Dimethylphthalate         ND         ug/l         0.490         0.115         1           2,6-Dinitrotoluene         ND         ug/l         0.490         0.165         1           Acenaphthene         1.54         ug/l         0.490         0.094         1           3-Nitroaniline         ND         ug/l         0.490         0.109         1           Dibenzofuran         2.92         ug/l         0.490         0.089         1	2-Nitroaniline	ND		ug/l	0.490	0.135	1
2,6-Dinitrotoluene ND ug/l 0.490 0.165 1  Acenaphthene 1.54 ug/l 0.490 0.094 1  3-Nitroaniline ND ug/l 0.490 0.109 1  Dibenzofuran 2.92 ug/l 0.490 0.089 1	Acenaphthylene	2.70		ug/l	0.490	0.110	1
Acenaphthene 1.54 ug/l 0.490 0.094 1 3-Nitroaniline ND ug/l 0.490 0.109 1 Dibenzofuran 2.92 ug/l 0.490 0.089 1	Dimethylphthalate	ND		ug/l	0.490	0.115	1
3-Nitroaniline ND ug/l 0.490 0.109 1 Dibenzofuran 2.92 ug/l 0.490 0.089 1	2,6-Dinitrotoluene	ND		ug/l	0.490	0.165	1
Dibenzofuran 2.92 ug/l 0.490 0.089 1	Acenaphthene	1.54		ug/l	0.490	0.094	1
	3-Nitroaniline	ND		ug/l	0.490	0.109	1
2,4-Dinitrotoluene ND ug/l 0.490 0.160 1	Dibenzofuran	2.92		ug/l	0.490	0.089	1
	2,4-Dinitrotoluene	ND		ug/l	0.490	0.160	1



10/13/22

**Project Name:** Lab Number: STEELWINDS ANNUAL/SEMI-ANNUAL L2250575

**Project Number:** 03.0033579.15

L2250575-03

**SAMPLE RESULTS** 

Date Collected: 09/15/22 10:30

Report Date:

Date Received: 09/15/22 MWN-03-091522 Field Prep: Sample Location: LACKAWANNA NY Not Specified

Sample Depth:

Lab ID:

Client ID:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS	- Mansfield Lab					
Fluorene	5.10		ug/l	0.490	0.102	1
Diethylphthalate	ND		ug/l	0.490	0.176	1
4-Nitroaniline	ND		ug/l	0.490	0.110	1
n-Nitrosodiphenylamine	ND		ug/l	0.490	0.071	1
Hexachlorobenzene	ND		ug/l	0.490	0.120	1
Phenanthrene	9.37		ug/l	0.490	0.109	1
Anthracene	1.38		ug/l	0.490	0.134	1
Carbazole	5.17		ug/l	0.490	0.140	1
Di-n-butylphthalate	ND		ug/l	0.490	0.098	1
Fluoranthene	3.19		ug/l	0.490	0.153	1
Pyrene	1.91		ug/l	0.490	0.167	1
Butylbenzylphthalate	ND		ug/l	0.490	0.083	1
3,3'-Dichlorobenzidine	ND		ug/l	0.490	0.189	1
Benz(a)anthracene	ND		ug/l	0.490	0.180	1
Chrysene	ND		ug/l	0.490	0.139	1
bis(2-Ethylhexyl)phthalate	ND		ug/l	0.490	0.079	1
Di-n-octylphthalate	ND		ug/l	0.980	0.077	1
Benzo(b)fluoranthene	ND		ug/l	0.490	0.064	1
Benzo(k)fluoranthene	ND		ug/l	0.490	0.158	1
Benzo(a)pyrene	ND		ug/l	0.490	0.059	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.490	0.088	1
Dibenz(a,h)anthracene	ND		ug/l	0.490	0.063	1
Benzo(g,h,i)perylene	ND		ug/l	0.490	0.107	1

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	34	15-115
Phenol-d5	22	15-115
Nitrobenzene-d5	67	30-130
2-Fluorobiphenyl	66	30-130
2,4,6-Tribromophenol	98	15-115
Terphenyl-d14	89	30-130



L2250575

10/13/22

09/17/22 08:30

**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

**Project Number:** 03.0033579.15

**SAMPLE RESULTS** 

Date Collected: 09/15/22 14:50

Lab Number:

Report Date:

L2250575-05 Date Received: 09/15/22 MWN-04-091522 Field Prep: LACKAWANNA NY Not Specified

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Extraction Method: EPA 3510C Matrix: Water **Extraction Date:** Analytical Method: 1,8270D Analytical Date: 09/19/22 17:08

Analyst: PS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor				
Semivolatile Organics by GC/MS - Mansfield Lab										
bis(2-Chloroethyl)ether	ND		ug/l	0.495	0.092	1				
1,3-Dichlorobenzene	ND		ug/l	0.495	0.078	1				
1,4-Dichlorobenzene	ND		ug/l	0.495	0.082	1				
1,2-Dichlorobenzene	ND		ug/l	0.495	0.067	1				
Benzyl alcohol	ND		ug/l	0.495	0.122	1				
bis(2-chloroisopropyl)ether	ND		ug/l	0.495	0.107	1				
Acetophenone	0.967	J	ug/l	0.990	0.205	1				
Hexachloroethane	ND		ug/l	0.495	0.101	1				
Nitrobenzene	ND		ug/l	0.495	0.101	1				
Isophorone	ND		ug/l	0.495	0.125	1				
bis(2-Chloroethoxy)methane	ND		ug/l	0.495	0.085	1				
1,2,4-Trichlorobenzene	ND		ug/l	0.495	0.095	1				
Naphthalene	11.2		ug/l	0.495	0.087	1				
4-Chloroaniline	ND		ug/l	0.495	0.127	1				
Hexachlorobutadiene	ND		ug/l	0.495	0.085	1				
2-Methylnaphthalene	2.49		ug/l	0.495	0.090	1				
1,2,4,5-Tetrachlorobenzene	ND		ug/l	0.495	0.079	1				
Hexachlorocyclopentadiene	ND		ug/l	0.495	0.151	1				
Biphenyl	0.394	J	ug/l	0.495	0.110	1				
2-Chloronaphthalene	ND		ug/l	0.495	0.089	1				
2-Nitroaniline	ND		ug/l	0.495	0.137	1				
Acenaphthylene	0.167	J	ug/l	0.495	0.111	1				
Dimethylphthalate	ND		ug/l	0.495	0.116	1				
2,6-Dinitrotoluene	ND		ug/l	0.495	0.166	1				
Acenaphthene	5.26		ug/l	0.495	0.095	1				
3-Nitroaniline	ND		ug/l	0.495	0.110	1				
Dibenzofuran	2.54		ug/l	0.495	0.090	1				
2,4-Dinitrotoluene	ND		ug/l	0.495	0.161	1				



10/13/22

Project Name: STEELWINDS ANNUAL/SEMI-ANNUAL Lab Number: L2250575

**Project Number:** 03.0033579.15

L2250575-05

**SAMPLE RESULTS** 

Date Collected: 09/15/22 14:50

Report Date:

Client ID: MWN-04-091522 Date Received: 09/15/22 Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Lab ID:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor				
Semivolatile Organics by GC/MS - Mansfield Lab										
Fluorene	4.37		ug/l	0.495	0.103	1				
Diethylphthalate	ND		ug/l	0.495	0.178	1				
4-Nitroaniline	ND		ug/l	0.495	0.111	1				
n-Nitrosodiphenylamine	ND		ug/l	0.495	0.071	1				
Hexachlorobenzene	ND		ug/l	0.495	0.121	1				
Phenanthrene	7.31		ug/l	0.495	0.110	1				
Anthracene	1.39		ug/l	0.495	0.136	1				
Carbazole	8.59		ug/l	0.495	0.142	1				
Di-n-butylphthalate	ND		ug/l	0.495	0.099	1				
Fluoranthene	1.55		ug/l	0.495	0.154	1				
Pyrene	1.90		ug/l	0.495	0.168	1				
Butylbenzylphthalate	ND		ug/l	0.495	0.084	1				
3,3'-Dichlorobenzidine	ND		ug/l	0.495	0.191	1				
Benz(a)anthracene	ND		ug/l	0.495	0.182	1				
Chrysene	ND		ug/l	0.495	0.140	1				
bis(2-Ethylhexyl)phthalate	ND		ug/l	0.495	0.080	1				
Di-n-octylphthalate	ND		ug/l	0.990	0.078	1				
Benzo(b)fluoranthene	0.125	J	ug/l	0.495	0.065	1				
Benzo(k)fluoranthene	ND		ug/l	0.495	0.159	1				
Benzo(a)pyrene	0.076	J	ug/l	0.495	0.060	1				
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.495	0.089	1				
Dibenz(a,h)anthracene	ND		ug/l	0.495	0.064	1				
Benzo(g,h,i)perylene	ND		ug/l	0.495	0.108	1				

Surrogate	% Recovery	Acceptance Qualifier Criteria
2-Fluorophenol	39	15-115
Phenol-d5	25	15-115
Nitrobenzene-d5	70	30-130
2-Fluorobiphenyl	69	30-130
2,4,6-Tribromophenol	91	15-115
Terphenyl-d14	84	30-130



L2250575

Project Name: STEELWINDS ANNUAL/SEMI-ANNUAL Lab Number:

**Project Number:** 03.0033579.15 **Report Date:** 10/13/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270D Extraction Method: EPA 3510C
Analytical Date: 09/19/22 10:43 Extraction Date: 09/17/22 08:30

