Groundwater Monitoring Report September 2025 Quarterly Monitoring Event

113 7th North Street Site
113 7th North Street, Liverpool, New York
Site No. C734152

Submitted to:
New York State Department of Environmental Conservation
Division of Environmental Remediation
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Submitted by:
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February 17, 2026
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1. Background and Site Description

1.1. Introduction

This report has been prepared on UniFirst Corporation's (Applicant's) behalf by GEI Consultants, Inc. (GEI) for the 113 7th North Street Site in Liverpool, New York (the Site; Figure 1). The Site is currently being managed as Site No. C734152 under a Brownfield Cleanup Agreement (BCA) executed on October 12, 2021. The BCA is being administered by the New York State Department of Environmental Conservation (NYSDEC) through the Brownfield Cleanup Program (BCP).

This report summarizes results from the quarterly groundwater monitoring event conducted September 23 through 25, 2025. Monitoring was performed in accordance with the Interim Site Management Plan (ISMP; GEI, 2025c) that was submitted to the NYSDEC on September 30, 2025.[1]

Additional samples were collected in December 2025 from three monitoring wells—installed by a third party on the Applicant's property and in the 7th North Street Right-of-Way—as part of the neighboring property spill investigation, outside the ISMP scope. The NYSDEC project manager approved sampling the additional wells (MW1C, MW1D, and MW1E) by email on June 9, 2025.

1.2. Background

A Remedial Investigation (RI), as well as additional environmental investigation activities in support of facility renovation/construction and two associated Interim Remedial Measures (IRMs) were conducted from 2020 through 2023. Environmental investigations identified an area of impacted soil and groundwater in the immediate vicinity of a former wastewater vault (Figure 2) beneath the former main facility building (the "source area"). The investigations identified the primary constituents of concern as chlorinated volatile organic compounds (CVOCs; tetrachloroethylene [PCE] and its degradation products).

The construction-related IRMs were implemented between 2021 and 2023 in conjunction with facility renovations (GEI, 2025a&b). The renovations included construction of a portion of a new expanded operations building, transition of operations to the new building, demolition of the pre-existing facility building, and construction of the remainder of the expanded facility over the footprint of the prior facility building. The IRMs generally included injection of treatment reagents (zero-valent iron [ZVI] and nutrients) into the source area and three adjacent permeable reactive barriers (PRBs) in May 2022 to address an area of soil and groundwater impacted by PCE (Figure 2), and installation of a sub-slab vapor mitigation system during construction of the new facility buildings. Due to the nature of the facility renovation and related construction schedule, the IRMs did not include source removal other than

[1] The ISMP was first submitted to NYSDEC and the Department of Health (NYSDOH) as draft in August 2024, as an attachment to the Construction Completion Report for Phase 2 Construction. The agencies subsequently provided minor comments in letters dated April 21 and May 21, 2025, respectively, including a request to make the ISMP a stand-alone document. The agencies' comments were addressed in the revised final ISMP submitted on September 30, 2025. (This final ISMP has not yet been approved by the agencies in writing.)

material (soil and groundwater) removed for the purposes of constructing the new facility. Instead, the IRMs focused on in-situ remedial measures intended to promote the attenuation of the CVOCs and to mitigate employees' potential exposure via the vapor intrusion pathway.

The groundwater monitoring component of the ISMP was developed to monitor the continued attenuation of CVOCs in groundwater within and downgradient from the Site after the IRMs were completed and until such time that a comprehensive Site Management Plan is prepared, as may be required in the future.

2. Quarterly Field Monitoring Activities

Quarterly field monitoring was conducted in September 2025, the monitoring well locations are on Figure 2.

2.1. Groundwater Elevation Gauging

The ISMP specifies groundwater-level gauging of 19 monitoring wells during each event to assess the direction of groundwater flow. Gauging was completed on September 23, 2025, at 22 wells (i.e., the 19 wells that are part of the ISMP, plus three third-party wells installed near MW1R in May 2025).

Table 1 and Figure 3 present the water-level data. As documented in Table 1, the highest groundwater elevation measurement among the water table wells was in the southern portion of the Site at MW3S (371.15 feet NAVD88). The lowest water table elevation measurement was in the northeastern-most monitoring well MW13S (366.89 feet NAVD88). The overall change in the elevation of the water table was 4.26 feet, which represents a hydraulic gradient of approximately 0.013 ft/ft.

The estimated groundwater flow (Figure 3) is inferred to be generally toward the northeast with certain components of flow toward the north. This is generally consistent with the December 2024 through June 2025 groundwater monitoring events performed following building completion.

2.2. Groundwater Sampling and Analysis

Groundwater samples were collected from 16 monitoring wells (13 wells included in the ISMP and three wells installed near MW1R (MW1C, MW1D and MW1E), associated with the petroleum-related spill at the Pilot property). Groundwater sampling was performed on September 23 through September 25, 2025, for the parameters in Table 2. Low flow groundwater samples were collected from 15 of the 16 monitoring wells. A sample was collected via volumetric purge from MW1R due to the persistence of non-Site-related NAPL in the well.

Samples from all 16 sampled wells were analyzed for volatile organic compounds (VOCs) using USEPA Method 8260D. In addition, depending on location, certain wells were also sampled for per- and polyfluoroalkyl substances (PFAS) by USEPA Method 1633 and monitored natural attenuation (MNA) parameters (including dissolved and total iron by USEPA Method 6020B, dissolved gases, and microbial populations using Microbial Insights' QuantArray-Chlor methodology), per the ISMP.

Groundwater samples were submitted for analysis to Eurofins Environment Testing of Buffalo, New York (Eurofins) and Microbial Insights of Knoxville Tennessee (MI; QuantArray-Chlor analysis only). Eurofins is certified to perform environmental analyses under the New York State Department of Health (NYSDOH) Environmental Laboratory Approval Program (ELAP). MI is a specialized laboratory that quantifies bacteria and functional genes to evaluate the potential for biodegradation of a spectrum of common CVOCs through anaerobic and aerobic (co)metabolic pathways.

Sample collection methods, analytical methods and the quality control measures performed during the September 2025 event were consistent with the Field Sampling Plan (FSP) and Quality Assurance Project Plan (QAPP) included in the ISMP, as well as the specifications provided by the NYSDEC Analytical Services Protocol (ASP).

2.3. Waste Management

Three 55-gallon drums of purge water generated during low-flow groundwater sampling in June 2025 and September 2025 were removed from the Site on November 20, 2025. The transporter and transport vehicles were compliant with 6 NYCRR, Chapter IV, Part 364 and the waste was properly disposed of at a permitted facility. Waste documentation is included in Appendix A.

3. Analytical Results

The analytical data, including VOCs, PFAS, and natural attenuation parameters except for the microbial populations, are tabulated in Table 3 and the associated laboratory data packages are included in Appendix B. In addition, the historical data associated with these sampling locations is summarized in Tables 4A, 4B, and 4C. The microbial population data report is in Appendix C.

VOC and PFAS groundwater analytical results were compared to the NYS Ambient Water Quality Standards (AWQS) or guidance values provided in the document titled "NYSDEC Division of Water Technical and Operational Guidance Series (TOGS) Number 1.1.1" (NYSDEC, 1998 with appendices through 2023).

The natural attenuation parameter results are not intended for comparison to standards, but for evaluation along with the VOC data to determine to what extent conditions favorable to natural attenuation are present and whether natural attenuation is occurring, as further discussed in the subsections below.

3.1. VOCs

The results of the September 2025 VOC analyses are summarized in Table 3 and historical data is summarized in Table 4A. Figure 4A presents the historical analytical data for each monitoring well. Figure 5A depicts a plume map of the total CVOC data, focusing on September 2025 quarterly monitoring results.[2] CVOC concentration trend graphs are in Appendix D.

3.1.1. September 2025 Data Discussion

A total of 16 wells (MW1R through MW8, MW11, MW12S, and three non-Applicant related wells [MW1C, MW1D and MW1E]) were analyzed for VOCs during the September 2025 event. The total CVOC concentrations detected at these wells ranged from non-detect (several perimeter wells) to 280,000 µg/L (source area MW3S).

As shown in Table 3, VOCs were detected above NYSDEC AWQS in September 2025 at eight of the sampled wells (MW2R, MW3S, MW4S, MW4I, MW1D, MW1E, MW7, MW5S):

- Four wells in the vicinity of the PCE source area (MW2R, MW3S, MW4S and MW4I) exceeded AWQS for Site-related CVOCs. While above AWQS, concentrations of COCs in these wells have generally decreased relative to prior sampling events.
- CVOCs were detected at concentrations above AWQS in MW1D and MW1E, downgradient from the source area, despite no detections above reporting limits of CVOCs in MW1R or MW1C.

[2] For the purposes of this report, Site-related CVOCs include tetrachloroethene (PCE), trichloroethene (TCE), cis 1,2 dichloroethene (cis 1,2 DCE), trans-1,2-dichloroethene (trans-1,2-DCE), and vinyl chloride.

- Vinyl chloride was detected above the AWQS in cross-gradient well MW7 adjacent to Luther Avenue. Vinyl chloride has consistently been greater than the AWQS since November 2023 but has been relatively stable in MW7.
- Several non-Site-related VOCs (benzene, toluene, ethylbenzene, xylenes (BTEX), isopropylbenzene, and acetone) were detected above AWQS in MW5S an upgradient well along Luther Avenue. These VOCs have consistently been detected in MW5S and are likely associated with an unrelated source. Further details regarding BTEX concentrations in MW5S are discussed in Section 3.5.

3.1.2. Trend Discussion

Trend monitoring graphs are in Appendix D-1 for the wells sampled as part of the ISMP program where CVOCs have been detected (MW-1/1R, MW2/2R, MW3S/TW6, MW3I, MW4S/TW9, MW4I, MW7, MW12S). As depicted in Appendix D-1, overall, evidence of reductive dechlorination of CVOCs is clear. Data for MW3S, which is located within the source area where remedial injections were performed in May 2022, show decreasing PCE and TCE concentrations and increasing intermediate breakdown product concentrations since the injections were performed. In general, CVOCs in the remaining source area/source adjacent wells[3] also continue to degrade.

The most downgradient well sampled as part of the quarterly program, MW12S, generally exhibits a stable trend for PCE and TCE, and a decreasing trend for cis-1,2-DCE and vinyl chloride (Appendix D-1), with concentrations of CVOCs consistently below AWQS in all samples collected in 2025.

CVOCs are generally not detected above laboratory reporting limits or are present at very low concentrations in upgradient wells MW5S and MW5I, and cross/down-gradient monitoring wells MW6, MW7, MW8 and MW11. Except for vinyl chloride in MW7, September 2025 CVOC concentrations are below AWQS. While vinyl chloride is above AWQS in MW7, it has been stable for the past four quarterly rounds of sampling and remains far lower than in source area wells, suggesting that dechlorination continues to occur in the source area and that the low levels of degraded CVOCs are generally delineated by these wells.

CVOCs have been detected above AWQS in MW1D and MW1E, downgradient from the source area, despite no detections above reporting limits of CVOCs in MW1R or MW1C, immediately upgradient. (These wells were installed as part of the petroleum-related spill investigation at the Pilot property.) For these wells, no apparent trend can be discerned at this time because they have only been sampled

[3] Note that the shallow groundwater data reported together for several monitoring locations in this area were unavoidably collected from a variety of wells and/or methods. This has likely resulted in variations in data that may not have existed if the sampling sources and methods were consistent. Specifically, for both the "MW1/1R" and "MW2/2R" locations, the February 2022 pre-remediation samples were collected from permanent 2-inch wells (MW1 and MW2) that were later decommissioned to allow for construction, while the February 2023 post-injection samples were collected during construction are represented by grab water samples from hydro punch samplers (MW1_G and MW2_G), and finally data generated from samples collected from July 2023 through present are from the 2-inch permanent replacement wells (MW1R and MW2R) that were installed in June 2023. A similar situation exists at MW3S and MW4S, except that the pre-remediation wells were temporary 1-inch well points (TW6 and TW9, respectively), not permanent wells like MW1 and MW2.

twice. Additional evaluation will be performed as more sampling events are conducted, which will help clarify any emerging trends, extent of the elevated CVOC contamination and inform future monitoring strategies.

VOCs will continue to be evaluated against applicable criteria and historical data during future monitoring events.

3.2. PFAS

The results of the September 2025 PFAS analyses are summarized in Table 3 and historical data is summarized in Table 4B. Figure 4B presents the historical analytical data for each monitoring well. Figures 4B and 5B present graphical representations of the PFOS and PFOA compounds at each well (an analytical data box figure of historical data and a plume map of the most recent concentrations, respectively). Trend graphs for PFOS, PFOA, and total PFAS concentrations are in Appendix D.

3.2.1. September 2025 Data Discussion

Six wells were analyzed for PFAS, including four in the source area (MW3S, MW3I, MW4S and MW4I) and two upgradient (MW5S and MW5I), during the September 2025 event. Detected PFOA concentrations ranged from non-detect (MW3I, MW4I, MW5I) to 24 ng/L (MW4S). The detected PFOS concentrations ranged from non-detect (MW3I, MW4I, MW5I) to 20 ng/L (MW5S).

As shown in Table 3, data were compared to NYSDEC AWQS, where available (i.e., for PFOS and PFOA). As shown in the table, PFOS and PFOA were above NYSDEC AWQS in samples from the three sampled shallow wells (MW3S, MW4S, and MW5S). Additional PFAS compounds were detected in groundwater but have not been assigned AWQS Guidance Values.

3.2.2. Trend Discussion

Trend monitoring graphs for PFOS, PFOA, and total PFAS concentrations are in Appendix D-2 for the sampled wells (MW3S/TW6, MW3I, MW4S, MW4I, MW5S, MW5I). These concentration ranges represent general stable or decreasing trends in PFOA and PFOS concentrations at these locations. Total PFAS concentrations have had generally decreasing trends in monitoring wells MW3S/I, MW4S/I, and MW5S, and a generally stable trend in MW5I.

Upgradient monitoring wells MW5S and MW5I contain PFAS likely attributable to an offsite source.

3.3. MNA Parameters

The results of September 2025 MNA parameter analyses are summarized in Table 3 and historical data is summarized in Table 4C. MNA parameters were analyzed in samples from monitoring wells within and surrounding the source area (MW2R, MW3S, MW3I, MW4S and MW4I). Analyses included dissolved iron, total iron, sulfate, and dissolved gases (ethane, ethene, methane). Data for these parameters have been collected as part of ongoing monitoring following the IRM injections for the purpose of establishing

geochemical trends indicative of MNA and its effects on CVOC degradation. Trend graphs for MNA parameters are in Appendix D-3.

3.3.1. Iron

When ZVI is injected, it readily oxidizes and releases Fe^{2+} ions into groundwater, resulting in a noticeable increase in dissolved iron levels (Lawrinenko, et. al., 2023). The primary purpose of the injected iron is to provide an electron donor to reduce CVOCs (i.e., degrade higher chlorinated VOCs to lower chlorinated VOCs through the process of chemical reduction). Further, under typical natural attenuation conditions, decreasing total iron concentrations, coupled with increasing dissolved iron concentrations suggest anaerobic dechlorination is occurring.

In September 2025, dissolved and total iron were present in all sampled wells, with highest concentrations in MW3S, as expected.[4] Post-injection concentrations of dissolved iron in MW3S increased throughout 2023, and appeared to have leveled off throughout 2024 and 2025.

While concentrations fluctuate, dissolved iron is consistently present in the five sampled wells, while total iron has decreased significantly relative to starting conditions, suggesting that degradation of CVOCs is occurring via chemical reduction as the result of the ZVI injections and natural attenuation, particularly in and immediately downgradient from the source area.

3.3.2. Sulfate

Depleted concentrations of sulfate relative to background/baseline indicate that conditions conducive to anaerobic dechlorination via sulfate reduction are present. Sulfate was detected at concentrations lower than baseline in the three shallow monitoring wells sampled in September 2025, and generally consistent with baseline in the two intermediate wells. Overall, sulfate trends were decreasing (in all but MW4I) during the first few sampling events after injections, but concentrations have leveled out since 2023. Because the reduction in sulfate is indicative of anaerobic degradation, this leveling-off of concentrations may suggest that biological degradation through sulfate reduction is slowing.