Analyst: PS

bis(2-Chloroethyl)ether   ND	Parameter	Result	Qualifier	Units	RL		MDL
1,3-Dichlorobenzene         ND         ug/l         0.500         0.078           1,4-Dichlorobenzene         ND         ug/l         0.500         0.083           1,2-Dichlorobenzene         ND         ug/l         0.500         0.068           Benzyl alcohol         ND         ug/l         0.500         0.123           bis(2-chloroisopropyl)ether         ND         ug/l         0.500         0.108           Acetophenone         ND         ug/l         1.00         0.207           Hexachloroethane         ND         ug/l         0.500         0.102           Nitrobenzene         ND         ug/l         0.500         0.102           Isophorone         ND         ug/l         0.500         0.126           bis(2-Chloroethoxy)methane         ND         ug/l         0.500         0.085           1,2,4-Trichlorobenzene         ND         ug/l         0.500         0.085           1,2,4-Trichlorobenzene         ND         ug/l         0.500         0.088           4-Chloroaniline         ND         ug/l         0.500         0.088           4-Chloroaniline         ND         ug/l         0.500         0.086           Hexachlorobutadiene	Semivolatile Organics by GC/MS	- Mansfield La	ab for sample	e(s):	01,03,05	Batch:	WG1688692-1
1,4-Dichlorobenzene         ND         ug/l         0.500         0.083           1,2-Dichlorobenzene         ND         ug/l         0.500         0.068           Benzyl alcohol         ND         ug/l         0.500         0.123           bis(2-chloroisopropyl)ether         ND         ug/l         0.500         0.108           Acetophenone         ND         ug/l         1.00         0.207           Hexachloroethane         ND         ug/l         0.500         0.102           Nitrobenzene         ND         ug/l         0.500         0.102           Isophorone         ND         ug/l         0.500         0.102           Isophorone         ND         ug/l         0.500         0.126           Isophorone         ND         ug/l         0.500         0.085           Ley-Trichlorobenzene         ND         ug/l         0.500         0.088           4-Chloroaphthalene         ND         ug/l         <	bis(2-Chloroethyl)ether	ND		ug/l	0.500	)	0.093
1,2-Dichlorobenzene         ND         ug/l         0.500         0.068           Benzyl alcohol         ND         ug/l         0.500         0.123           bis(2-chloroisopropyl)ether         ND         ug/l         0.500         0.108           Acetophenone         ND         ug/l         1.00         0.207           Hexachloroethane         ND         ug/l         0.500         0.102           Nitrobenzene         ND         ug/l         0.500         0.102           Isophorone         ND         ug/l         0.500         0.126           bis(2-Chloroethoxy)methane         ND         ug/l         0.500         0.085           1,2,4-Trichlorobenzene         ND         ug/l         0.500         0.085           1,2,4-Trichlorobenzene         ND         ug/l         0.500         0.086           Naphthalene         ND         ug/l         0.500         0.088           4-Chloroaniline         ND         ug/l         0.500         0.088           Hexachlorobutadiene         ND         ug/l         0.500         0.086           2-Methylnaphthalene         ND         ug/l         0.500         0.091           1,2,4,5-Tetrachlorobenzene <td>1,3-Dichlorobenzene</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.500</td> <td>)</td> <td>0.078</td>	1,3-Dichlorobenzene	ND		ug/l	0.500	)	0.078
Benzyl alcohol   ND   ug/l   0.500   0.123	1,4-Dichlorobenzene	ND		ug/l	0.500	)	0.083
bis(2-chloroisopropyl)ether         ND         ug/l         0.500         0.108           Acetophenone         ND         ug/l         1.00         0.207           Hexachloroethane         ND         ug/l         0.500         0.102           Nitrobenzene         ND         ug/l         0.500         0.102           Isophorone         ND         ug/l         0.500         0.126           bis(2-Chloroethoxy)methane         ND         ug/l         0.500         0.085           1,2,4-Trichlorobenzene         ND         ug/l         0.500         0.096           Naphthalene         ND         ug/l         0.500         0.088           4-Chloroaniline         ND         ug/l         0.500         0.088           4-Chloroaniline         ND         ug/l         0.500         0.086           2-Methylnaphthalene         ND         ug/l         0.500         0.091           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.500         0.080           Hexachlorocyclopentadiene         ND         ug/l         0.500         0.153           Biphenyl         ND         ug/l         0.500         0.111           2-Chloronaphthalene	1,2-Dichlorobenzene	ND		ug/l	0.500	)	0.068
Acetophenone         ND         ug/l         1.00         0.207           Hexachloroethane         ND         ug/l         0.500         0.102           Nitrobenzene         ND         ug/l         0.500         0.102           Isophorone         ND         ug/l         0.500         0.126           bis(2-Chloroethoxy)methane         ND         ug/l         0.500         0.085           1,2,4-Trichlorobenzene         ND         ug/l         0.500         0.096           Naphthalene         ND         ug/l         0.500         0.098           4-Chloroaniline         ND         ug/l         0.500         0.088           4-Chloroaniline         ND         ug/l         0.500         0.086           2-Methylnaphthalene         ND         ug/l         0.500         0.091           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.500         0.091           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.500         0.080           Hexachlorocyclopentadiene         ND         ug/l         0.500         0.113           Biphenyl         ND         ug/l         0.500         0.111           2-Chloronaphthalene	Benzyl alcohol	ND		ug/l	0.500	)	0.123
Hexachloroethane   ND	bis(2-chloroisopropyl)ether	ND		ug/l	0.500	)	0.108
Nitrobenzene         ND         ug/l         0.500         0.102           Isophorone         ND         ug/l         0.500         0.126           bis(2-Chloroethoxy)methane         ND         ug/l         0.500         0.085           1,2,4-Trichlorobenzene         ND         ug/l         0.500         0.096           Naphthalene         ND         ug/l         0.500         0.088           4-Chloroaniline         ND         ug/l         0.500         0.128           Hexachlorobutadiene         ND         ug/l         0.500         0.086           2-Methylnaphthalene         ND         ug/l         0.500         0.091           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.500         0.080           Hexachlorocyclopentadiene         ND         ug/l         0.500         0.153           Biphenyl         ND         ug/l         0.500         0.111           2-Chloronaphthalene         ND         ug/l         0.500         0.138           Acenaphthylene         ND         ug/l         0.500         0.112           Dimethylphthalate         ND         ug/l         0.500         0.168           Acenaphthene	Acetophenone	ND		ug/l	1.00		0.207
Sophorone   ND	Hexachloroethane	ND		ug/l	0.500	)	0.102
bis(2-Chloroethoxy)methane         ND         ug/l         0.500         0.085           1,2,4-Trichlorobenzene         ND         ug/l         0.500         0.096           Naphthalene         ND         ug/l         0.500         0.088           4-Chloroaniline         ND         ug/l         0.500         0.128           Hexachlorobutadiene         ND         ug/l         0.500         0.086           2-Methylnaphthalene         ND         ug/l         0.500         0.091           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.500         0.080           Hexachlorocyclopentadiene         ND         ug/l         0.500         0.153           Biphenyl         ND         ug/l         0.500         0.111           2-Chloronaphthalene         ND         ug/l         0.500         0.111           2-Chloronaphthylene         ND         ug/l         0.500         0.138           Acenaphthylene         ND         ug/l         0.500         0.112           Dimethylphthalate         ND         ug/l         0.500         0.117           2,6-Dinitrotoluene         ND         ug/l         0.500         0.168           Acenaphthene <td>Nitrobenzene</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.500</td> <td>)</td> <td>0.102</td>	Nitrobenzene	ND		ug/l	0.500	)	0.102
1,2,4-Trichlorobenzene         ND         ug/l         0.500         0.096           Naphthalene         ND         ug/l         0.500         0.088           4-Chloroaniline         ND         ug/l         0.500         0.128           Hexachlorobutadiene         ND         ug/l         0.500         0.086           2-Methylnaphthalene         ND         ug/l         0.500         0.091           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.500         0.080           Hexachlorocyclopentadiene         ND         ug/l         0.500         0.153           Biphenyl         ND         ug/l         0.500         0.111           2-Chloronaphthalene         ND         ug/l         0.500         0.111           2-Chloronaphthylene         ND         ug/l         0.500         0.118           Acenaphthylene         ND         ug/l         0.500         0.1112           Dimethylphthalate         ND         ug/l         0.500         0.1168           Acenaphthene         ND         ug/l         0.500         0.168           Acenaphthene         ND         ug/l         0.500         0.111           Dibenzofuran         ND<	Isophorone	ND		ug/l	0.500	)	0.126
Naphthalene         ND         ug/l         0.500         0.088           4-Chloroaniline         ND         ug/l         0.500         0.128           Hexachlorobutadiene         ND         ug/l         0.500         0.086           2-Methylnaphthalene         ND         ug/l         0.500         0.091           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.500         0.080           Hexachlorocyclopentadiene         ND         ug/l         0.500         0.153           Biphenyl         ND         ug/l         0.500         0.111           2-Chloronaphthalene         ND         ug/l         0.500         0.111           2-Nitroaniline         ND         ug/l         0.500         0.138           Acenaphthylene         ND         ug/l         0.500         0.112           Dimethylphthalate         ND         ug/l         0.500         0.117           2,6-Dinitrotoluene         ND         ug/l         0.500         0.168           Acenaphthene         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND	bis(2-Chloroethoxy)methane	ND		ug/l	0.500	)	0.085
4-Chloroaniline         ND         ug/l         0.500         0.128           Hexachlorobutadiene         ND         ug/l         0.500         0.086           2-Methylnaphthalene         ND         ug/l         0.500         0.091           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.500         0.080           Hexachlorocyclopentadiene         ND         ug/l         0.500         0.153           Biphenyl         ND         ug/l         0.500         0.111           2-Chloronaphthalene         ND         ug/l         0.500         0.090           2-Nitroaniline         ND         ug/l         0.500         0.1138           Acenaphthylene         ND         ug/l         0.500         0.112           Dimethylphthalate         ND         ug/l         0.500         0.117           2,6-Dinitrotoluene         ND         ug/l         0.500         0.168           Acenaphthene         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND <td>1,2,4-Trichlorobenzene</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.500</td> <td>)</td> <td>0.096</td>	1,2,4-Trichlorobenzene	ND		ug/l	0.500	)	0.096
Hexachlorobutadiene         ND         ug/l         0.500         0.086           2-Methylnaphthalene         ND         ug/l         0.500         0.091           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.500         0.080           Hexachlorocyclopentadiene         ND         ug/l         0.500         0.153           Biphenyl         ND         ug/l         0.500         0.111           2-Chloronaphthalene         ND         ug/l         0.500         0.090           2-Nitroaniline         ND         ug/l         0.500         0.138           Acenaphthylene         ND         ug/l         0.500         0.112           Dimethylphthalate         ND         ug/l         0.500         0.117           2,6-Dinitrotoluene         ND         ug/l         0.500         0.168           Acenaphthene         ND         ug/l         0.500         0.111           3-Nitroaniline         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND         ug/l         0.500         0.163	Naphthalene	ND		ug/l	0.500	)	0.088
2-Methylnaphthalene         ND         ug/l         0.500         0.091           1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.500         0.080           Hexachlorocyclopentadiene         ND         ug/l         0.500         0.153           Biphenyl         ND         ug/l         0.500         0.111           2-Chloronaphthalene         ND         ug/l         0.500         0.090           2-Nitroaniline         ND         ug/l         0.500         0.138           Acenaphthylene         ND         ug/l         0.500         0.112           Dimethylphthalate         ND         ug/l         0.500         0.117           2,6-Dinitrotoluene         ND         ug/l         0.500         0.168           Acenaphthene         ND         ug/l         0.500         0.096           3-Nitroaniline         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND         ug/l         0.500         0.091	4-Chloroaniline	ND		ug/l	0.500	)	0.128
1,2,4,5-Tetrachlorobenzene         ND         ug/l         0.500         0.080           Hexachlorocyclopentadiene         ND         ug/l         0.500         0.153           Biphenyl         ND         ug/l         0.500         0.111           2-Chloronaphthalene         ND         ug/l         0.500         0.090           2-Nitroaniline         ND         ug/l         0.500         0.138           Acenaphthylene         ND         ug/l         0.500         0.112           Dimethylphthalate         ND         ug/l         0.500         0.117           2,6-Dinitrotoluene         ND         ug/l         0.500         0.168           Acenaphthene         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND         ug/l         0.500         0.163	Hexachlorobutadiene	ND		ug/l	0.500	)	0.086
Hexachlorocyclopentadiene         ND         ug/l         0.500         0.153           Biphenyl         ND         ug/l         0.500         0.111           2-Chloronaphthalene         ND         ug/l         0.500         0.090           2-Nitroaniline         ND         ug/l         0.500         0.138           Acenaphthylene         ND         ug/l         0.500         0.112           Dimethylphthalate         ND         ug/l         0.500         0.117           2,6-Dinitrotoluene         ND         ug/l         0.500         0.168           Acenaphthene         ND         ug/l         0.500         0.096           3-Nitroaniline         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND         ug/l         0.500         0.163	2-Methylnaphthalene	ND		ug/l	0.500	)	0.091
Biphenyl         ND         ug/l         0.500         0.111           2-Chloronaphthalene         ND         ug/l         0.500         0.090           2-Nitroaniline         ND         ug/l         0.500         0.138           Acenaphthylene         ND         ug/l         0.500         0.112           Dimethylphthalate         ND         ug/l         0.500         0.117           2,6-Dinitrotoluene         ND         ug/l         0.500         0.168           Acenaphthene         ND         ug/l         0.500         0.096           3-Nitroaniline         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND         ug/l         0.500         0.163	1,2,4,5-Tetrachlorobenzene	ND		ug/l	0.500	)	0.080
2-Chloronaphthalene         ND         ug/l         0.500         0.090           2-Nitroaniline         ND         ug/l         0.500         0.138           Acenaphthylene         ND         ug/l         0.500         0.112           Dimethylphthalate         ND         ug/l         0.500         0.117           2,6-Dinitrotoluene         ND         ug/l         0.500         0.168           Acenaphthene         ND         ug/l         0.500         0.096           3-Nitroaniline         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND         ug/l         0.500         0.163	Hexachlorocyclopentadiene	ND		ug/l	0.500	)	0.153
2-Nitroaniline         ND         ug/l         0.500         0.138           Acenaphthylene         ND         ug/l         0.500         0.112           Dimethylphthalate         ND         ug/l         0.500         0.117           2,6-Dinitrotoluene         ND         ug/l         0.500         0.168           Acenaphthene         ND         ug/l         0.500         0.096           3-Nitroaniline         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND         ug/l         0.500         0.163	Biphenyl	ND		ug/l	0.500	)	0.111
Acenaphthylene         ND         ug/l         0.500         0.112           Dimethylphthalate         ND         ug/l         0.500         0.117           2,6-Dinitrotoluene         ND         ug/l         0.500         0.168           Acenaphthene         ND         ug/l         0.500         0.096           3-Nitroaniline         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND         ug/l         0.500         0.163	2-Chloronaphthalene	ND		ug/l	0.500	)	0.090
Dimethylphthalate         ND         ug/l         0.500         0.117           2,6-Dinitrotoluene         ND         ug/l         0.500         0.168           Acenaphthene         ND         ug/l         0.500         0.096           3-Nitroaniline         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND         ug/l         0.500         0.163	2-Nitroaniline	ND		ug/l	0.500	)	0.138
2,6-Dinitrotoluene       ND       ug/l       0.500       0.168         Acenaphthene       ND       ug/l       0.500       0.096         3-Nitroaniline       ND       ug/l       0.500       0.111         Dibenzofuran       ND       ug/l       0.500       0.091         2,4-Dinitrotoluene       ND       ug/l       0.500       0.163	Acenaphthylene	ND		ug/l	0.500	)	0.112
Acenaphthene         ND         ug/l         0.500         0.096           3-Nitroaniline         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND         ug/l         0.500         0.163	Dimethylphthalate	ND		ug/l	0.500	)	0.117
3-Nitroaniline         ND         ug/l         0.500         0.111           Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND         ug/l         0.500         0.163	2,6-Dinitrotoluene	ND		ug/l	0.500	)	0.168
Dibenzofuran         ND         ug/l         0.500         0.091           2,4-Dinitrotoluene         ND         ug/l         0.500         0.163	Acenaphthene	ND		ug/l	0.500	)	0.096
2,4-Dinitrotoluene ND ug/l 0.500 0.163	3-Nitroaniline	ND		ug/l	0.500	)	0.111
·	Dibenzofuran	ND		ug/l	0.500	)	0.091
Fluorene ND ug/l 0.500 0.104	2,4-Dinitrotoluene	ND		ug/l	0.500	)	0.163
	Fluorene	ND		ug/l	0.500	)	0.104



L2250575

Project Name: STEELWINDS ANNUAL/SEMI-ANNUAL Lab Number:

**Project Number:** 03.0033579.15 **Report Date:** 10/13/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270D Extraction Method: EPA 3510C
Analytical Date: 09/19/22 10:43 Extraction Date: 09/17/22 08:30

Analyst: PS

Parameter	Result	Qualifier	Units	RL		MDL
Semivolatile Organics by GC/MS - M	/lansfield La	b for samp	le(s):	01,03,05	Batch:	WG1688692-1
Diethylphthalate	ND		ug/l	0.500	)	0.180
4-Nitroaniline	ND		ug/l	0.500	)	0.112
n-Nitrosodiphenylamine	ND		ug/l	0.500	)	0.072
Hexachlorobenzene	ND		ug/l	0.500	)	0.122
Phenanthrene	ND		ug/l	0.500	)	0.111
Anthracene	ND		ug/l	0.500	)	0.137
Carbazole	ND		ug/l	0.500	)	0.143
Di-n-butylphthalate	ND		ug/l	0.500	)	0.100
Fluoranthene	ND		ug/l	0.500	)	0.156
Pyrene	ND		ug/l	0.500	)	0.170
Butylbenzylphthalate	ND		ug/l	0.500	)	0.085
3,3'-Dichlorobenzidine	ND		ug/l	0.500	)	0.193
Benz(a)anthracene	ND		ug/l	0.500	)	0.184
Chrysene	ND		ug/l	0.500	)	0.142
bis(2-Ethylhexyl)phthalate	ND		ug/l	0.500	)	0.081
Di-n-octylphthalate	ND		ug/l	1.00		0.079
Benzo(b)fluoranthene	ND		ug/l	0.500	)	0.066
Benzo(k)fluoranthene	ND		ug/l	0.500	)	0.161
Benzo(a)pyrene	ND		ug/l	0.500	)	0.060
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.500	)	0.090
Dibenz(a,h)anthracene	ND		ug/l	0.500	)	0.064
Benzo(g,h,i)perylene	ND		ug/l	0.500	)	0.109



Project Name: STEELWINDS ANNUAL/SEMI-ANNUAL Lab Number: L2250575

**Project Number:** 03.0033579.15 **Report Date:** 10/13/22

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8270D Extraction Method: EPA 3510C
Analytical Date: 09/19/22 10:43 Extraction Date: 09/17/22 08:30

Analyst: PS

Parameter Result Qualifier Units RL MDL

Semivolatile Organics by GC/MS - Mansfield Lab for sample(s): 01,03,05 Batch: WG1688692-1

Surrogate	%Recovery Qualifie	Acceptance r Criteria
2-Fluorophenol	64	15-115
Phenol-d5	42	15-115
Nitrobenzene-d5	93	30-130
2-Fluorobiphenyl	90	30-130
2,4,6-Tribromophenol	95	15-115
Terphenyl-d14	106	30-130



**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

**Project Number:** 03.0033579.15

Lab Number: L2250575

**Report Date:** 10/13/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	%Recovery Qual Limits	RPD	RPD Qual Limits
Semivolatile Organics by GC/MS - Mansfield	Lab Associate	d sample(s):	01,03,05 Batch:	WG1688692-2 WG168869	02-3	
bis(2-Chloroethyl)ether	84		88	40-140	5	20
1,3-Dichlorobenzene	44		53	40-140	19	20
1,4-Dichlorobenzene	46		54	40-140	16	20
1,2-Dichlorobenzene	48		56	40-140	15	20
bis(2-chloroisopropyl)ether	75		80	40-140	6	20
Acetophenone	82		85	40-140	4	20
Hexachloroethane	40		49	10-97	20	20
Nitrobenzene	83		88	40-140	6	20
Isophorone	82		86	40-140	5	20
bis(2-Chloroethoxy)methane	89		93	40-140	4	20
1,2,4-Trichlorobenzene	50		57	40-140	13	20
Naphthalene	65		71	40-140	9	20
4-Chloroaniline	93		97	40-140	4	20
Hexachlorobutadiene	41		47	40-140	14	20
2-Methylnaphthalene	64		69	40-140	8	20
1,2,4,5-Tetrachlorobenzene	56		61	40-140	9	20
Hexachlorocyclopentadiene	40		48	10-109	18	20
Biphenyl	75		74	40-140	1	20
2-Chloronaphthalene	68		73	40-140	7	20
2-Nitroaniline	93		98	40-140	5	20
Acenaphthylene	77		81	40-140	5	20
Dimethylphthalate	84		80	40-140	5	20
2,6-Dinitrotoluene	101		106	40-140	5	20



**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

**Project Number:** 03.0033579.15

Lab Number: L2250575

**Report Date:** 10/13/22

Parameter	LCS %Recovery	Qual	LCSD %Recove	ry	% Qual	Recovery Limits	RPD	Qual	RPD Limits
Semivolatile Organics by GC/MS - Mansfield	Lab Associated	sample(s):	01,03,05 B	atch:	WG1688692-	2 WG168869	2-3		
Acenaphthene	76		80			40-140	5		20
3-Nitroaniline	100		106			40-140	6		20
Dibenzofuran	84		87			40-140	4		20
2,4-Dinitrotoluene	106		110			40-140	4		20
Fluorene	86		89			40-140	3		20
Diethylphthalate	92		92			40-140	0		20
4-Nitroaniline	108		113			40-140	5		20
n-Nitrosodiphenylamine	99		103			40-140	4		20
Hexachlorobenzene	95		98			40-140	3		20
Phenanthrene	93		96			40-140	3		20
Anthracene	96		100			40-140	4		20
Carbazole	100		106			40-140	6		20
Di-n-butylphthalate	101		104			40-140	3		20
Fluoranthene	98		103			40-140	5		20
Pyrene	94		99			40-140	5		20
Butylbenzylphthalate	97		101			40-140	4		20
3,3'-Dichlorobenzidine	92		99			40-140	7		20
Benz(a)anthracene	91		97			40-140	6		20
Chrysene	99		105			40-140	6		20
bis(2-Ethylhexyl)phthalate	99		105			40-140	6		20
Di-n-octylphthalate	103		108			40-140	5		20
Benzo(b)fluoranthene	89		93			40-140	4		20
Benzo(k)fluoranthene	105		111			40-140	6		20



**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

Lab Number:

L2250575

**Project Number:** 03.0033579.15

Report Date:

10/13/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Semivolatile Organics by GC/MS - Mansfield	Lab Associated	d sample(s): (	01,03,05 Batch	: WG16886	92-2 WG168869	2-3		
Benzo(a)pyrene	97		101		40-140	4		20
Indeno(1,2,3-cd)pyrene	102		109		40-140	7		20
Dibenz(a,h)anthracene	101		105		40-140	4		20
Benzo(g,h,i)perylene	91		95		40-140	4		20

	LCS	LCSD	Acceptance
Surrogate	%Recovery Qu	ual %Recovery Qua	al Criteria
2-Fluorophenol	58	61	15-115
Phenol-d5	42	43	15-115
Nitrobenzene-d5	86	89	30-130
2-Fluorobiphenyl	80	81	30-130
2,4,6-Tribromophenol	99	102	15-115
Terphenyl-d14	95	100	30-130

### **METALS**



Project Name:STEELWINDS ANNUAL/SEMI-ANNUALLab Number:L2250575

**Project Number:** 03.0033579.15 **Report Date:** 10/13/22

**SAMPLE RESULTS** 

 Lab ID:
 L2250575-01
 Date Collected:
 09/15/22 08:30

 Client ID:
 MWN-03D-091522
 Date Received:
 09/15/22

Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Matrix: Water

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	Date Prepared	Date Analyzed	Prep Method	Analytical Method	Analyst
Dissolved Metals - N	Mansfield	Lab									
Barium, Dissolved	0.779		mg/l	0.0100	0.0021	1	09/21/22 23:00	0 10/13/22 07:28	EPA 3005A	1,6010D	NB
Manganese, Dissolved	0.333		mg/l	0.0100	0.0016	1	09/21/22 23:00	0 10/13/22 07:28	EPA 3005A	1,6010D	NB



Project Name:STEELWINDS ANNUAL/SEMI-ANNUALLab Number:L2250575

**Project Number:** 03.0033579.15 **Report Date:** 10/13/22

**SAMPLE RESULTS** 

 Lab ID:
 L2250575-02
 Date Collected:
 09/15/22 09:45

 Client ID:
 MWN-03B-091522
 Date Received:
 09/15/22

Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Matrix: Water

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	Date Prepared	Date Analyzed	Prep Method	Analytical Method	Analyst
Dissolved Metals - N	Mansfield	Lab									
Arsenic, Dissolved	ND		mg/l	0.0050	0.0019	1	09/21/22 23:00	10/13/22 07:31	EPA 3005A	1,6010D	NB
Barium, Dissolved	1.32		mg/l	0.0100	0.0021	1	09/21/22 23:00	10/13/22 07:31	EPA 3005A	1,6010D	NB
Chromium, Dissolved	0.0032	J	mg/l	0.0100	0.0021	1	09/21/22 23:00	10/13/22 07:31	EPA 3005A	1,6010D	NB
Manganese, Dissolved	0.178		mg/l	0.0100	0.0016	1	09/21/22 23:00	10/13/22 07:31	EPA 3005A	1,6010D	NB



Project Name:STEELWINDS ANNUAL/SEMI-ANNUALLab Number:L2250575

**Project Number:** 03.0033579.15 **Report Date:** 10/13/22

**SAMPLE RESULTS** 

 Lab ID:
 L2250575-04
 Date Collected:
 09/15/22 13:45

 Client ID:
 MWN-02D-091522
 Date Received:
 09/15/22

Sample Location: LACKAWANNA NY Field Prep: Not Specified

Sample Depth:

Matrix: Water

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	Date Prepared	Date Analyzed	Prep Method	Analytical Method	Analyst
Dissolved Metals -	Mansfield	Lab									
Arsenic, Dissolved	ND		mg/l	0.0050	0.0019	1	09/21/22 23:00	10/13/22 07:35	EPA 3005A	1,6010D	NB
Barium, Dissolved	0.860		mg/l	0.0100	0.0021	1	09/21/22 23:00	10/13/22 07:35	EPA 3005A	1,6010D	NB
Chromium, Dissolved	ND		mg/l	0.0100	0.0021	1	09/21/22 23:00	10/13/22 07:35	EPA 3005A	1,6010D	NB



Project Name:STEELWINDS ANNUAL/SEMI-ANNUALLab Number:L2250575

**Project Number:** 03.0033579.15 **Report Date:** 10/13/22

# Method Blank Analysis Batch Quality Control

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	Date Prepared	Date Analyzed	Analytical Method	Analyst
Dissolved Metals - Mar	nsfield Lab	for sample	(s): 01-0	2,04 Ba	tch: W	G1690348-	1			
Arsenic, Dissolved	ND		mg/l	0.0050	0.0019	1	09/21/22 23:00	10/05/22 07:21	1,6010D	EW
Barium, Dissolved	ND		mg/l	0.0100	0.0021	1	09/21/22 23:00	10/05/22 07:2	1,6010D	EW
Chromium, Dissolved	ND		mg/l	0.0100	0.0021	1	09/21/22 23:00	10/05/22 07:2	1,6010D	EW
Manganese, Dissolved	ND		mg/l	0.0100	0.0016	1	09/21/22 23:00	10/05/22 07:21	1,6010D	EW

**Prep Information** 

Digestion Method: EPA 3005A



## Lab Control Sample Analysis Batch Quality Control

**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

Lab Number: L2250575

**Project Number:** 03.0033579.15 Report Date: 10/13/22

Parameter	LCS %Recovery	Qual %	LCSD Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Dissolved Metals - Mansfield Lab Associated	sample(s): 01-02,0	4 Batch: WG	1690348-2					
Arsenic, Dissolved	108		-		80-120	-		
Barium, Dissolved	100		-		80-120	-		
Chromium, Dissolved	95		-		80-120	-		
Manganese, Dissolved	98		-		80-120	-		

### Matrix Spike Analysis Batch Quality Control

Project Name: STEELWINDS ANNUAL/SEMI-ANNUAL

**Project Number:** 03.0033579.15

Lab Number: L2250575

**Report Date:** 10/13/22

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	MSD Qual Found	MSD %Recovery Q	Recovery ual Limits	RPD Qual	RPD Limits
Dissolved Metals - Mansfield	Lab Associated	d sample(s):	01-02,04	QC Batch ID:	WG1690348-3	QC Sample: L224	9555-13 Client	t ID: MS Sam	ple
Arsenic, Dissolved	ND	0.12	0.128	107	-	-	75-125	-	20
Barium, Dissolved	ND	2	1.92	96	-	-	75-125	-	20
Chromium, Dissolved	ND	0.2	0.182	91	-	-	75-125	-	20
Manganese, Dissolved	ND	0.5	0.453	91	-	-	75-125	-	20

## Lab Duplicate Analysis Batch Quality Control

**Project Name:** STEELWINDS ANNUAL/SEMI-ANNUAL

**Project Number:** 03.0033579.15 Lab Number:

L2250575

Report Date:

10/13/22

Parameter	Native Sa	mple	Duplicate Samp	ole Units	RPD	Qual	RPD Limits
Dissolved Metals - Mansfield Lab Associated sample(s):	01-02,04	QC Batch ID:	WG1690348-4	QC Sample:	L2249555-13	Client ID:	DUP Sample
Arsenic, Dissolved	ND		ND	mg/l	NC		20
Barium, Dissolved	ND		ND	mg/l	NC		20
Chromium, Dissolved	ND		ND	mg/l	NC		20
Manganese, Dissolved	ND		0.0048J	mg/l	NC		20



Serial\_No:10132213:50 *Lab Number:* L2250575

**Report Date:** 10/13/22

Project Name: STEELWINDS ANNUAL/SEMI-ANNUAL

Project Number: 03.0033579.15

### Sample Receipt and Container Information

Were project specific reporting limits specified?

**Cooler Information** 

Cooler Custody Seal

A Absent

Container Info	ormation		Initial	Final	Temp			Frozen	
Container ID	Container Type	Cooler	рН	pН	deg C	Pres	Seal	Date/Time	Analysis(*)
L2250575-01A	Vial HCI preserved	Α	NA		4.3	Υ	Absent		NYCP51-8260(14)
L2250575-01B	Vial HCl preserved	Α	NA		4.3	Υ	Absent		NYCP51-8260(14)
L2250575-01C	Vial HCl preserved	Α	NA		4.3	Υ	Absent		NYCP51-8260(14)
L2250575-01D	Plastic 250ml unpreserved	Α	<2	<2	4.3	Υ	Absent		-
L2250575-01E	Amber 1000ml unpreserved	Α	7	7	4.3	Υ	Absent		A2-SVOC-8270(7)
L2250575-01F	Amber 1000ml unpreserved	Α	7	7	4.3	Υ	Absent		A2-SVOC-8270(7)
L2250575-01X	Plastic 120ml HNO3 preserved Filtrates	Α	NA		4.3	Υ	Absent		BA-SI(180),MN-SI(180)
L2250575-02A	Plastic 250ml unpreserved	Α	<2	<2	4.3	Υ	Absent		-
L2250575-02X	Plastic 120ml HNO3 preserved Filtrates	Α	NA		4.3	Υ	Absent		BA-SI(180),MN-SI(180),AS-SI(180),CR-SI(180)
L2250575-03A	Vial HCl preserved	Α	NA		4.3	Υ	Absent		NYCP51-8260(14)
L2250575-03B	Vial HCI preserved	Α	NA		4.3	Υ	Absent		NYCP51-8260(14)
L2250575-03C	Vial HCI preserved	Α	NA		4.3	Υ	Absent		NYCP51-8260(14)
L2250575-03D	Amber 1000ml unpreserved	Α	7	7	4.3	Υ	Absent		A2-SVOC-8270(7)
L2250575-03E	Amber 1000ml unpreserved	Α	7	7	4.3	Υ	Absent		A2-SVOC-8270(7)
L2250575-04A	Plastic 250ml unpreserved	Α	<2	<2	4.3	Υ	Absent		-
L2250575-04X	Plastic 120ml HNO3 preserved Filtrates	Α	NA		4.3	Υ	Absent		BA-SI(180),AS-SI(180),CR-SI(180)
L2250575-05A	Vial HCI preserved	Α	NA		4.3	Υ	Absent		NYCP51-8260(14)
L2250575-05B	Vial HCl preserved	Α	NA		4.3	Υ	Absent		NYCP51-8260(14)
L2250575-05C	Vial HCl preserved	Α	NA		4.3	Υ	Absent		NYCP51-8260(14)
L2250575-05D	Amber 1000ml unpreserved	Α	7	7	4.3	Υ	Absent		A2-SVOC-8270(7)
L2250575-05E	Amber 1000ml unpreserved	Α	7	7	4.3	Υ	Absent		A2-SVOC-8270(7)
L2250575-06A	Vial HCl preserved	Α	NA		4.3	Υ	Absent		NYCP51-8260(14)



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#### **GLOSSARY**

#### **Acronyms**

**EDL** 

LOQ

MS

DL - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

 Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis

of PAHs using Solid-Phase Microextraction (SPME).

EMPC - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case estimate of the concentration.

EPA - Environmental Protection Agency.

LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LCSD - Laboratory Control Sample Duplicate: Refer to LCS.

LFB - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LOD - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

 - Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

MDL - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

 Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.

MSD - Matrix Spike Sample Duplicate: Refer to MS.

NA - Not Applicable.

NC - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's

reporting unit.