3.3.3. Dissolved Gases

Elevated concentrations of methane, as well as ethane and ethene, suggest that conditions are conducive to anaerobic degradation.[5] Methane was detected in most monitoring wells sampled, with overall increasing or steadily elevated trends (particularly within the source area, especially injection zone well MW3S) from 2023 through 2024, before temporarily falling through early 2025, then increasing again in September 2025. This suggests conditions that support reductive dechlorination persist, particularly in the source zone.

[4] MW3S is (the only well) located within the source area injection zone and has displayed the greatest rates of CVOC degradation to date among sampled wells.

[5] Although higher levels of ethane and ethene may also imply ongoing anaerobic dechlorination, methane typically serves as the principal indicator because ethene and ethane are subject to rapid degradation; thus, their absence does not necessarily indicate that CVOC biodegradation is not taking place.

3.4. QuantArray-Chlor Data

The results of the September 2025 QuantArray-Chlor analyses are included in Appendix C. The QuantArray-Chlor data continue to show that mid to high microbial populations remain present in MW3S (i.e., in the source injection zone), while low to mid microbial populations remain present in the other analyzed wells (MW3I, MW4S, and MW4I, immediately adjacent to the source injection zone).

The remainder of this summary focuses specifically on dehalococcoides (DHC), as this is the only known bacterial group capable of complete reductive dechlorination of PCE and TCE to ethene and therefore provides the most relevant/representative data with respect to natural attenuation conditions at the Site.

DHC was detected at a range of concentrations in the four analyzed wells (MW3S, MW3I, MW4S, and MW4I). Consistent with prior quarterly sampling events, the highest DHC concentration was observed at MW3S, which is within the source injection zone, at 3,100,000 cells/mL. Concentrations at other wells were lower (MW3I at 480 cells/mL, MW4S at 100 cells/mL, and MW4I at 2,800 cells/mL). Concentrations above 10,000 cells/ml are generally indicative of increased biological degradation potential. The DHC concentration in MW3S increased by approximately 1,900,000 cells/mL between June 2025 and September 2025. Sustained high levels of DHC in MW3S are ideal for degradation of PCE and TCE in the source zone. The lower population densities of DHC in the other wells are consistent with results of other analyses that suggest degradation continues to occur outside the injection zone, but at a lower rate than within the source area.

3.5. BTEX Detections

BTEX and/or petroleum-related non-aqueous phase liquid (NAPL) persist in two wells (upgradient well MW5S and down/cross gradient well MW1R). However, there is no evidence that their presence is attributable to a site-related source. Review of historical and current site information indicates that the facility is and was historically heated using natural gas, and petroleum products are not stored or used on the Site in significant quantities[6].

3.5.1. Off-Site Pilot Spill Response

In March 2024, because of the NYSDEC's Spill Response Team's tracing of a spill reported in Ley Creek and Onondaga Lake, petroleum NAPL was discovered in the new stormwater collection system located in the southeastern portion of the Property (Figure 2). The source of the spill was determined to be petroleum that migrated from the adjacent property, a Pilot fuel station, at 107 7th North Street. NYSDEC assigned NYSDEC Spill Number 2310035 and ultimately linked it to an existing Pilot spill (NYSDEC Spill Number 1511262).

[6] A pair of heating fuel oil underground storage tanks (USTs) were formerly located on a property located southwest of the Site, which the Applicant purchased as part of their facility renovation project. (This property is currently a parking area for the new facility.) The applicant properly removed these USTs and surrounding soil under NYSDEC Spill No. 2302335 as part of the facility renovation. The applicant received an No Further Action determination from NYSDEC on December 19, 2023.

Although Pilot had taken measures to mitigate the release, GEI continues to observe NAPL in MW1R.

GEI first notified the Applicant, Pilot, and the NYSDEC of this observation in December 2024. At that time, GEI initially bailed the product from the well, recovering approximately 500 milliliters of diesel-like material, which was sampled for petroleum hydrocarbon identification analyses. The laboratory analysis confirmed the NAPL to be diesel fuel.

Between December 2024 and present, Pilot's representatives have performed several monitoring and vacuum extraction events in MW1R. Post-vacuum extraction monitoring indicated little to no measurable NAPL in MW1R following the vacuum extraction events. However, NAPL continues to accumulate in the well over time, which Pilot purges periodically (most recently in September 2025).

Sovereign Consulting of Oxford Connecticut, on behalf of Pilot, installed five monitoring wells around MW1R in May 2025 to further characterize the spill conditions; one of which is on Applicant property (MW1C), and two of which are in the 7th North Street right-of-way (MW1D and MW1E). The remaining two (MW1A and MW1B) are on the adjacent 109 7th North Street property (Sposato Flooring Showroom). The five monitoring wells are four inches in diameter and screened approximately from 2 to 12 feet (same as MW1R). Pilot's representatives did not observe evidence of petroleum impacts during or after installation of these five wells and has not observed NAPL in these new wells to date. Similarly, GEI has not observed NAPL in MW1C, MW1D, or MW1E during any sampling event to date.

As water is purged from MW1R during quarterly sampling events, NAPL is drawn in from the surrounding formation. Depending on the NAPL conditions during each sampling event, GEI has employed different approaches during sampling, including low-flow sampling as planned per the GEI's standard operating procedure in March 2025, forgoing sampling in June 2025 (because NAPL could not be separated from groundwater), and using a volumetric purge technique (September 2025). Specific to the September 2025 event, approximately 4.75 gallons of NAPL and groundwater were purged from the well. To avoid fouling of the flow-through cell, field parameters were not measured with a water quality meter during purging. After approximately three well volumes of water were purged, the well was allowed to recharge for approximately 25 minutes before the sample was collected. The sample was collected via peristaltic pump, similar to prior sampling methods for this monitoring well.

3.5.2. Unknown Upgradient Source

Based on the inferred direction of groundwater flow, as well as the knowledge that gasoline is and was not historically stored on site, the impacts in MW5S are assumed to come from an upgradient off-site source in the vicinity of Luther Avenue. Sheen and fuel-like odor were observed in MW5S during the March 2025 sampling event but were not observed during the June 2025 event. Slight fuel-like odor was observed in MW5S during the September 2025 event, and sheen was not observed. Total BTEX concentrations in MW5S have been slightly decreasing from February 2023 through September 2025.

4. Summary

Quarterly field monitoring was conducted in September 2025, including gauging at 22 wells and sampling for VOCs, PFAS and MNA parameters at 16 wells, depending on location. Overall, the data support continued attenuation of CVOCs and continue to document the petroleum-related impacts from offsite source(s).

Analytical results indicate CVOC concentrations were detected above New York State AWQS in seven wells, mainly near the source area but most PCE concentrations are declining. Analytical results for MNA parameters and microbial populations support the declining trends finding, indicating that conditions conducive for ongoing anaerobic degradation and reductive dechlorination are present, especially in the source area.

CVOCs were again reported in MW1D and MW1E[7] at concentrations greater than surrounding wells.

The groundwater monitoring program will continue on a quarterly basis. In addition to the regular ISMP sampling program, GEI will continue to sample groundwater from Pilot wells MW1C, MW1D and MW1E and analyze for VOCs, similar to the June and September 2025 events.

[7] These are two of the five the new wells installed by others to further investigate/characterize the off-site petroleum impacts and then voluntarily sampled for VOCs in June 2025 and again in September 2025.