NDPA/DPA - N-Nitrosodiphenylamine/Diphenylamine.

NI - Not Ignitable.

NP - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.

NR - No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile

Organic TIC only requests.

RL - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL

includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the

values; although the RPD value will be provided in the report.

SRM - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the

associated field samples.

STLP - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.

TEF - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.

TEQ - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF

and then summing the resulting values.

TIC - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.

Report Format: DU Report with 'J' Qualifiers



Project Name:STEELWINDS ANNUAL/SEMI-ANNUALLab Number:L2250575Project Number:03.0033579.15Report Date:10/13/22

#### **Footnotes**

1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

#### **Terms**

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Chlordane: The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA,this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

Difference: With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

Final pH: As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

Frozen Date/Time: With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

Gasoline Range Organics (GRO): Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert but

Gasoline Range Organics (GRO): Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

Initial pH: As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

PAH Total: With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benza(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

PFAS Total: With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

Total: With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

#### Data Qualifiers

- A Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentrations of the analyte at less than ten times (10x) the concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G The concentration may be biased high due to matrix interferences (i.e, co-elution) with non-target compound(s). The result should be considered estimated.
- H The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I The lower value for the two columns has been reported due to obvious interference.
- J Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively

Report Format: DU Report with 'J' Qualifiers



Project Name:STEELWINDS ANNUAL/SEMI-ANNUALLab Number:L2250575Project Number:03.0033579.15Report Date:10/13/22

#### **Data Qualifiers**

Identified Compounds (TICs).

- M Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- ND Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.
- **NJ** Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- ${f P}$  The RPD between the results for the two columns exceeds the method-specified criteria.
- Q The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- RE Analytical results are from sample re-extraction.
- S Analytical results are from modified screening analysis.
- The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits.
   (Applicable to MassDEP DW Compliance samples only.)

Report Format: DU Report with 'J' Qualifiers



Project Name:STEELWINDS ANNUAL/SEMI-ANNUALLab Number:L2250575Project Number:03.0033579.15Report Date:10/13/22

#### REFERENCES

Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

#### **LIMITATION OF LIABILITIES**

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Alpha Analytical, Inc. Facility: Company-wide

Department: Quality Assurance

Title: Certificate/Approval Program Summary

ID No.:17873 Revision 19

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#### Certification Information

#### The following analytes are not included in our Primary NELAP Scope of Accreditation:

#### Westborough Facility

EPA 624/624.1: m/p-xylene, o-xylene, Naphthalene

EPA 625/625.1: alpha-Terpineol

EPA 8260C/8260D: NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene;

EPA 8270D/8270E: NPW: Dimethylnaphthalene,1,4-Diphenylhydrazine, alpha-Terpineol; SCM: Dimethylnaphthalene,1,4-Diphenylhydrazine.

SM4500: NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO2, NO3.

#### **Mansfield Facility**

**SM 2540D:** TSS

EPA 8082A: NPW: PCB: 1, 5, 31, 87,101, 110, 141, 151, 153, 180, 183, 187.

EPA TO-15: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

Biological Tissue Matrix: EPA 3050B

#### The following analytes are included in our Massachusetts DEP Scope of Accreditation

#### Westborough Facility:

#### **Drinking Water**

EPA 300.0: Chloride, Nitrate-N, Fluoride, Sulfate; EPA 353.2: Nitrate-N, Nitrite-N; SM4500NO3-F: Nitrate-N, Nitrite-N; SM4500F-C, SM4500CN-CE,

EPA 180.1, SM2130B, SM4500CI-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B

EPA 332: Perchlorate; EPA 524.2: THMs and VOCs; EPA 504.1: EDB, DBCP.

Microbiology: SM9215B; SM9223-P/A, SM9223B-Colilert-QT,SM9222D.

#### Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, EPA 350.1: Ammonia-N, LACHAT 10-107-06-1-B: Ammonia-N, EPA 351.1, SM4500NO3-F, EPA 353.2: Nitrate-N, SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300: Chloride, Sulfate, Nitrate. EPA 624.1: Volatile Halocarbons & Aromatics,

EPA 608.3: Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan II, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

EPA 625.1: SVOC (Acid/Base/Neutral Extractables), EPA 600/4-81-045: PCB-Oil.

Microbiology: SM9223B-Colilert-QT; Enterolert-QT, SM9221E, EPA 1600, EPA 1603, SM9222D.

#### Mansfield Facility:

#### **Drinking Water**

EPA 200.7: Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. EPA 200.8: Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. EPA 245.1 Hg. EPA 522, EPA 537.1.

#### Non-Potable Water

EPA 200.7: Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.

EPA 200.8: Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.

EPA 245.1 Hg

SM2340B

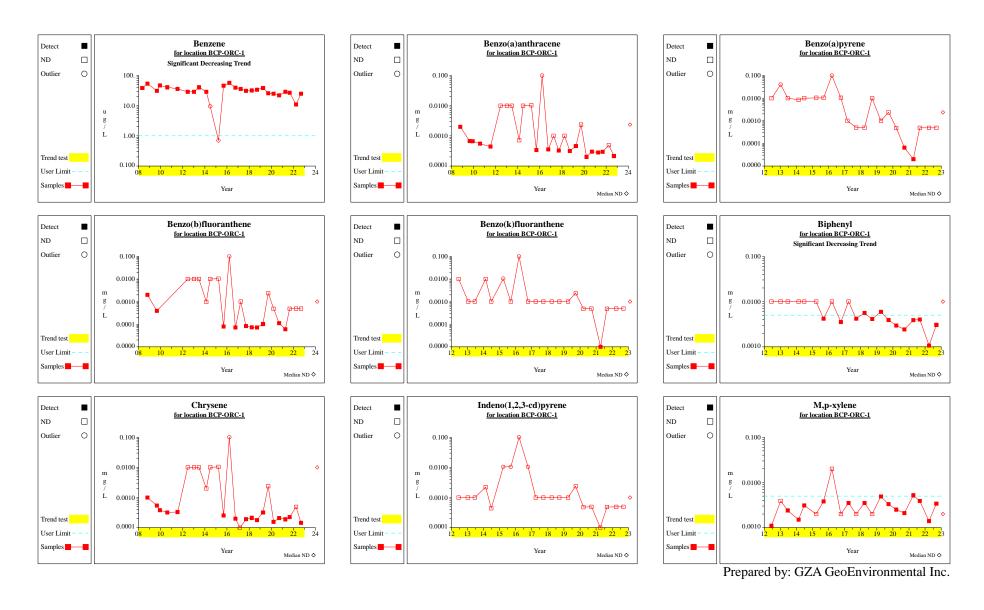
For a complete listing of analytes and methods, please contact your Alpha Project Manager.

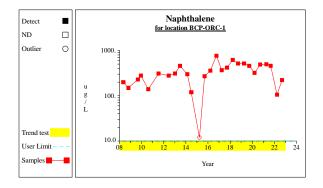
Pre-Qualtrax Document ID: 08-113 Document Type: Form

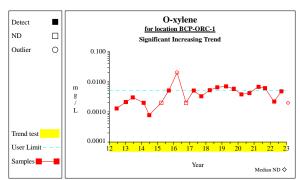
Westborough, MA 01581 8 Walkup Dr. TEL: 508-898-9193  Client Information Client: G-ZA  Address: 300 Peor 1  Buffalo, NU Phone: 716) 51	7 14202	Service Centers Mahwah, NJ 07430: 35 Whitne Albany, NY 12205: 14 Walker I Tonawanda, NY 14150: 275 Co  Project Information  Project Name: Saco  Project Location: Loc  Project # 03 003  (Use Project name as P  Project Manager: DA  ALPHAQuote #:  Turn-Around Time  Standare	Way poper Ave, Suite 1  Was 5  Chawan, 3579, 15  Project #)	Annal) A NY	Spiritor	,	Delivi	Date I in L erables ASP-/ EQuis Other latery NY TO AWQ S NY Re	A (1 Fill A) Requir GS Standar stricted	le) - CA reinten ds Use		ASP-I	B S (4 File rt 375	e)	ALPHA Job # L
	BOY DOZA COA	Rush (only if pre approved		Due Date	333										
	The second secon		u) [	# of Days	S			_	ewer D	ischarg	je			-	Other:
	een previously analyze						ANA	LYSIS	_	_		_	-	-	Sample Filtration
Please specify Metals	c requirements/comm	ents:					STARS	12	· · · · · ·	Δ	D 3k		D CR		Done t a a a a a a a a a a a a a a a a a a
ALPHA Lab ID	Sa	mple ID	Coll	ection	Sample	Sampler's	9260	8270	60 10	900	0009	2	6010		1
(Lab Use Only)		pio io	Date	Time	Matrix	Initials	92	82	9	3	3	3	ĕ		Sample Specific Comments
50575-01	MWN-030-0	91522	9-15-22	0830	6w	FN	×	×			×	×			JLAB FILLET
02	MWN-038 -09	11522	1	0945	1	4				×	×	×	×		Metal Samples
03	MWN-03-091	522		1030			×	X							
04	MWN-020-09	11522		1345					X						- LAB FILLER
05	MWN-04-09	1522		1450	1		¥	X							
06	TRIP BLANK-	3	1		W		×								
							0.50								
Preservative Code: A = None B = HCl C = HNO <sub>3</sub> D = H <sub>2</sub> SO <sub>4</sub> E = NaOH	Container Code P = Plastic A = Amber Glass V = Vial G = Glass B = Bacteria Cup	Westboro: Certification N Mansfield: Certification N			_	tainer Type									Please print clearly, legibly and completely. Samples can not be logged in and turnaround time clock will not start until any ambiguities are
F = MeOH	C = Cube O = Other	Relinquished	By:	Date	/Time		Receiv	red By				Date	Time		resolved. BY EXECUTING
$G = NaHSO_4$ $H = Na_2S_2O_3$ K/E = Zn Ac/NaOH O = Other	E = Encore D = BOD Bottle	What I LIAL		9-15-22		MA	1		912	_	9/1	-	200	_	THIS COC, THE CLIENT HAS READ AND AGREES TO BE BOUND BY ALPHA'S TERMS & CONDITIONS.
Form No: 01-25 HC (rev. 3	0-Sept-2013)														(See reverse side.)