5. References

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Tables

Table 1. Groundwater Elevation Data
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

Well Location	Screened Interval	Reference Elevation	Groundwater Measurements	
			9/23/2025	
			DTW	Elevation
MW1R	4 - 14	374.01	4.38	369.63
MW1C	2 - 12	374.51	4.94	369.57
MW1D	2 - 12	373.26	4.62	368.64
MW1E	2 - 12	373.10	4.99	368.11
MW2R	4 - 14	376.48	5.70	370.78
MW3S	4 - 14	376.35	5.20	371.15
MW3I	26 - 36	376.24	5.50	370.74
MW4S	4 - 14	376.45	5.50	370.95
MW4I	26 - 36	376.32	6.25	370.07
MW5S	5 - 15	374.66	3.60	371.06
MW5I	34 - 44	374.66	4.19	370.47
MW6	8 - 18	374.06	4.84	369.22
MW7	8 - 18	375.26	6.18	369.08
MW8	5 - 15	375.64	6.89	368.75
MW9S	5 - 15	374.10	3.08	371.02
MW9I	26 - 36	374.21	3.85	370.36
MW10	8 - 18	373.72	2.80	370.92
MW11	6 - 16	373.50	6.29	367.21
MW12S	5 - 15	371.91	4.90	367.01
MW12I	25 - 35	371.98	2.23	369.75
MW13S	5 - 15	372.71	5.82	366.89
MW13I	20 - 30	372.59	5.5	367.09

Notes:
Screened interval reported in feet below top of well casing
DTW = depth to water (feet below measuring point)
Elevation = groundwater elevation in feet above mean sea level (NAVD83)

Table 2. Groundwater Sampling Scope of Work
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

Well ID	Screened Interval (ft bgs)	Water Level[1]	VOCs (Method 8260) Alpha	PFAS (Method 1633) Alpha	MNA Parameters (Various Methods[2]) Alpha	DHC Census (qPCR) Microbial Insights
MW1R	4 - 14	X	X	--	X	--
MW2R	4 - 14	X	X	--	X	--
MW3S	4 - 14	X	X	X	X	X
MW3I	26 - 36	X	X	X	X	X
MW4S	4 - 14	X	X	X	X	X
MW4I	26 - 36	X	X	X	X	X
MW5S	5 - 15	X	X	X	--	--
MW5I	34 - 44	X	X	X	--	--
MW6[4]	8 - 18	X	X	--	--	--
MW7	8 - 18	X	X	--	--	--
MW8[4]	5 - 15	X	X	--	--	--
MW9S[4]	5 - 15	X	--	--	--	--
MW9I[4]	26 - 36	X	--	--	--	--
MW10[4]	8 - 18	X	--	--	--	--
MW11	6 - 16	X	X	--	--	--
MW12S	5 - 15	X	X	--	--	--
MW12I	25 - 35	X	--	--	--	--
MW13S	5 - 15	X	--	--	--	--
MW13I	20 - 30	X	--	--	--	--
MW-1C[3]	*2 - 12*	*X*	*X*	--	--	--
MW-1D[3]	*2 - 12*	*X*	*X*	--	--	--
MW-1E[3]	*2 - 12*	*X*	*X*	--	--	--

NOTES:

1. A full round of water levels are performed at start of event, prior to sampling.
2. MNA parameters consist of dissolved and total iron by Method 6010B, ethene, ethane and methane by Method RSK-175, and sulfate by Method 121,4500SO4-E. The dissolved iron samples are field-filtered using a disposable 0.45 micron in-line filters.
3. MW1C, MW1D, and MW1E were installed by others in May 2025, associated with a 3rd party petroleum spill (NYSDEC Spill No. 2310035), and voluntarily sampled by UniFirst for VOCs, with NYSDEC project manager's (Michael Belveg's) approval via email dated June 9, 2025.

Table 3. Groundwater Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

Analyte	Units	CAS No.	AWQS	MW1C 9/25/2025		MW1D 9/25/2025		MW1E 9/25/2025		MW1R 9/24/2025		MW2R 9/24/2025		MW3S 9/24/2025		MW3I 9/24/2025		MW4S 9/24/2025		MW4I 9/24/2025		MW5S 9/24/2025		MW5I 9/24/2025		MW6 9/23/2025		MW7 9/23/2025		MW8 9/24/2025		MW11 9/24/2025		MW12S 9/24/2025	
Volatile Organic Compounds	ug/L																																
Acetone		67-64-1	50*	**5.6**	J	200	U	1000	U	50	U	10	U	20,000	U	10	U	20,000	U	400	U	**86**	J	10	U	10	U	10	U	**3.5**	J	10	U	10	U
Benzene		71-43-2	1	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	**18**		1	U	1	U	1	U	1	U	1	U	1	U
Bromodichloromethane		75-27-4	50*	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Bromoform		75-25-2	50*	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Bromomethane		74-83-9	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Carbon disulfide		75-15-0	60*	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Carbon tetrachloride		56-23-5	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Chlorobenzene		108-90-7	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Chloroethane		75-00-3	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Chloroform (Trichloromethane)		67-66-3	7	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Chloromethane		74-87-3	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Cyclohexane		110-82-7	NE	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	**370**		1	U	1	U	1	U	1	U	1	U	1	U
1,2-Dibromo-3-chloropropane		96-12-8	0.04	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Dibromochloromethane		124-48-1	50*	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
1,1-Dichloroethane		75-34-3	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
1,2-Dichloroethane		107-06-2	0.6	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
1,1-Dichloroethene		75-35-4	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	**36**	J	10	U	1	U	1	U	1	U	1	U	1	U	1	U
cis-1,2-Dichloroethene		156-59-2	5	1	U	**990**		**2,900**		5	U	**39**		**160,000**		**2.9**		**8,100**		**3,800**	F1	10	U	1	U	1	U	1	U	1	U	**2.3**	
trans-1,2-Dichloroethene		156-60-5	5	1	U	20	U	100	U	5	U	**1.7**		2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
1,2-Dichloropropane		78-87-5	1	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
cis-1,3-Dichloropropene		10061-01-5	0.4	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
trans-1,3-Dichloropropene		10061-02-6	0.4	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Ethylbenzene		100-41-4	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	**260**		1	U	1	U	1	U	1	U	1	U	1	U
2-Hexanone		591-78-6	50*	5	U	100	U	500	U	25	U	5	U	10,000	U	5	U	10,000	U	200	U	50	U	5	U	5	U	5	U	5	U	5	U	5	U
Isopropylbenzene		98-82-8	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	**110**		1	U	1	U	1	U	1	U	1	U	1	U
Methyl acetate		79-20-9	NE	2.5	U	50	U	250	U	13	U	2.5	U	5,000	U	2.5	U	5,000	U	100	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	10	U	200	U	1000	U	50	U	10	U	20,000	U	10	U	20,000	U	400	U	**34**	J	10	U	10	U	10	U	10	U	10	U	10	U
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	5	U	100	U	500	U	25	U	5	U	10,000	U	5	U	10,000	U	200	U	50	U	5	U	5	U	5	U	5	U	5	U	5	U
Methylcyclohexane		108-87-2	NE	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	**150**		1	U	1	U	1	U	1	U	1	U	1	U
Methylene chloride		75-09-2	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Styrene		100-42-5	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
1,1,2,2-Tetrachloroethane		79-34-5	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Tetrachloroethene (PCE)		127-18-4	5	1	U	**540**		**3,600**		5	U	**6.7**		2,000	U	1	U	**85,000**		**20**	J	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Toluene		108-88-3	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	**6.4**	J	1	U	1	U	1	U	1	U	1	U	1	U
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)		76-13-1	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
1,2,4-Trichlorobenzene		120-82-1	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
1,1,2-Trichloroethane		79-00-5	1	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Trichloroethene (TCE)		79-01-6	5	1	U	**210**		**1,100**		5	U	**4.4**		2,000	U	**0.53**	J	**4,100**		40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Trichlorofluoromethane (Freon 11)		75-69-4	5	1	U	20	U	100	U	5	U	1	U	2,000	U	1	U	2,000	U	40	U	10	U	1	U	1	U	1	U	1	U	1	U	1	U
Vinyl chloride		75-01-4	2	1	U	**28**		100	U	5	U	**3.9**		**120,000**		**8.4**		2,000	U	**550**		10	U	1	U	1	U	**13**		1	U	1	U	**1.5**	
Total Xylene		1330-20-7	5	2	U	40	U	200	U	**5.1**	J	2	U	4,000	U	2	U	4,000	U	80	U	**23**		2	U	2	U	2	U	2	U	2	U	2	U
Total BTEX		N/A	NE	ND		ND		ND		**5.1**		ND		ND		ND		ND		ND		**307.4**		ND		ND		ND		ND		ND	
Total VOCs		N/A	NE	**5.6**	J	**1,768**		**7,600**		**5.1**	J	**55.7**		**280,000**		**11.83**	J	**97,200**		**4,406**	J	**1,057.40**	J	ND		ND		**13**		**3.5**		ND		**3.8**	