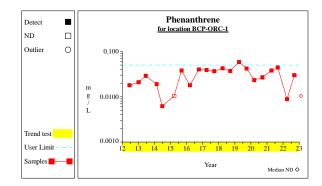


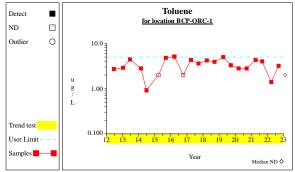
# APPENDIX C TIME SERIES PLOTS

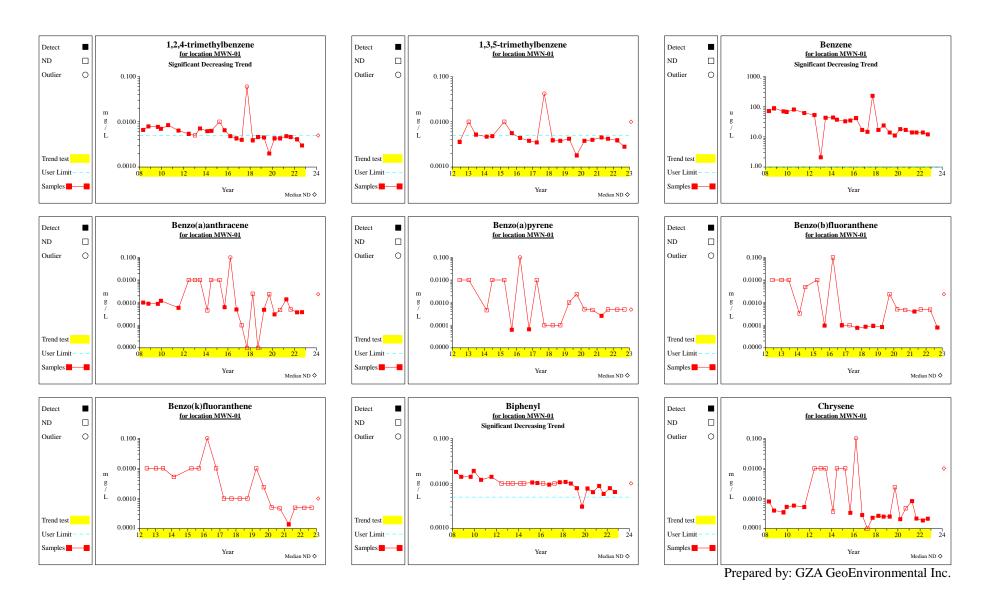


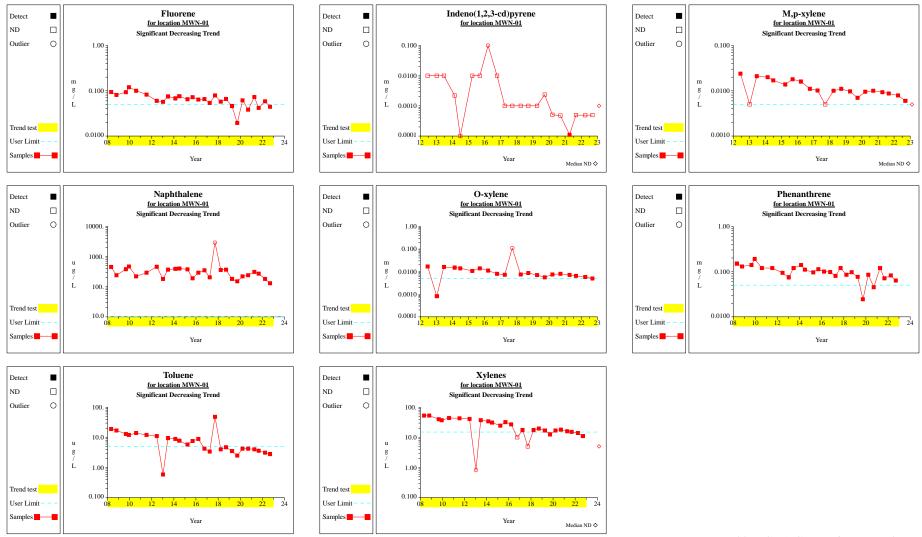




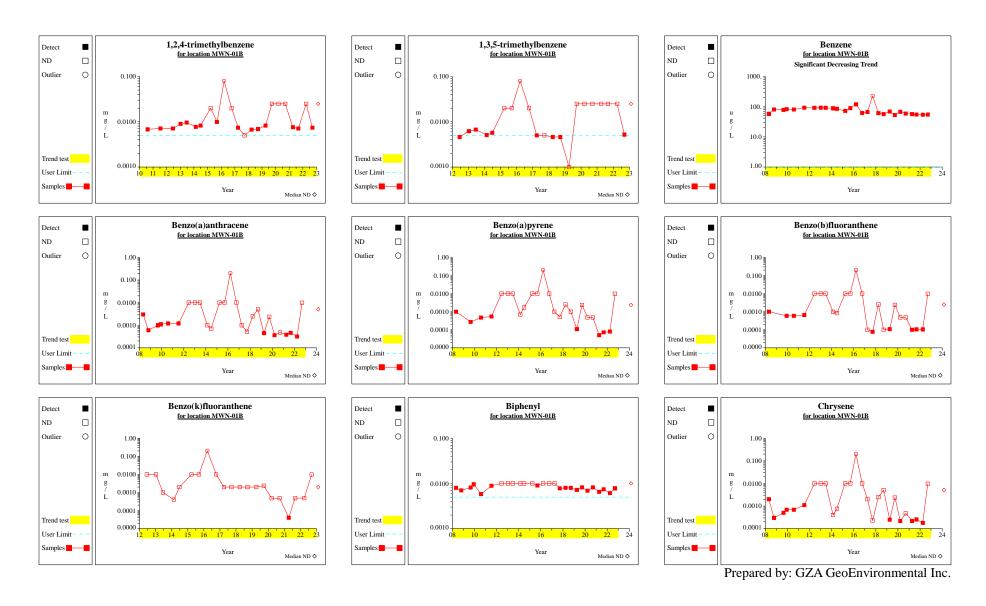








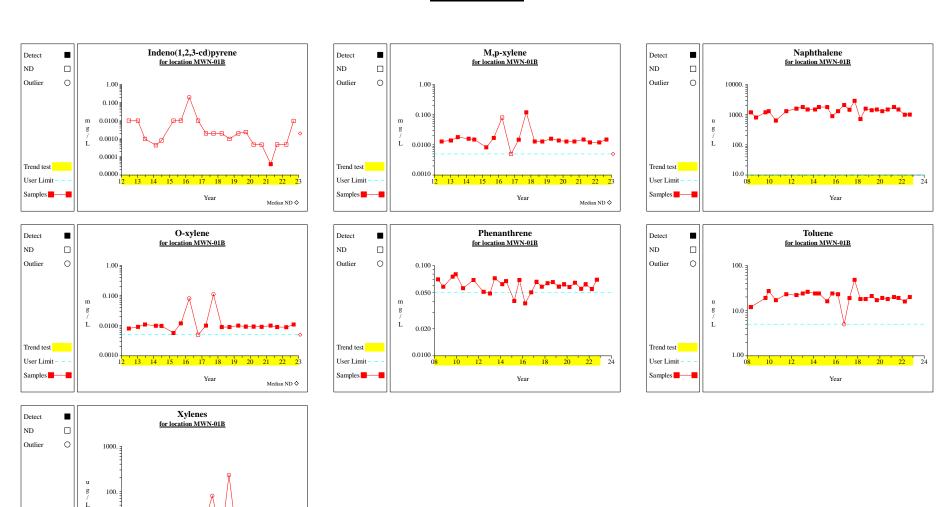
Prepared by: GZA GeoEnvironmental Inc.

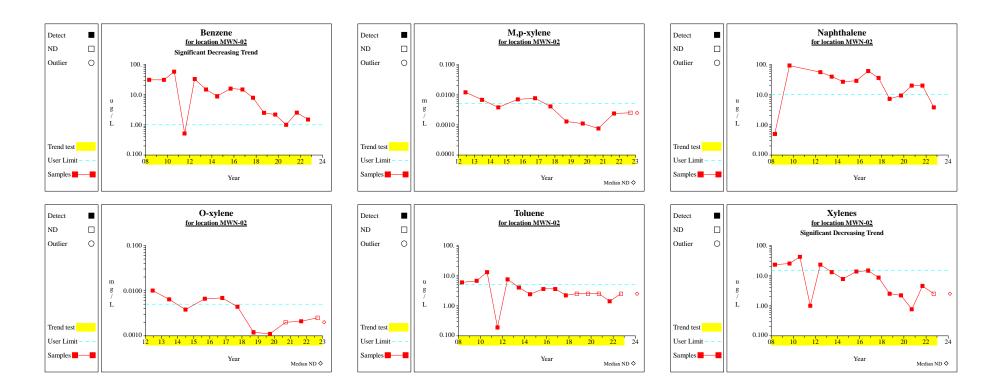


Trend test

Year

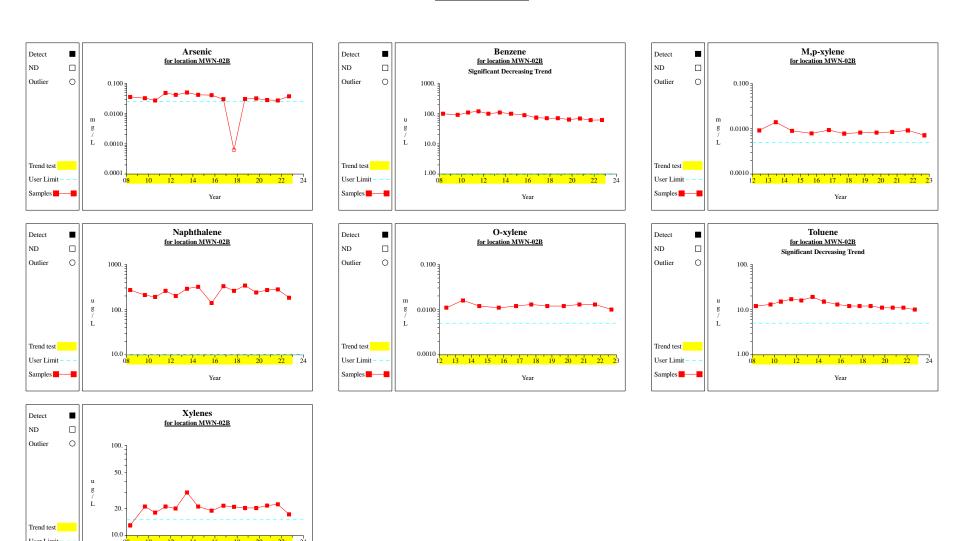
Median ND ♦



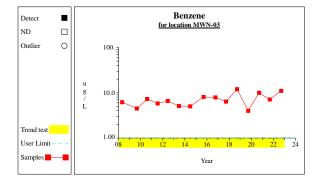


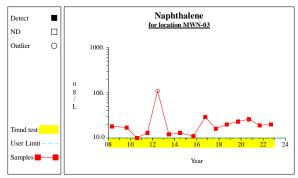
Samples \_\_\_\_

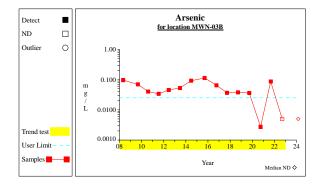
Year

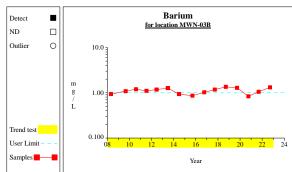


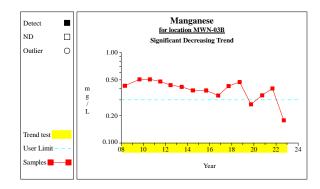
Prepared by: GZA GeoEnvironmental Inc.