Table 3. Groundwater Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

Analyte	Units	CAS No.	AWQS	MW1C 9/25/2025	MW1D 9/25/2025	MW1E 9/25/2025	MW1R 9/24/2025	MW2R 9/24/2025	MW3S 9/24/2025	MW3I 9/24/2025	MW4S 9/24/2025	MW4I 9/24/2025	MW5S 9/24/2025	MW5I 9/24/2025	MW6 9/23/2025	MW7 9/23/2025	MW8 9/24/2025	MW11 9/24/2025	MW12S 9/24/2025
PFAS	ng/L																		
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)		763051-92-9	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
2H,2H,3H,3H-Perfluorooctanoic Acid (5:3FTCA)		914637-49-3	NE	--	--	--	--	--	8.3 U	7.8 U	7.7 U	7.6 U	8.8 U	8.2 U	--	--	--	--	--
3-Perfluoroheptyl Propanoic Acid (7:3FTCA)		812-70-4	NE	--	--	--	--	--	8.3 U	7.8 U	7.7 U	7.6 U	8.8 U	8.2 U	--	--	--	--	--
3-Perfluoropropyl Propanoic Acid (3:3FTCA)		356-02-5	NE	--	--	--	--	--	3.3 U	3.1 U	3.1 U	3 U	3.5 U	3.3 U	--	--	--	--	--
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)		919005-14-4	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)		756426-58-1	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA)		4151-50-2	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
N-Ethyl Perfluorooctanesulfonamido Ethanol (NEtFOSE)		1691-99-2	NE	--	--	--	--	--	8.2 U	7.8 U	7.7 U	7.6 U	8.8 U	8.2 U	--	--	--	--	--
N-ethyl perfluorooctanesulfonamidoacetic acid (NEtFOSAA)		2991-50-6	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
N-Methyl Perfluorooctane Sulfonamide (NMeFOSA)		31506-32-8	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
N-Methyl Perfluorooctanesulfonamido Ethanol (NMeFOSE)		24448-09-7	NE	--	--	--	--	--	8.2 U	7.8 U	7.7 U	7.6 U	8.8 U	8.2 U	--	--	--	--	--
N-methyl perfluorooctanesulfonamidoacetic acid (NMeFOSAA)		2355-31-9	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)		151772-58-6	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA)		113507-82-7	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluoro(2-Propoxypropanoic) Acid		13252-13-6	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluoro-3-methoxypropanoic acid (PFMPA)		377-73-1	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluoro-4-methoxybutanoic acid (PFMBA)		863090-89-5	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluorobutanesulfonic acid (PFBS)		375-73-5	NE	--	--	--	--	--	**4.5**	1.6 U	**4.4**	1.5 U	**0.85 J**	1.6 U	--	--	--	--	--
Perfluorobutanoic acid (PFBA)		375-22-4	NE	--	--	--	--	--	**5.1**	**0.97 J**	**12**	**1.1 J**	**9.1**	**1.7 J**	--	--	--	--	--
Perfluorodecanesulfonic acid (PFDS)		335-77-3	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluorodecanoic acid (PFDA)		335-76-2	NE	--	--	--	--	--	1.7 U	1.6 U	**1.5**	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluorododecanesulfonic Acid (PFDoS)		79780-39-5	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluorododecanoic acid (PFDoA)		307-55-1	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluoroheptanesulfonic acid (PFHpS)		375-92-8	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluoroheptanoic acid (PFHpA)		375-85-9	NE	--	--	--	--	--	**5.7**	1.6 U	**11**	1.5 U	**2.6**	**1.5 J**	--	--	--	--	--
Perfluorohexanoic acid (PFHxA)		307-24-4	NE	--	--	--	--	--	**12**	1.6 U	**18**	1.5 U	**5**	**4.4**	--	--	--	--	--
Perfluorononanesulfonic Acid (Pfns)		68259-12-1	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluorooctanesulfonamide (FOSA)		754-91-6	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	**0.48 J**	1.6 U	--	--	--	--	--
Perfluoropentanesulfonic Acid (Pfpes)		2706-91-4	NE	--	--	--	--	--	**1.7**	1.6 U	**0.8 J**	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluoropentanoic Acid (PFPeA)		2706-90-3	NE	--	--	--	--	--	**7.3**	1.6 U	**25**	1.5 U	**8.1**	**4.9**	--	--	--	--	--
Perfluorotetradecanoic acid (PFTA/PFTeDA)		376-06-7	NE	--	--	--	--	--	1.6 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluorotridecanoic acid (PFTriA/PFTrDA)		72629-94-8	NE	--	--	--	--	--	1.6 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluoroundecanoic acid (PFUnA)		2058-94-8	NE	--	--	--	--	--	1.7 U	1.6 U	1.5 U	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Sodium 1H,1H,2H,2H-Perfluorodecane Sulfonate (8:2)		39108-34-4	NE	--	--	--	--	--	3.3 U	3.1 U	3.1 U	3 U	3.5 U	3.3 U	--	--	--	--	--
Sodium 1H,1H,2H,2H-Perfluorohexane Sulfonate (4:2)		757124-72-4	NE	--	--	--	--	--	3.3 U	3.1 U	3.1 U	3 U	3.5 U	3.3 U	--	--	--	--	--
Sodium 1H,1H,2H,2H-Perfluorooctane Sulfonate (6:2)		27619-97-2	NE	--	--	--	--	--	3.3 U	3.1 U	3.1 U	3 U	3.5 U	**39**	--	--	--	--	--
Perfluorohexanesulfonic acid (PFHxS)		355-46-4	NE	--	--	--	--	--	**3.4**	1.6 U	**3.7**	1.5 U	**1.4 J**	**1.1 J**	--	--	--	--	--
Perfluorononanoic Acid (PFNA)		375-95-1	NE	--	--	--	--	--	1.7 U	1.6 U	**4.1**	1.5 U	1.8 U	1.6 U	--	--	--	--	--
Perfluorooctanesulfonic acid (PFOS)		1763-23-1	2.7	--	--	--	--	--	**6.6 I**	1.6 U	**17**	1.5 U	**20**	1.6 U	--	--	--	--	--
Perfluorooctanoic Acid (PFOA)		335-67-1	6.7	--	--	--	--	--	**15**	1.6 U	**24**	1.5 U	**14**	1.6 U	--	--	--	--	--
Monitored Natural Attenuation Parameters	ug/L																		
Ethane		74-84-0	NE	--	--	--	**5.2 J**	7.5 U	**1,100**	7.5 U	**390**	**15**	--	--	--	--	--	--	--
Ethene		74-85-1	NE	--	--	--	7 U	7 U	**29,000**	7 U	**13**	**54**	--	--	--	--	--	--	--
Methane		74-82-8	NE	--	--	--	**1,800**	**890**	**6,800**	**140**	**4,600**	**1,400**	--	--	--	--	--	--	--
Sulfate		14808-79-8	250000	--	--	--	**30,600**	**27,100**	**25,000**	**79,900**	**131,000**	**128,000**	--	--	--	--	--	--	--
Iron, Dissolved		7439-89-6	300	--	--	--	**210**	**1,200**	**88,000**	**1000**	**81**	**1,300**	--	--	--	--	--	--	--
Iron, Total		7439-89-6	300	--	--	--	**610**	**1000**	**91,600**	**1,100**	**1,500**	**1,200**	--	--	--	--	--	--	--

Table 3. Groundwater Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

Notes:
Data for these sampling events have not been validated. Qualifiers are Lab Qualifiers.
Analytes in blue are not detected in any sample
ng/L = nanogram per liter (ppt)
ug/L = micrograms per liter or parts per billion (ppb)

BTEX = Benzene, Toluene, Ethylbenzene, and Xylenes
PFAS = Per- and polyfluoroalkyl substances
VOC = Volatile Organic Compound

Total VOCs are calculated using detects only.

AWQS = New York State Ambient Water Quality Standards and Guidance Values for GA groundwater
* indicates the value is a guidance value and not a standard

CAS No. = Chemical Abstracts Service Number
ND = Not Detected
NE = Not Established
N/A = Not Applicable

Bolding indicates a detected result concentration
Shading and bolding indicates that the detected concentration is above the NYSDOH guidance to which it was compared

Laboratory Qualifiers:
F1 = MS and/or MSD recovery exceeds the control limits
I = The lower value for the two columns has been reported due to obvious interference.
J = The result is an estimated value.
U = The result was not detected above the reporting limit.