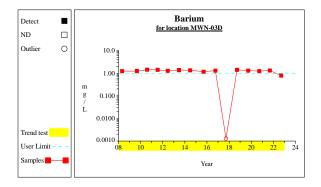


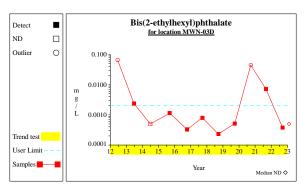


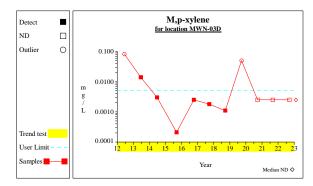


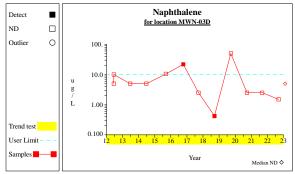


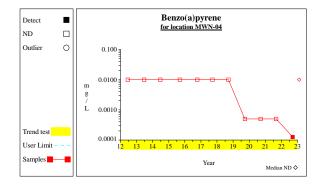


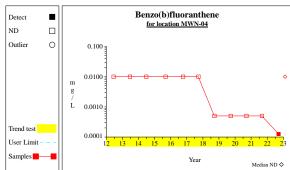


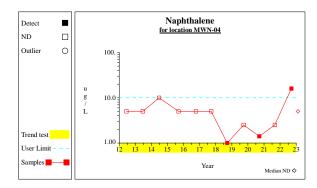


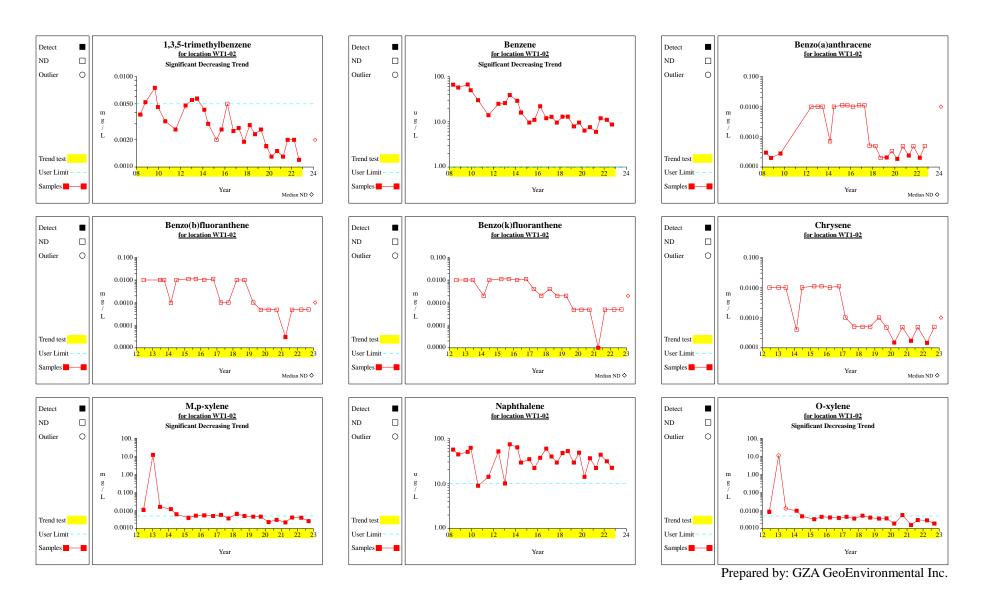


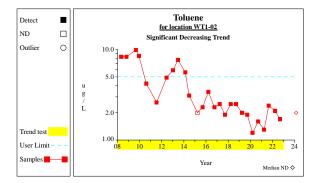


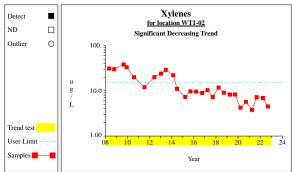


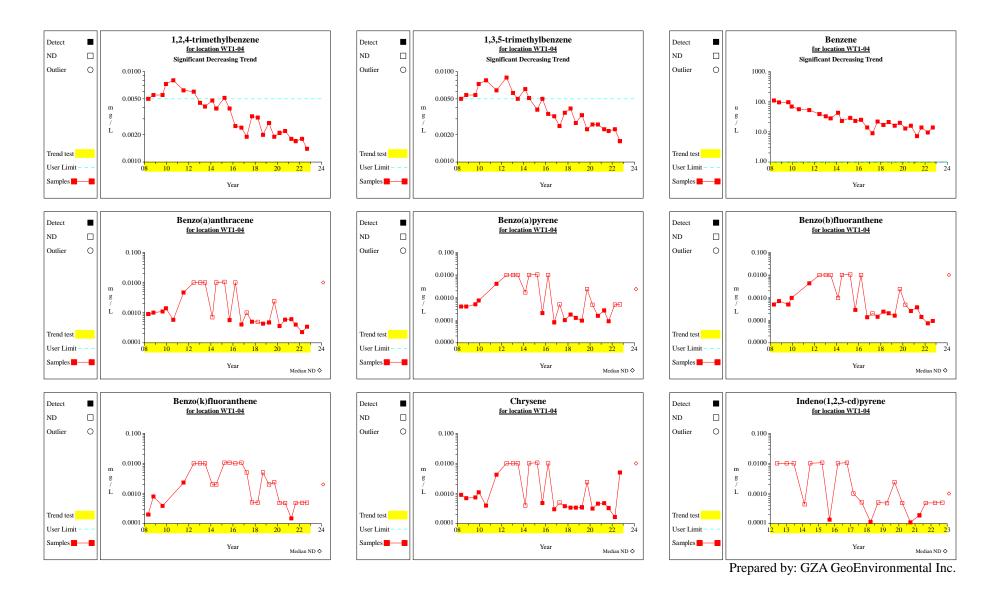


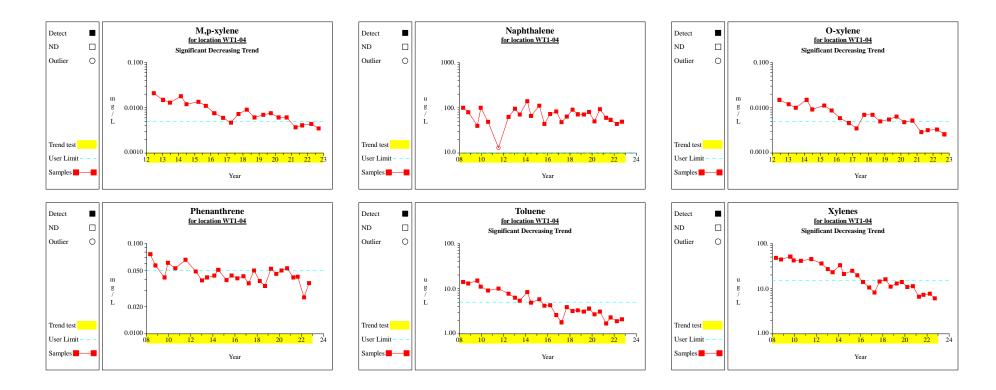


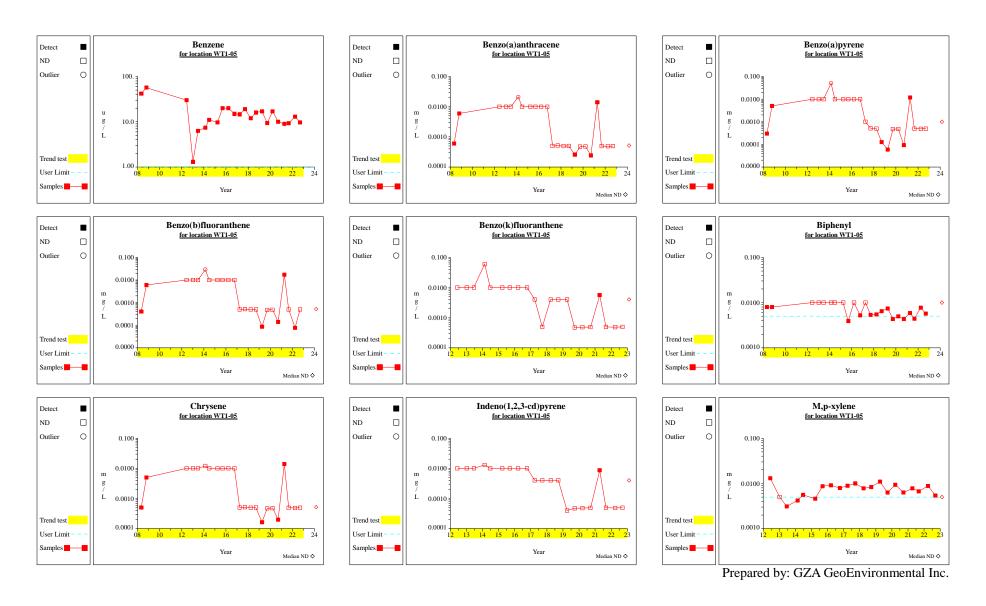


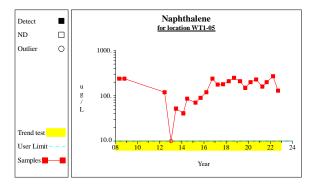


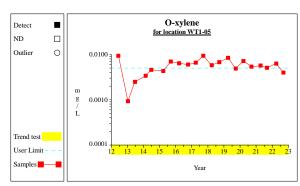


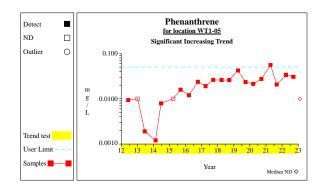


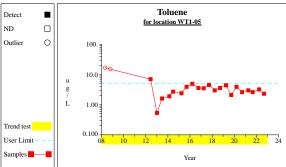


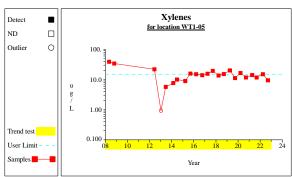














# APPENDIX D WELL DEVELOPMENT FORMS

						Historic Info	ormation						
Boring Log A	vailable (v	ves/no/attac	:hed):			Thotorio iini	Jillianoll						
Installation L			•										
motanation E	og / wallak	510 ( <b>ycs</b> /110/0	attaoricaj			Summ	arv.						
Monitoring W	/oll ·	MWN-01		Ground Sur	rface Elevation:	582.99	,	Dicor/So	reen Material	· D\/C			
Installation D		8/30/90		•	er Elevation:	569.92			creen Depth:				
Installed By:	ale.	Turnkey			Point Elevation:			•	of Screen Dep				
iristalied by.		Turrikey		Elevation D		303.14		Dollonic	n Screen Dep	<u>un. 19.15</u>			
Provious Fiel	ld maacur	ement Infor	mation Availa										
r revious i lei	iu ilieasui	ement inion	Tialion Availa	Jie (yes/110/	,	of Dravious F	iold Moonin	romonto					
Don'th to	<b>VA</b> / = 1 = =	1	- 1.1	0		s of Previous F				I	0.1		
Depth to			pH	•	Conductance	Tempera			urbidity		Color		
(ft)		`	ard Units)	`	mS/cm)	(°C		,	NTU)		Olaran		
11.5	i9	11.53	3 - 11.67	1.0	51 -1.450	10.4 -1	4.4	1.2	23 - 2.9		Clear		
Notes:													
			F	ield Observa	ations				Parame	eters +/-	Sampling Information		
Exterior Obse	ervations:	OK							pН	+/- 0.1	Sample ID: <b>MWN-01</b> -091322		
		-							Conductivity		Sample Time: 09:05		
nterior Obse	ervations								# of Sample Containers: Five				
									Turbidity	+/- 10%	Duplicate Sample ID: NA		
									ORP	+/- 10mV	Sample Analysis:		
Signs of Dan			None						DO	+/- 10%	VOC STARS List via EPA 8260C		
Locked (y	/es/ <b>no</b> )	Well Ca	ıp ( <b>yes</b> /no)	Sur	face Seal Intact (		PID Measi	urement:	ND	Odors:	SVOC B/N Via EPA 8270D		
						Well Quali	ty Data						
										_			
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes		
		Water	Volume	(Standard	Conductance	(°C)	(NTU)		Oxygen	Reduction			
			Purged (Gal)		(mS/cm)				%	Potential			
9/13/2022	8:40	15.49	0	11.71	1.244	11.5	3.48	Clear	9.8	22.7	Depth of Water: 15.22		
	8:50	15.52	10	11.8	1.25	11.9	2.64	Clear	6.5	-61.5	Length of Water Column: 3.93		
	8:55	15.52	15	11.8	1.252	12.0	2.84	Clear	6.2	-90.1	Depth of Well: 19.15		
	9:00	15.52	20	11.81	1.255	12.0	2.83	Clear	6.0	-98.7	Sheen Observed: Y N		
	9:05	15.52	25	11.81	1.258	12.0	2.80	Clear	5.9	-104.5	DNAPL Observed: Y N		
											Did Well Go Dry: Y N		
											Other:		
											1 Well Volume = 2.5 gal		
											Dome: 4		

						Historic Inf	ormation						
Boring Log Av	vailable ( <b>v</b>	<b>/es</b> /no/attac	:hed):			1							
nstallation Lo	1.5		,										
	og / tranac	) ( <b>) 00</b> /110/0	attaonoay			Summ	narv						
Monitoring W	اما ·	MWN-01B		Ground Su	rface Elevation:		iary	Riser/Sc	reen Material	· P\/C			
Installation Da		11/2/92		1	er Elevation:	571.81		_	creen Depth:				
Installed By:	ato.	Turnkey		•	Point Elevation:			_ :	of Screen Dep				
motanou by:		ranney		Elevation D		007.00			. 00.00 Dop	02.2 1			
Previous Field	d measure	ement Infori	mation Availal										
	<u> </u>			() 55/116/	,	s of Previous F	ield Meası	ırements					
Depth to \	Water		рН	Specific	Conductance	Tempera			urbidity		Color		
(ft)	vvaloi		ard Units)	•	nS/cm)	(°C			NTU)		30101		
15.5		`	-11.40	`	31 - 1.01	9.8 - 1		,	0 - 7.67		Clear		
	)	11.1	-11.40	0.6	51 - 1.01	9.0 - 1	١.٥	1.2	.0 - 7.07		Clear		
Notes:													
				ald Observe	tions				Parage	otoro I	Compositor or traffic area attach		
E tarian Olam		01/	FI	eld Observa	ations					eters +/-	Sampling Information		
Exterior Obse	ervations:	<u>OK</u>							pH Conductivity	+/- 0.1	Sample ID: MWN-01B-091322 Sample Time: 10:15		
Interior Obse	m rotions	OK							Temperature		•		
interior Obsei	rvations	OK							Turbidity	+/- 10%	# of Sample Containers: Five		
									ORP	+/- 10% +/- 10mV	Duplicate Sample ID: NA Sample Analysis:		
Signs of Dam	ago/Tami	noring:	None						DO	+/- 101117	VOC STARS List via EPA 8260C		
Locked (ye			p ( <b>yes</b> /no)	Surf	ace Seal Intact (	(ves/no)	PID Meas	urement.		Odors:	SVOC B/N Via EPA 8270D		
Locked (ye	G3/11 <b>U</b> )	vven oa	ip ( <b>yes</b> /110)	Ouri	ace Sear Illiact	Well Qual		urement.	ND	Odors.	SVOC B/N VIA ET A 0270B		
T		I	I			I Well Qual	l Data	l l	T T	1			
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes		
24.0		Water	Volume	(Standard	Conductance	(°C)	(NTU)		Oxygen	Reduction	110.00		
			Purged (Gal)	`	(mS/cm)	( )	(1110)		%	Potential			
9/13/2022	9:50	15.29	0	11.21	0.776	10.8	78.2	Clear	7.2	-28.5	Depth of Water: 15.22		
5. 15,2522	10:00	15.29	4	11.42	0.865	10.7	35.88	Clear	10.4	-98.7	Length of Water Column: 17.02		
	10:05	15.29	6	11.44	0.872	10.6	27.12	Clear	10.8	-107.2	Depth of Well: 32.24		
	10:10	15.29	8	11.46	0.888	10.6	26.32	Clear	11.2	-114.2	Sheen Observed: Y N		
	10:15	15.29	10	11.46	0.891	10.6	22.18	Clear	11.3	-118.8	DNAPL Observed: Y N		
											Did Well Go Dry: Y N		
											Other:		
											1 Well Volume = 2.7 gal		

							•				
						Historic Info	ormation				
Boring Log A	vailable (	<b>yes</b> /no/attac	ched):								
nstallation L	og Availat	ole ( <b>yes</b> /no/a	attached)								
						Summ	ary				
Monitoring W	/ell :	WT1-02		Ground Su	rface Elevation:	598.5		Riser/Sc	reen Materia	I: PVC	
Installation D		6/11/07		Groundwat	er Elevation:	572.96		Top of S	creen Depth:	27.78	
Installed By:		Turnkey		Monitoring	Point Elevation:	600.78		Bottom o	of Screen De	oth: 37.78	
				Elevation D	atum:						
Previous Fiel	ld measur	ement Infor	mation Availal	ole (yes/ <b>no</b> /	attached)						
					Range	s of Previous F	ield Measu	rements			
Depth to	Water		рН	Specific	Conductance	Tempera	ature	Tu	ırbidity		Color
(ft)		(Standa	ard Units)	· ·	nS/cm)	, (°C			NTU)		
26.9		` `	5 - 12.19	· ·	7 - 2.116	11.5 - 1		`	67 - 16		Clear
Notes:										<u> </u>	
			Fi	ield Observa	ations				Parame	eters +/-	Sampling Information
Exterior Obs	ervations:	OK							pH	+/- 0.1	Sample ID: <b>WT1-02</b> -091322
									Conductivity		Sample Time: 14:15
nterior Obse	ervations	OK							Temperature +/- 109		# of Sample Containers: Five
									Turbidity	+/- 10%	Duplicate Sample ID: NA
									ORP	+/- 10mV	Sample Analysis:
Signs of Dan	nage/Tam	pering:	None						DO	+/- 10%	VOC STARS List via EPA 8260C
Locked (y	/es/ <b>no</b> )	Well Ca	ap ( <b>yes</b> /no)	Surf	ace Seal Intact (	( <b>yes</b> /no)	PID Measu	urement:	ND	Odors:	SVOC B/N Via EPA 8270D
						Well Quali	ity Data				
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes
		Water	Volume	(Standard	Conductance	(°C)	(NTU)		Oxygen	Reduction	
			Purged (Gal)		(mS/cm)				%	Potential	
9/13/2022	13:55	28.45	0	11.93	1.605	13	1.75	Clear	24.1	28.3	Depth of Water: 27.82
	14:05	28.67	2	11.98	1.588	13.1	1.46	Clear	8.0	-30.7	Length of Water Column: 9.96
	14:10	28.67	3	11.97	1.589	13.1	1.40	Clear	7.7	-36.4	Depth of Well: 37.78
	14:15	28.67	4	11.98	1.592	13.1	1.43	Clear	7.6	-41.2	Sheen Observed: Y N
											DNAPL Observed: Y N
											Did Well Go Dry: Y N
											Other:
											1 Well Volume = 6.5 gal.

						Historic Inf	ormation				
Boring Log A	vailable (	yes/no/attac	:hed):								
Installation L			,								
	- 9 - 11 - 11 - 11	( <b>)</b> CO. 113.				Summ	narv				
Monitoring W	/ell :	WT1-04		Ground Su	rface Elevation		··· ,	Riser/Sc	reen Materi	al: PVC	
Installation D		5/21/07			er Elevation:	572.69		_	creen Depth		
Installed By:		Turnkey			Point Elevation				of Screen De		
,				Elevation D				•			
Previous Fiel	ld measur	ement Infor	mation Availal	ole (ves/ <b>no</b> /	attached)						
				()		s of Previous F	ield Measu	rements			
Depth to	Water		pН	Specific (	Conductance	Tempera			rbidity		Color
(ft)			ard Units)	•	S/cm)	O°C			NTU)		20.0.
12.9		`	- 11.99	,	6 - 1.550	10.2 -1:		`	1 - 8.3		Clear
Notes:	, ı	11.31	11.55	1.520	1.000	10.2 -1.	0.21	1.5	1 0.0	ļ	Olcai
INULES.											
			Fiz	eld Observa	tions				Param	eters +/-	Sampling Information
Exterior Obs	orvotions:	OK	1 10	du Observa	tions				pH	+/- 0.1	Sample ID: <b>WT1-04</b> -091322
Exterior Obse	ervalions.	Conductivity +/- 3 <sup>s</sup>									Sample Time: 10:55
nterior Obse	arvations	OK Temperature +/- 1								# of Sample Containers: Five	
intenoi Obse	o valions	OK							Turbidity		Duplicate Sample ID: NA
									Sample Analysis:		
Signs of Dan	nage/Tam	nering:									VOC STARS List via EPA 8260C
Locked (y		<del></del>	p (yes/no)	Surfa	ace Seal Intact	(ves/no)	PID Measi	urement:	ND	Odors:	SVOC B/N Via EPA 8270D
	, 33, 110,		( <b>) (3)</b>	3		Well Qual				10 00.0.	
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes
		Water	Volume	(Standard	Conductance	( °C)	(NTU)		Oxygen	Reduction	
			Purged (Gal)	`	(mS/cm)				%	Potential	
9/13/2022	10:35	14.22	0	11.73	1.682	13.1	4.50	Clear	7.6	-18	Depth of Water: 13.76
	10:40	14.30	1	11.78	1.400	13.0	2.62	Clear	6.1	-62.1	Length of Water Column: 11.76
	10:45	14.30	1.5	11.77	1.343	13.1	3.67	Clear	5.7	-112.2	Depth of Well: 25.52
	10:50	14.30	2	11.75	1.336	13.2	3.57	Clear	5.6	-116.2	Sheen Observed: Y N
	10:55	14.30	2.5	11.75	1.326	13.3	3.80	Clear	5.5	-118.5	DNAPL Observed: Y N
											Did Well Go Dry: Y N
											Other:
											1 Well Volume = 1.9 gal

						Historic Inf	ormation				
Boring Log A	vailable (	yes/no/attac	hed):								
Installation L			,								
	- 9	· · · (J 30/				Summ	narv				
Monitoring W	/ell :	WT1-05		Ground Su	rface Elevation		,	Riser/Sc	reen Materi	al: PVC	
Installation D		5/29/07			er Elevation:	571.93		_	creen Depth		
Installed By:		Turnkey			Point Elevation				of Screen De		
,				Elevation D				•			
Previous Fiel	ld measur	ement Infor	mation Availal	ole (yes/ <b>no</b> /	attached)						
					Range	s of Previous F	ield Measu	rements			
Depth to	Water		рН	Specific (	Conductance	Tempera			rbidity		Color
(ft)			ard Units)	-	hos/cm)	O°C			NTU)		
11.8		,	5 - 11.93		) - 1.490	10.4 - 1		`	4 - 5.3		Clear
Notes:		11.00	7 11.00	1.20	1.100	10.1	2.01	1.7	1 0.0		Oldai
110153.											
			Fie	eld Observa	tions				Param	eters +/-	Sampling Information
Exterior Obs	ervations:	OK	110	old Obool va	110110				Hq	+/- 0.1	Sample ID: <b>WT1-05</b> -091322
	J. VALIO113.										Sample Time: 7:55
Interior Obse	ervations	OK									# of Sample Containers: Five
									Turbidity		Duplicate Sample ID: NA
									ORP		Sample Analysis:
Signs of Dan	nage/Tam	pering:	None						DO	+/- 10%	VOC STARS List via EPA 8260C
Locked (y		<del></del>	ıp ( <b>yes</b> /no)	Surfa	ace Seal Intact	(yes/no)	PID Measi	urement:	ND	Odors:	SVOC B/N Via EPA 8270D
						Well Qual	ity Data				
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes
		Water	Volume	(Standard	Conductance	( °C)	(NTU)		Oxygen	Reduction	
		ft bgs	Purged (Gal)	Units)	(mS/cm)				%	Potential	
9/13/2022	7:36	12.50	0	11.51	1.347	12.8	1.52	Clear	8.8	194.6	Depth of Water: 12.48
	7:45	12.50	9	11.53	1.306	13.0	1.00	Clear	6.2	-58.6	Length of Water Column: 10.82
	7:50	12.50	14	11.56	1.299	13.0	1.11	Clear	5.9	-60.5	Depth of Well: 23.3
	7:55	12.50	19	11.61	1.292	13.0	0.98	Clear	5.7	-68.7	Sheen Observed: Y N
											DNAPL Observed: Y N
											Did Well Go Dry: Y N
											Other:
											1 Well Volume = 1.7 gal
								-	-	ļ	

						Historic Inf	ormation				
Boring Log A	vailable ( <b>v</b>	es/no/attac	hed).			THOUGHO IIII	omation				
Installation L	1.5		,								
motanation E	og Avallak	olo ( <b>yes</b> /110/6	attacrica)			Summ	narv				
Monitoring W	Iall :	BCP-ORC	1	Cround Su	rface Elevation		iaiy	Dicor/Co	reen Materia	ol: DVC	
Installation D		10/3/07	- 1		er Elevation:	572.74			creen Depth		
Installed By:	ale.	Turnkey			Point Elevation				of Screen Depti		
ilistalied by.		Turrikey		Elevation D		391.91		_ Dolloin C	n Scieen De	spiii. 54.00	
Provious Fiel	ld moocur	omont Infor	mation Availal								
r levious i lei	iu ilieasui	ement inton	nation Availat	ole (yes/110/		s of Previous F	iold Massu				
Don'th to	10/-1		- 1.1	0					.l. 1.19	ı	2.1.
Depth to			pH	•	Conductance	Tempera			rbidity		Color
(ft)		`	ard Units)	,	iS/cm)	O°)			NTU)		
18.3	80	11.21	- 11.57	0.95	7 - 1.360	10.0 - 1	2.02	2.1	- 11.4		Clear
Notes:											
			Fi∈	eld Observa	tions					eters +/-	Sampling Information
Exterior Obse	ervations:	OK							рН	+/- 0.1	Sample ID: BCP-ORC-091322
		Conductivity +/- 3%								Sample Time: 13:25	
Interior Obse	ervations								# of Sample Containers: Five		
		Turbidity +/- 10							Duplicate Sample ID: NA		
									ORP		Sample Analysis:
Signs of Dam			None				1		DO	+/- 10%	VOC STARS List via EPA 8260C
Locked (y	/es/ <b>no</b> )	Well Ca	p ( <b>yes</b> /no)	Surfa	ace Seal Intact		PID Meas	urement:	ND	Odors:	SVOC B/N Via EPA 8270D
						Well Qual	ity Data				
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes
		Water	Volume	`	Conductance	(°C)	(NTU)		Oxygen	Reduction	
			Purged (Gal)		(mS/cm)				%	Potential	
9/13/2022	13:05	20.55	0	11.78	1.105	11.7	1.12	Clear	15.4	21.7	Depth of Water: 19.23
	13:15	20.60	1	11.61	1.041	11.5	1.74	Clear	10.9	34.0	Length of Water Column: 15.45
	13:20	20.60	1.5	11.61	1.052	11.5	1.52	Clear	11.3	28.2	Depth of Well: 34.68
	13:25	20.60	2	11.60	1.060	11.5	1.56	Clear	11.0	20.7	Sheen Observed: Y N
								ļ			DNAPL Observed: Y N
											Did Well Go Dry: Y N
											Other: Sulfur odor.
											1 Well Volume = 2.5 gal.
		l									

						Historic Inf	ormation				
Boring Log A	vailable (y	<b>/es</b> /no/attac	hed):								
nstallation Lo	1.5		,								
	- 0	()	,			Summ	narv				
Monitoring W	'ell :	MWN-02		Ground Su	rface Elevation		,	Riser/Sc	reen Materi	al: PVC	
nstallation D		9/10/90			er Elevation:	572.52		_	creen Depth		
nstalled By:		Turnkey			Point Elevation				of Screen De		
·				Elevation D				•		•	
Previous Fiel	d measure	ement Inforr	mation Availal	ole (yes/ <b>no</b> /	attached)						
				-	Range	s of Previous F	ield Measu	rements			
Depth to	Water		рН	Specific (	Conductance	Tempera			rbidity		Color
· (ft)			ard Units)	•	s/cm)	Ċ°C			NTU)		
27.7		`	) - 12.31	,	3 - 2.06	11.3 - 1		`	9 - 38.6		Clear
Notes:					•	-				!	
			Fie	eld Observa	tions				Param	eters +/-	Sampling Information
xterior Obse	ervations:	OK							рН	+/- 0.1	Sample ID: <b>MWN-02</b> -091422
									Conductivi		Sample Time: 08:00
nterior Obse	rvations	OK									# of Sample Containers: Five
									Turbidity		Duplicate Sample ID: NA
									ORP		Sample Analysis:
Signs of Dam	nage/Tamp		None						DO	+/- 10%	VOC STARS List via EPA 8260C
Locked (y	es/ <b>no</b> )	Well Ca	ıp ( <b>yes</b> /no)	Surfa	ace Seal Intact	7	PID Measi	urement:	ND	Odors:	SVOC B/N Via EPA 8270D
						Well Qual	ity Data				
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	,	Color	Dissolved	Oxygen	Notes
		Water	Volume	(Standard	Conductance	( °C)	(NTU)		Oxygen	Reduction	
			Purged (Gal)		(mS/cm)				%	Potential	
9/14/2022	7:40	28.58	0	11.71	1.853	12.7	2.16	Clear	28.9	182.6	Depth of Water: 28.49
	7:50	28.58	5	11.79	1.934	12.5	2.50	Clear	14.2	154.9	Length of Water Column: 5.13
	7:55	28.58	1.5	11.81	1.954	12.6	2.62	Clear	13.8	146.7	Depth of Well: 33.62
	8:00	28.58	10	11.85	1.965	12.6	2.54	Clear	13.6	137.8	Sheen Observed: Y N
											DNAPL Observed: Y N Did Well Go Dry: Y N
											Did Well Go Dry: Y N Other:
											1 Well Volume = 3.3 gal.
										 	i vven volunie – 3.3 gai.
		I				ĺ					

						Historic Inf	ormation				
Boring Log A	vailable (	ves/no/attac	hed).			THSTOTIC IIII	Offialion				
Installation L			,								
IIIStallation L	og Avallai	ole ( <b>yes</b> /110/	attacrieu)			Cuma	2011				
N d = usid = usid = usid = usid	/all a B	Summary  MWN-02B Ground Surface Elevation: 599.00 Riser/Screen Material: PVC									
Monitoring W		/WN-02B			rrace Elevation er Elevation:			_			
	nstallation Date: 11/2/92								creen Depth		
Installed By:		Turnkey			Point Elevation	601.28 Bott		_Bottom C	of Screen De	ptn: 56.28	
D				Elevation D							
Previous Fiel	a measur	ement infori	mation Availal	ole (yes/ <b>no</b> /							
		T				s of Previous F				•	
Depth to			pН	•	Conductance	Tempera			rbidity		Color
(ft)		•	ard Units)	,	S/cm)	O°)		,	NTU)		
28.0	5	11.30	) - 11.75	0.91	0 - 1.13	12.1 - 1	13.4	1.7	6 - 6.9		Clear
Notes:	_										
			Fie	eld Observa	tions				Param	eters +/-	Sampling Information
Exterior Obse	ervations:	OK							рН	+/- 0.1	Sample ID: <b>MWN-02B</b> -091422
									Conductivi	ty +/- 3%	Sample Time: 8:45
Interior Obse	rvations	OK							Temperatu	re +/- 10%	# of Sample Containers:
									Turbidity	+/- 10%	Duplicate Sample ID: NA
									ORP	+/- 10mV	Sample Analysis: Arsenic
Signs of Dam	nage/Tam	pering:	None						DO	+/- 10%	VOC STARS List via EPA 8260C
Locked (y	/es/ <b>no</b> )	Well Ca	ıp ( <b>yes</b> /no)	Surfa	ace Seal Intact	(yes/no)	PID Meas	urement:	ND	Odors:	SVOC B/N Via EPA 8270D, arsenic
						Well Qual	ity Data				
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes
		Water	Volume	(Standard	Conductance	(°C)	(NTU)		Oxygen	Reduction	
		ft bgs	Purged (Gal)	Units)	(uMhos/cm)				%	Potential	
9/14/2022	8:24	29.65	0	11.07	0.833	13.2	1.48	Clear	11.0	114.0	Depth of Water: 28.73
	8:35	29.33	5	11.18	0.902	13.7	2.88	Clear	7.0	-42.6	Length of Water Column: 27.55
	8:40	29.33	7.5	11.17	0.897	13.9	2.61	Clear	6.8	-48.9	Depth of Well: 56.28
	8:45	29.33	10	11.17	0.902	13.9	2.57	Clear	6.5	-56.1	Sheen Observed: Y N
											DNAPL Observed: Y N
											Did Well Go Dry: Y N
											Other: Sulfur odor.
											1 Well Volume = 4.4 gal.

						Historic Inf	ormation					
Boring Log A	vailable (v	ves/no/attac	hed).			Thistoric init	omation					
Installation L		•	,									
IIIStallation L	og Avallai	ole ( <b>yes</b> /110/8	attached)			Summ	oorv.					
Monitoring W	Summary g Well: MWN-02D Ground Surface Elevation: 600.61 Riser/Screen Material: PVC											
Installation D		8/4/95			er Elevation:	573.65		_				
Installed By:	ale.	Turnkey			Point Elevation			Top of Screen Depth: 74.34  Bottom of Screen Depth: 79.34				
ilistalled by.		Turrikey		Elevation D				_ Bollom C	JI SCIECTI DE	pin. 79.34		
Drovious Fiel	ld magazir	omont Infor	mation Availat									
FIEVIOUS FIEI	iu illeasui	ement inion	nation Availat	ne (yes/110/		<del></del>	"alal Manas					
5 4 4	147			0 :::		s of Previous F			1.1.1.			
Depth to		1	рН	•	Conductance	Tempera			rbidity		Color	
(ft)		`	ard Units)		nS/cm)	O°)		,	NTU)			
28.8	_		- 7.42		54 - 2.08	12.6 - 1	-		- 15.1		Clear	
Notes:	**Monso	on pump ur	nable to purg	e water. 3 v	well volumes p	ourged with de	edicated ba	ailer.				
			Fi∈	eld Observa	tions				Param	eters +/-	Sampling Information	
Exterior Obse	ervations:	OK							рН	+/- 0.1	Sample ID: <b>MWN-02D</b> -091422	
									Conductivi	•	Sample Time: 13:45	
Interior Obse	ervations	OK									# of Sample Containers: One	
									Turbidity		Duplicate Sample ID: NA	
									ORP		Sample Analysis: Barium, Arsenic	
Signs of Dam			None				1		DO		Chromium	
Locked (y	res/ <b>no</b> )	Well Ca	p ( <b>yes</b> /no)	Surfa	ace Seal Intact	( <b>J</b> /	PID Meas	urement:	ND	Odors: Sulf	ur	
						Well Qual	ity Data					
_										_		
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	,	Color	Dissolved	Oxygen	Notes	
		Water	Volume	`	Conductance	( °C)	(NTU)		Oxygen	Reduction		
			Purged (Gal)		(mS/cm)				%	Potential		
9/14/2022	13:45	32.62	24**	7.86	2.027	13.8	189.3	Gray	29.9	49.8	Depth of Water: 29.30	
											Length of Water Column: 50.04	
											Depth of Well: 79.34	
											Sheen Observed: Y N	
											DNAPL Observed: Y N	
											Did Well Go Dry: Y N	
											Other: Sulfur odor.	
											1 Well Volume = 8.0 gal.	
								1				