Table 4A. Historical Groundwater VOC Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW1		MW1		MW1		MW1		MW1		MW1_G		MW1R		MW1R		MW1R		MW1R		MW1R	
			Sample Date	5/28/2020		3/21/2021		7/10/2021		11/22/2021		2/28/2022		2/7/2023		7/6/2023		11/30/2023		12/11/2024		3/12/2025		9/24/2025	
Analyte	**Units**	**CAS No.**	**NYS AWQS**																						
Volatile Organic Compounds	ug/L																								
Acetone		67-64-1	50*	**36**	**J**	5	U	5	U	5	UJ	5	U	5	U	5	U	5	U	**2.5**	**J**	25	U	50	U
Benzene		71-43-2	1	5	U	0.5	U	0.5	U	0.5	U	0.5	U	**0.18**	**J**	0.5	U	**0.62**		0.5	U	2.5	U	5	U
Bromochloromethane		74-97-5	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	--	
Bromodichloromethane		75-27-4	50*	5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	2.5	U	5	U
Bromoform		75-25-2	50*	20	U	2	U	2	U	2	U	2	U	2	U	2	UJ	2	U	2	U	10	U	5	U
Bromomethane		74-83-9	5	25	U	2.5	UJ	2.5	UJ	2.5	U	2.5	U	2.5	UJ	2.5	U	2.5	U	2.5	U	12	U	5	U
Carbon disulfide		75-15-0	60*	50	U	5	U	5	U	5	U	5	U	5	UJ	5	U	5	U	5	U	25	U	5	U
Carbon tetrachloride		56-23-5	5	5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	2.5	U	5	U
Chlorobenzene		108-90-7	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
Chloroethane		75-00-3	5	25	UJ	2.5	UJ	2.5	UJ	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
Chloroform (Trichloromethane)		67-66-3	7	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
Chloromethane		74-87-3	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	UJ	2.5	U	12	U	5	U
Cyclohexane		110-82-7	NE	100	U	10	U	10	U	10	UJ	10	U	10	U	10	U	10	U	10	U	50	U	5	U
1,2-Dibromo-3-chloropropane		96-12-8	0.04	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
Dibromochloromethane		124-48-1	50*	5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	UJ	0.5	U	0.5	U	2.5	U	5	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	20	U	2	U	2	U	2	U	2	U	2	U	2	U	2	U	2	U	10	U	5	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	50	U	5	U	5	UJ	5	U	5	U	5	U	5	U	5	U	5	U	25	U	5	U
1,1-Dichloroethane		75-34-3	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
1,2-Dichloroethane		107-06-2	0.6	5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	2.5	U	5	U
1,1-Dichloroethene		75-35-4	5	5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	**0.18**	**J**	0.5	U	0.5	U	2.5	U	5	U
cis-1,2-Dichloroethene		156-59-2	5	**80**		**17**		**20**		**29**		**24**	**J**	**37**		**130**		**14**		**0.78**	**J**	12	U	5	U
trans-1,2-Dichloroethene		156-60-5	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	**0.71**	**J**	2.5	U	2.5	U	12	U	5	U
1,2-Dichloropropane		78-87-5	1	10	U	1	U	1	U	1	U	1	U	1	U	1	U	1	U	1	U	5	U	5	U
cis-1,3-Dichloropropene		10061-01-5	0.4	5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	2.5	U	5	U
trans-1,3-Dichloropropene		10061-02-6	0.4	5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	UJ	0.5	U	0.5	U	0.5	U	2.5	U	5	U
1,4-Dioxane		123-91-1	NE	2,500	UJ	250	UJ	250	U	250	U	250	U	250	U	250	U	250	U	250	U	1,200	U	--	
Ethylbenzene		100-41-4	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	**15**		**14**		**7.2**	**J**	5	U
2-Hexanone		591-78-6	50*	50	U	5	U	5	U	5	U	5	U	5	U	5	U	5	U	5	U	25	U	25	U
Isopropylbenzene		98-82-8	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	**3.2**		**5.4**		12	U	5	U
Methyl acetate		79-20-9	NE	20	U	2	UJ	2	U	2	U	2	U	2	U	2	U	2	UJ	2	U	10	U	13	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	50	U	5	U	5	U	5	U	5	U	5	U	5	U	5	U	5	U	25	U	50	U
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	50	U	5	U	5	U	5	U	5	U	5	U	5	U	5	UJ	5	U	25	U	25	U
Methylcyclohexane		108-87-2	NE	100	U	10	U	10	U	10	U	10	U	10	U	**0.41**	**J**	**0.57**	**J**	**0.58**	**J**	50	U	5	U
Methylene chloride		75-09-2	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
Styrene		100-42-5	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
1,1,2,2-Tetrachloroethane		79-34-5	5	5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	2.5	U	5	U
Tetrachloroethene (PCE)		127-18-4	5	**1,400**		**29**		**8.7**		**22**		**25**		**52**		**91**		0.5	U	0.5	U	**1.5**	**J**	5	U
Tetrahydrofuran		109-99-9	50*	--		--		--		--		--		--		--		5	U	--		--		--	
Toluene		108-88-3	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113		76-13-1	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
1,2,3-Trichlorobenzene		87-61-6	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	--	
1,2,4-Trichlorobenzene		120-82-1	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
1,1,2-Trichloroethane		79-00-5	1	15	U	1.5	U	1.5	U	1.5	U	1.5	U	1.5	U	1.5	U	1.5	U	1.5	U	7.5	U	5	U
Trichloroethene (TCE)		79-01-6	5	**100**		**3.8**		**1.9**		**3.7**		**4**	**J**	**8.7**		**27**		**0.29**	**J**	0.5	U	2.5	U	5	U
Trichlorofluoromethane (Freon 11)		75-69-4	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	5	U
Vinyl chloride		75-01-4	2	**3.7**	**J**	**0.73**	**J**	**1.4**	**J**	**2**	**J**	**2**		**2**		**54**		**14**		**0.57**	**J**	**0.61**	**J**	5	U
o-Xylene		95-47-6	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	**0.92**	**J**	**35**		**13**		--	
m/p-Xylene		179601-23-1	5	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	**13**		12	U	--	
Total Xylene				25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	**0.92**	**J**	**48**		**13**		**5.1**	**J**
Total VOCs		--	NE	**1,619.70**	**J**	**50.53**	**J**	**32**	**J**	**56.7**	**J**	**55**	**J**	**99.88**	**J**	**303.3**	**J**	**48.6**	**J**	**71.83**	**J**	**22.31**	**J**	**5.1**	**J**