						Historic Inf	ormation				
Boring Log A	vailahle ( <b>v</b>	ves/no/attac	hed).			Thistoric irii	Omalion				
Installation Log		•	,								
IIIStallation E	og Avallat	oic ( <b>yes</b> /110/6	attacrica)			Summ	narv				
Monitoring W	ام/	MWN-03		Ground Su	rface Elevation		шу	Riser/Sc	reen Materia	al· P\/C	
Installation Date: 9/6/90					er Elevation:	572.47					
Installed By:	ato.	Turnkey		_ Monitoring Point Elevation				Top of Screen Depth: 39.17  Bottom of Screen Depth: 49.17			
motanou by:		Tarritoy		Elevation D		0.11.00			. 00.00 2	<u> </u>	
Previous Fiel	ld measur	ement Inforr	mation Availat								
				()		s of Previous F	ield Measu	rements			
Depth to	Water		рН	Specific (	Conductance	Tempera			rbidity		Color
(ft)			ard Units)	•	iS/cm)	(°C			NTU)		30101
38.8		`	) - 12.51		24 - 3.04	12.8 - 1		,	- 8.41		Clear
	_				vell volumes p	_	_		0.41	ļ	Oleai
Notes:	MOUSO	on pump ur	iable to purg	e water. 3 V	well volulles t	Jurgeu with at	ulcaleu Da	ailei.			
			Eid	eld Observa	tions				Param	eters +/-	Sampling Information
Exterior Obse	orvations:	OK	1 10	du Observa	tions				pH	+/- 0.1	Sample ID: MWN-03-091522
EXTERIOR ODS	ei valioi is.	OK							Conductivi		Sample Time: 10:30
Interior Obse	rvations	OK								•	# of Sample Containers:
	n valiono	<u> </u>									Duplicate Sample ID: NA
									Turbidity ORP		Sample Analysis:
Signs of Dam	nage/Tam	perina:	None						DO	+/- 10%	VOC STARS List via EPA 8260C
Locked (y			p ( <b>yes</b> /no)	Surfa	ace Seal Intact	(ves/no)	PID Meas	urement:	ND	Odors:	SVOC B/N Via EPA 8270D
· ·						Well Qual	ity Data			•	
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes
		Water	Volume	(Standard	Conductance	( °C)	(NTU)		Oxygen	Reduction	
		ft bgs	Purged (Gal)	Units)	(mS/cm)	, ,			%	Potential	
9/15/2022	10:30	40.63	19**	12.49	3.058	14.3	4.06	Clear	43.0	-39.3	Depth of Water: 39.49
											Length of Water Column: 9.68
											Depth of Well: 49.17
											Sheen Observed: Y N
											DNAPL Observed: Y N
											Did Well Go Dry: Y N
								ļ			Other:
											1 Well Volume = 6.3 gal.

						Historic Info	ormotion				
Boring Log A	vailable (1	res/no/attac	hod):			HISTORIC INIC	ormation				
	,,,		•								
nstallation Lo	og Avallat	ne ( <b>yes</b> /no/a	allached)			C					
Summary  Monitoring Well: MWN-03B Ground Surface Elevation: 609.57 Riser/Screen Material: PVC											
Monitoring Well : MWN-03B Installation Date: 11/5/92					rrace Elevation er Elevation:	572.65					
	ale.	11/5/92							creen Depth		
nstalled By:		Turnkey		Elevation D	Point Elevation	612.29		ם וווטווו ט	f Screen De	pin. 70.72	
Provious Fiel	ld maacur	omont Inform	mation Availab								
Tevious Fiel	iu measur	ement mion	Hation Availat	ne (yes/110/		o of Drovious C	iold Mosou	romonto			
Danth to	11/2422	Ι .	-11	O:f: - /		s of Previous F			.l. ! .l!4		O a la r
Depth to			pH	-	Conductance	Tempera			bidity		Color
(ft)		•	ard Units)		S/cm)	( °C)		,	ITU)		Olean
39.4			- 7.80		3 - 3.139	13.2 - 1			- 38.04		Clear
Notes:	**Monso	on pump un	nable to purg	e water. 3 v	well volumes p	ourged with de	dicated ba	ailer.			
					· · · · ·				l D	-1	
	.,	014	FIE	eld Observa	tions					eters +/-	Sampling Information
xterior Obse	ervations:	OK							pH	+/- 0.1	Sample ID: <b>MWN-03B</b> -091522
- t - :: - : O		OV							Conductivit		Sample Time: 9:45
nterior Obse	ervations	OK							Turbidity	mperature +/- 10% # of Sample Containers: One rbidity +/- 10% Duplicate Sample ID: NA	
		-							ORP		Duplicate Sample ID: NA
Signs of Don	oogo/Tom		None					DO		Sample Analysis: Arsenic, Barium Chromium, Manganese	
<u> </u>				Surface Seal Intact ( <b>yes</b> /no) PID Measurement:					טטו	T/- 10 /0	
Locked (y			n (voc/no)	Surfa	aca Saal Intact	(vac/na)	DID Massi	iromont.	ND		Chromiani, Manganese
	(es/ <b>no</b> )		p ( <b>yes</b> /no)	Surfa	ace Seal Intact	\ <b>3</b>		urement:	ND	Odors:	pomonium, wanganese
	res/ <b>no</b> )		p ( <b>yes</b> /no)	Surfa	ace Seal Intact	( <b>yes</b> /no) Well Quali		urement:	ND		Johnston and Managariese
Date		Well Ca				Well Qual	ity Data			Odors:	
Date	Time	Well Ca	Cumulative	рН	Specific	Well Quali	ity Data Turbidity	Color	Dissolved	Odors: Oxygen	Notes
Date		Well Ca  Depth to  Water	Cumulative Volume	pH (Standard	Specific Conductance	Well Qual	ity Data		Dissolved Oxygen	Odors: Oxygen Reduction	
	Time	Well Ca  Depth to  Water ft bgs	Cumulative Volume Purged (Gal)	pH (Standard Units)	Specific Conductance (mS/cm)	Well Quali Temperature (°C)	Turbidity (NTU)	Color	Dissolved Oxygen %	Odors: Oxygen Reduction Potential	Notes
Date 9/15/2022		Well Ca  Depth to  Water	Cumulative Volume	pH (Standard	Specific Conductance	Well Quali	ity Data Turbidity		Dissolved Oxygen	Odors: Oxygen Reduction	Notes  Depth of Water: 39.64
	Time	Well Ca  Depth to  Water ft bgs	Cumulative Volume Purged (Gal)	pH (Standard Units)	Specific Conductance (mS/cm)	Well Quali Temperature (°C)	Turbidity (NTU)	Color	Dissolved Oxygen %	Odors: Oxygen Reduction Potential	Notes  Depth of Water: 39.64  Length of Water Column: 31.08
	Time	Well Ca  Depth to  Water ft bgs	Cumulative Volume Purged (Gal)	pH (Standard Units)	Specific Conductance (mS/cm)	Well Quali Temperature (°C)	Turbidity (NTU)	Color	Dissolved Oxygen %	Odors: Oxygen Reduction Potential	Notes  Depth of Water: 39.64  Length of Water Column: 31.08  Depth of Well: 70.72
	Time	Well Ca  Depth to  Water ft bgs	Cumulative Volume Purged (Gal)	pH (Standard Units)	Specific Conductance (mS/cm)	Well Quali Temperature (°C)	Turbidity (NTU)	Color	Dissolved Oxygen %	Odors: Oxygen Reduction Potential	Notes  Depth of Water: 39.64  Length of Water Column: 31.08  Depth of Well: 70.72
	Time	Well Ca  Depth to  Water ft bgs	Cumulative Volume Purged (Gal)	pH (Standard Units)	Specific Conductance (mS/cm)	Well Quali Temperature (°C)	Turbidity (NTU)	Color	Dissolved Oxygen %	Odors: Oxygen Reduction Potential	Notes  Depth of Water: 39.64  Length of Water Column: 31.08  Depth of Well: 70.72  Sheen Observed: Y N
	Time	Well Ca  Depth to  Water ft bgs	Cumulative Volume Purged (Gal)	pH (Standard Units)	Specific Conductance (mS/cm)	Well Quali Temperature (°C)	Turbidity (NTU)	Color	Dissolved Oxygen %	Odors: Oxygen Reduction Potential	Notes  Depth of Water: 39.64  Length of Water Column: 31.08  Depth of Well: 70.72  Sheen Observed: Y N  DNAPL Observed: Y N
	Time	Well Ca  Depth to  Water ft bgs	Cumulative Volume Purged (Gal)	pH (Standard Units)	Specific Conductance (mS/cm)	Well Quali Temperature (°C)	Turbidity (NTU)	Color	Dissolved Oxygen %	Odors: Oxygen Reduction Potential	Notes  Depth of Water: 39.64  Length of Water Column: 31.08  Depth of Well: 70.72  Sheen Observed: Y N  DNAPL Observed: Y N  Did Well Go Dry: Y N
	Time	Well Ca  Depth to  Water ft bgs	Cumulative Volume Purged (Gal)	pH (Standard Units)	Specific Conductance (mS/cm)	Well Quali Temperature (°C)	Turbidity (NTU)	Color	Dissolved Oxygen %	Odors: Oxygen Reduction Potential	Notes  Depth of Water: 39.64  Length of Water Column: 31.08  Depth of Well: 70.72  Sheen Observed: Y N  DNAPL Observed: Y N  Did Well Go Dry: Y N
	Time	Well Ca  Depth to  Water ft bgs	Cumulative Volume Purged (Gal)	pH (Standard Units)	Specific Conductance (mS/cm)	Well Quali Temperature (°C)	Turbidity (NTU)	Color	Dissolved Oxygen %	Odors: Oxygen Reduction Potential	Notes  Depth of Water: 39.64  Length of Water Column: 31.08  Depth of Well: 70.72  Sheen Observed: Y N  DNAPL Observed: Y N  Did Well Go Dry: Y N

						Historic Inf	ormation					
Boring Log A			•									
nstallation L	og Availat	ole ( <b>yes</b> /no/a	attached)									
		Summary										
Monitoring W		MWN-03D Ground Surface Elevation: 610.75 Riser/Screen Material: PVC										
nstallation D	ate:	7/29/94			er Elevation:	574.02			creen Depth			
nstalled By:		Turnkey	y Monitorin		Monitoring Point Elevation 613.			Bottom o	of Screen De	epth: 121.26		
				Elevation D								
Previous Fiel	ld measur	ement Inforr	mation Availal	ole (yes/ <b>no</b> /	attached)							
					Range	s of Previous F	ield Measu	rements				
Depth to	Water		рН	Specific (	Conductance	Tempera	ature	Tu	rbidity		Color	
(ft)		(Standa	ard Units)	(m	S/cm)	( °C)		(1)	NTU)			
41.7		6.17	' - 7.31	24.4	1 - 26.69	12.9 - 1		1.97	- 35.83		Clear	
Notes:	** At app	roximately	65' monsoon	pump una	ble to advance	e down well. 3	well volur	nes pura	ed with pur	mp and		
						with dedicate		p g	,	1. *****		
				eld Observa	-				Param	eters +/-	Sampling Information	
xterior Obs	ervations:	OK							рН	+/- 0.1	Sample ID: <b>MWN-03D</b> -091522	
									Conductivi		Sample Time: 8:30	
nterior Obse	ervations	OK								•	# of Sample Containers: Six	
											Duplicate Sample ID: NA	
									ORP	+/- 10mV	Sample Analysis: Barium, Manganese	
Signs of Dan	nage/Tam	pering:	** See note a	bove					DO	+/- 10%	VOC STARS List via EPA 8260C	
Locked (y	/es/ <b>no</b> )	Well Ca	p ( <b>yes</b> /no)	Surfa	ace Seal Intact	(yes/no)	PID Measi	urement:	ND	Odors:	SVOC B/N Via EPA 8270D, Barium. Manganese	
						Well Qual	ity Data					
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	1	Color	Dissolved	Oxygen	Notes	
		Water	Volume	`	Conductance	(°C)	(NTU)		Oxygen	Reduction		
		ft bgs	Purged (Gal)	Units)	(mS/cm)				%	Potential		
9/15/2022	8:30	41.60	39**	7.26	26.11	13.5	165.2	Brown	16.2	50.8	Depth of Water: 39.49	
											Length of Water Column: 81.77	
											Depth of Well: 121.26	
											Sheen Observed: Y N	
											DNAPL Observed: Y N	
											Did Well Go Dry: Y N	
											Other:	
											1 Well Volume = 13.0 gal.	

						Historic Inf	ormation				
Boring Log A	vailable (	vos/no/attac	hod):			HISTORIC IIII	omation				
		•	,								
Installation Lo	og Avallat	ole (yes/no/a	allached)			0					
B. 4 '( ' \ A	/ - II	Summary  MWN-04 Ground Surface Elevation: 621.02 Riser/Screen Material: PVC									
Monitoring Well : MWN-04											
Installation Date: 9/12/90					er Elevation:				creen Depth		
nstalled By:		Turnkey			Point Elevation	623.45		Bottom o	of Screen De	eptn: 58.53	
D		( ] . (	(' A'1-1	Elevation D							
Previous Fiel	a measur	ement infori	mation Availal	oie (yes/ <b>no</b> /	,						
						s of Previous F					
Depth to		1	рН	•	Conductance	Tempera			rbidity		Color
(ft)		,	ard Units)	,	S/cm)	( °C)		,	NTU)		
50.6	1	11.57	' - 12.05	2.31	1 - 2.72	15.7 - 1	16.7	1.98	8 - 2.6		Clear
Notes:	** Monso	on pump u	nable to purç	ge water. 6	gallons purge	d to dry.		·			
					<u> </u>						
			Fie	eld Observa	tions				Param	eters +/-	Sampling Information
xterior Obse	ervations:	OK							рН	+/- 0.1	Sample ID: <b>MWN-04</b> -091522
									Conductivi	ty +/- 3%	Sample Time: 14:50
nterior Obse	rvations	OK							Temperatu		# of Sample Containers: Five
									Turbidity	+/- 10%	Duplicate Sample ID: NA
									ORP		Sample Analysis:
Signs of Dam	nage/Tam								DO	+/- 10%	VOC STARS List via EPA 8260C
Locked (y	es/ <b>no</b> )	Well Ca	p ( <b>yes</b> /no)	Surfa	ace Seal Intact	73				Odors:	SVOC B/N Via EPA 8270D
						Well Qual	ity Data				
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	,	Color	Dissolved	Oxygen	Notes
		Water	Volume	(Standard	Conductance	( °C)	(NTU)		Oxygen	Reduction	
		ft bgs	Purged	Units)	(mS/cm)				%	Potential	
9/15/2022	14:50	56.12	6**	11.35	3.54	17.3	33.47	Brown	69.6	35.4	Depth of Water: 51.36
											Length of Water Column: 7.17
											Depth of Well: 58.53
											Sheen Observed: Y N
											DNAPL Observed: Y N
											Did Well Go Dry: Y N
											Other:
											1 Well Volume = 4.7 gal.
								ļ			



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