Table 4A. Historical Groundwater VOC Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW1C		MW1C		MW1D		MW1D		MW1E		MW1E	
			Sample Date	6/18/2025		9/25/2025		6/18/2025		9/25/2025		6/18/2025		9/25/2025	
Analyte	**Units**	**CAS No.**	**NYS AWQS**												
Volatile Organic Compounds	ug/L														
Acetone		67-64-1	50*	**5.9**		**5.6**	**J**	50	U	200	U	120	U	1000	U
Benzene		71-43-2	1	0.5	U	1	U	5	U	20	U	12	U	100	U
Bromochloromethane		74-97-5	5	2.5	U	--		25	U	--		62	U	--	
Bromodichloromethane		75-27-4	50*	0.5	U	1	U	5	U	20	U	12	U	100	U
Bromoform		75-25-2	50*	2	U	1	U	20	U	20	U	50	U	100	U
Bromomethane		74-83-9	5	2.5	U	1	U	25	U	20	U	62	U	100	U
Carbon disulfide		75-15-0	60*	5	U	1	U	50	U	20	U	120	U	100	U
Carbon tetrachloride		56-23-5	5	0.5	U	1	U	5	U	20	U	12	U	100	U
Chlorobenzene		108-90-7	5	2.5	U	1	U	25	U	20	U	62	U	100	U
Chloroethane		75-00-3	5	2.5	U	1	U	25	U	20	U	62	U	100	U
Chloroform (Trichloromethane)		67-66-3	7	2.5	U	1	U	25	U	20	U	62	U	100	U
Chloromethane		74-87-3	5	2.5	U	1	U	25	U	20	U	62	U	100	U
Cyclohexane		110-82-7	NE	10	U	1	U	100	U	20	U	250	U	100	U
1,2-Dibromo-3-chloropropane		96-12-8	0.04	2.5	U	1	U	25	U	20	U	62	U	100	U
Dibromochloromethane		124-48-1	50*	0.5	U	1	U	5	U	20	U	12	U	100	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	2	U	1	U	20	U	20	U	50	U	100	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	2.5	U	1	U	25	U	20	U	62	U	100	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	2.5	U	1	U	25	U	20	U	62	U	100	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	2.5	U	1	U	25	U	20	U	62	U	100	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	5	U	1	U	50	U	20	U	120	U	100	U
1,1-Dichloroethane		75-34-3	5	2.5	U	1	U	25	U	20	U	62	U	100	U
1,2-Dichloroethane		107-06-2	0.6	0.5	U	1	U	5	U	20	U	12	U	100	U
1,1-Dichloroethene		75-35-4	5	0.5	U	1	U	5	U	20	U	12	U	100	U
cis-1,2-Dichloroethene		156-59-2	5	2.5	U	1	U	**230**		**990**		**430**		**2,900**	
trans-1,2-Dichloroethene		156-60-5	5	2.5	U	1	U	25	U	20	U	62	U	100	U
1,2-Dichloropropane		78-87-5	1	1	U	1	U	10	U	20	U	25	U	100	U
cis-1,3-Dichloropropene		10061-01-5	0.4	0.5	U	1	U	5	U	20	U	12	U	100	U
trans-1,3-Dichloropropene		10061-02-6	0.4	0.5	U	1	U	5	U	20	U	12	U	100	U
1,4-Dioxane		123-91-1	NE	250	U	--		2,500	U	--		6,200	U	--	
Ethylbenzene		100-41-4	5	2.5	U	1	U	25	U	20	U	62	U	100	U
2-Hexanone		591-78-6	50*	5	U	5	U	50	U	100	U	120	U	500	U
Isopropylbenzene		98-82-8	5	2.5	U	1	U	25	U	20	U	62	U	100	U
Methyl acetate		79-20-9	NE	2	U	2.5	U	20	U	50	U	50	U	250	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	5	U	10	U	50	U	200	U	120	U	1000	U
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	2.5	U	1	U	25	U	20	U	62	U	100	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	5	U	5	U	50	U	100	U	120	U	500	U
Methylcyclohexane		108-87-2	NE	**1.1**	**J**	1	U	**12**	**J**	20	U	250	U	100	U
Methylene chloride		75-09-2	5	2.5	U	1	U	25	U	20	U	62	U	100	U
Styrene		100-42-5	5	2.5	U	1	U	25	U	20	U	62	U	100	U
1,1,2,2-Tetrachloroethane		79-34-5	5	0.5	U	1	U	5	U	20	U	12	U	100	U
Tetrachloroethene (PCE)		127-18-4	5	**1.1**		1	U	**1000**		**540**		**3,600**		**3,600**	
Tetrahydrofuran		109-99-9	50*	--		--		--		**--**		--		**--**	
Toluene		108-88-3	5	2.5	U	1	U	25	U	20	U	62	U	100	U
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113		76-13-1	5	2.5	U	1	U	25	U	20	U	62	U	100	U
1,2,3-Trichlorobenzene		87-61-6	5	2.5	U	--		1	U	--		20	U	--	
1,2,4-Trichlorobenzene		120-82-1	5	2.5	U	1	U	25	U	20	U	62	U	100	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	2.5	U	1	U	25	U	20	U	62	U	100	U
1,1,2-Trichloroethane		79-00-5	1	1.5	U	1	U	15	U	20	U	38	U	100	U
Trichloroethene (TCE)		79-01-6	5	0.5	U	1	U	**210**		**210**		**760**		**1,100**	
Trichlorofluoromethane (Freon 11)		75-69-4	5	2.5	U	1	U	25	U	20	U	62	U	100	U
Vinyl chloride		75-01-4	2	**0.13**	**J**	1	U	**16**		**28**		**0.64**	**J**	100	U
o-Xylene		95-47-6	5	2.5	U	--		25	U	**--**		62	U	--	
m/p-Xylene		179601-23-1	5	2.5	U	--		25	U	**--**		62	U	--	
Total Xylene				2.5	U	2	U	25	U	40	U	62	U	200	U
Total VOCs		--	NE	**8.23**	**J**	**5.6**	**J**	**1,468**	**J**	**1,768**		**4,802**	**J**	**7,600**	

Table 4A. Historical Groundwater VOC Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW2		MW2		MW2 (DUP)		MW2		MW2		MW2		MW2_G		MW2R		MW2R		MW2R		MW2R		MW2R		MW2R	
			Sample Date	5/28/2020		3/20/2021		3/20/2021		7/11/2021		11/22/2021		2/28/2022		2/7/2023		7/6/2023		12/1/2023		12/11/2024		3/12/2025		6/17/2025		9/24/2025	
Analyte	**Units**	**CAS No.**	**NYS AWQS**																										
Volatile Organic Compounds	ug/L																												
Acetone		67-64-1	50*	**20**		**17**		**17**		5	U	5	UJ	**6.8**		**16**	**J**	**7.8**		5	U	5	U	5	U	**6.4**		10	U
Benzene		71-43-2	1	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	2.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Bromochloromethane		74-97-5	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	--	
Bromodichloromethane		75-27-4	50*	0.5	U	**0.21**	**J**	**0.25**	**J**	**0.93**		**0.92**		**0.57**		2.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Bromoform		75-25-2	50*	2	U	2	U	2	U	2	U	2	U	2	U	10	U	2	U	2	U	2	U	2	U	2	U	1	U
Bromomethane		74-83-9	5	2.5	UJ	2.5	UJ	2.5	UJ	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Carbon disulfide		75-15-0	60*	5	U	5	U	5	U	5	U	5	U	5	U	25	U	5	U	5	U	5	U	5	U	5	U	1	U
Carbon tetrachloride		56-23-5	5	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	2.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Chlorobenzene		108-90-7	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloroethane		75-00-3	5	2.5	UJ	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloroform (Trichloromethane)		67-66-3	7	**3.1**		**36**		**36**		**28**		**8.3**		**4.3**		12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloromethane		74-87-3	5	2.5	UJ	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	UJ	2.5	U	2.5	U	2.5	U	1	U
Cyclohexane		110-82-7	NE	10	U	10	U	10	U	10	U	10	UJ	10	U	50	U	10	U	10	U	10	U	10	U	10	U	1	U
1,2-Dibromo-3-chloropropane		96-12-8	0.04	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Dibromochloromethane		124-48-1	50*	0.5	U	0.5	U	0.5	U	**0.24**	**J**	0.5	U	0.5	U	2.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	2	U	2	U	2	U	2	U	2	U	2	U	10	U	2	U	2	U	2	U	2	U	2	U	1	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	5	UJ	5	U	5	U	5	U	5	U	5	U	25	U	5	U	5	U	5	U	5	U	5	U	1	U
1,1-Dichloroethane		75-34-3	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2-Dichloroethane		107-06-2	0.6	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	2.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,1-Dichloroethene		75-35-4	5	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	2.5	U	**0.27**	**J**	**0.18**	**J**	0.5	U	0.5	U	0.5	U	1	U
cis-1,2-Dichloroethene		156-59-2	5	**3.4**		**8.5**		**7.8**		**0.92**	**J**	2.5	U	2.5	U	**130**		**38**		**51**		**28**		**21**		**37**		**39**	
trans-1,2-Dichloroethene		156-60-5	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	**0.85**	**J**	**1.7**	
1,2-Dichloropropane		78-87-5	1	1	U	1	U	1	U	1	U	1	U	1	U	5	U	1	U	1	U	1	U	1	U	1	U	1	U
cis-1,3-Dichloropropene		10061-01-5	0.4	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	2.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
trans-1,3-Dichloropropene		10061-02-6	0.4	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	2.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,4-Dioxane		123-91-1	NE	250	UJ	250	UJ	250	UJ	250	UJ	250	U	250	U	1,200	U	250	U	250	U	250	U	250	U	250	U	--	
Ethylbenzene		100-41-4	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
2-Hexanone		591-78-6	50*	5	U	5	U	5	U	5	U	5	U	5	U	25	U	5	U	5	U	5	U	5	U	5	U	5	U
Isopropylbenzene		98-82-8	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Methyl acetate		79-20-9	NE	2	U	2	UJ	2	UJ	2	U	2	U	2	U	10	U	2	U	2	UJ	2	U	2	U	2	U	2.5	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	5	U	5	U	5	U	5	U	5	U	5	U	25	U	5	U	5	U	**1.9**	**J**	5	U	5	U	10	U
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	5	U	5	U	5	U	5	U	5	U	5	U	25	U	5	U	5	UJ	5	U	5	U	5	U	5	U
Methylcyclohexane		108-87-2	NE	10	U	10	U	10	U	10	U	10	U	10	U	50	U	10	U	10	U	10	U	10	U	10	U	1	U
Methylene chloride		75-09-2	5	2.5	U	**0.73**	**J**	**0.73**	**J**	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Styrene		100-42-5	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2,2-Tetrachloroethane		79-34-5	5	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	2.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Tetrachloroethene (PCE)		127-18-4	5	**4**		**64**		**61**		**0.76**		**2.1**		**0.43**	**J**	**470**		**150**	**J**	**42**		**1.7**		**16**		**0.27**	**J**	**6.7**	
Tetrahydrofuran		109-99-9	50*	--		--		--		--		--		--		--		--		5	U	--		--		--		--	
Toluene		108-88-3	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	**530**		**9.8**		2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113		76-13-1	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2,3-Trichlorobenzene		87-61-6	5	62	U	2.5	U	2.5	U	2.5	U	2.5	UJ	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	--	
1,2,4-Trichlorobenzene		120-82-1	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2-Trichloroethane		79-00-5	1	1.5	U	1.5	U	1.5	U	1.5	U	1.5	U	1.5	U	7.5	U	1.5	U	1.5	U	1.5	U	1.5	U	1.5	U	1	U
Trichloroethene (TCE)		79-01-6	5	**0.91**		**7.1**		**6.7**		**0.34**	**J**	**0.35**	**J**	0.5	U	**120**		**18**		**17**		**2.8**		**7.2**		**3.7**		**4.4**	
Trichlorofluoromethane (Freon 11)		75-69-4	5	2.5	UJ	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride		75-01-4	2	**0.44**	**J**	**0.66**	**J**	**0.64**	**J**	1	U	**0.09**	**J**	1	U	**1**	**J**	**0.79**	**J**	**1.2**		**1.4**		**0.4**	**J**	**1.7**		**3.9**	
o-Xylene		95-47-6	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U
m/p-Xylene		179601-23-1	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	--	
Total Xylene				2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	--	
Total VOCs		--	NE	**31.85**	**J**	**134.2**	**J**	**130.12**	**J**	**31.19**	**J**	**11.76**	**J**	**12.1**	**J**	**1,267**	**J**	**224.66**	**J**	**111.38**	**J**	**35.8**		**44.6**	**J**	**49.92**	**J**	**55.7**	

Table 4A. Historical Groundwater VOC Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	TW6		TW6		TW6		TW6		TW6 (DUP)		MW3S_G		MW3S		MW3S		MW3S		MW3S		MW3S		MW3S	
			Sample Date	3/20/2021		7/11/2021		11/21/2021		2/27/2022		2/27/2022		2/7/2023		7/6/2023		12/1/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
Analyte	**Units**	**CAS No.**	**NYS AWQS**																								
Volatile Organic Compounds	ug/L																										
Acetone		67-64-1	50*	200	U	1000	U	1000	U	1,200	U	1,200	U	2,500	U	2,500	U	5,000	U	5,000	U	5,000	U	5,000	U	20,000	U
Benzene		71-43-2	1	20	U	100	U	100	U	120	U	120	U	250	U	250	U	500	U	500	U	500	U	500	U	2,000	U
Bromochloromethane		74-97-5	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U		
Bromodichloromethane		75-27-4	50*	20	U	100	U	100	U	120	U	120	U	250	U	250	U	500	U	500	U	500	U	500	U	2,000	U
Bromoform		75-25-2	50*	80	U	400	U	400	U	500	U	500	U	1000	U	1000	UJ	2,000	U	2,000	U	2,000	U	2,000	U	2,000	U
Bromomethane		74-83-9	5	100	UJ	500	UJ	500	U	620	U	620	UJ	1,200	UJ	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
Carbon disulfide		75-15-0	60*	200	U	1000	U	1000	U	1,200	U	1,200	U	2,500	UJ	2,500	U	5,000	U	5,000	U	5,000	U	5,000	U	2,000	U
Carbon tetrachloride		56-23-5	5	20	U	100	U	100	U	120	U	120	U	250	U	250	U	500	U	500	U	500	U	500	U	2,000	U
Chlorobenzene		108-90-7	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
Chloroethane		75-00-3	5	100	U	500	UJ	500	U	620	U	620	UJ	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
Chloroform (Trichloromethane)		67-66-3	7	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
Chloromethane		74-87-3	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	UJ	2,500	U	2,500	U	2,500	U	2,000	U
Cyclohexane		110-82-7	NE	400	U	2,000	U	2,000	U	2,500	U	2,500	U	5,000	U	5,000	U	10,000	U	10,000	U	10,000	U	10,000	U	2,000	U
1,2-Dibromo-3-chloropropane		96-12-8	0.04	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
Dibromochloromethane		124-48-1	50*	20	U	100	U	100	U	120	U	120	U	250	U	250	UJ	500	U	500	U	500	U	500	U	2,000	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	80	U	400	U	400	U	500	U	500	U	1000	U	1000	U	2,000	U	2,000	U	2,000	U	2,000	U	2,000	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	200	U	1000	UJ	1000	U	1,200	U	1,200	UJ	2,500	U	2,500	U	5,000	U	5,000	U	5,000	U	5,000	U	2,000	U
1,1-Dichloroethane		75-34-3	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
1,2-Dichloroethane		107-06-2	0.6	20	U	100	U	100	U	120	U	120	U	250	U	250	U	500	U	500	U	500	U	500	U	2,000	U
1,1-Dichloroethene		75-35-4	5	**40**		100	U	100	U	120	U	120	U	**100**	**J**	**200**	**J**	**220**	**J**	500	U	500	U	500	U	2,000	U
cis-1,2-Dichloroethene		156-59-2	5	**16,000**		**11,000**		**8,400**		**14,000**		**10,000**		**57,000**		**90,000**		**190,000**		**92,000**		**100,000**		**86,000**		**160,000**	
trans-1,2-Dichloroethene		156-60-5	5	**110**		**170**	**J**	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
1,2-Dichloropropane		78-87-5	1	40	U	200	U	200	U	250	U	250	U	500	U	500	U	1000	U	1000	U	1000	U	1000	U	2,000	U
cis-1,3-Dichloropropene		10061-01-5	0.4	20	U	100	U	100	U	120	U	120	U	250	U	250	U	500	U	500	U	500	U	500	U	2,000	U
trans-1,3-Dichloropropene		10061-02-6	0.4	20	U	100	U	100	U	120	U	120	U	250	UJ	250	U	500	U	500	U	500	U	500	U	2,000	U
1,4-Dioxane		123-91-1	NE	10,000	UJ	50,000	U	50,000	U	62,000	U	62,000	U	120,000	U	120,000	U	250,000	U	250,000	U	250,000	U	250,000	U	--	
Ethylbenzene		100-41-4	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
2-Hexanone		591-78-6	50*	200	U	1000	U	1000	U	1,200	U	1,200	U	2,500	U	2,500	U	5,000	U	5,000	U	5,000	U	5,000	U	10,000	U
Isopropylbenzene		98-82-8	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
Methyl acetate		79-20-9	NE	80	UJ	400	U	400	U	500	U	500	U	1000	U	1000	U	2,000	UJ	2,000	U	2,000	U	2,000	U	5,000	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	200	U	1000	U	1000	U	1,200	U	1,200	U	2,500	U	2,500	U	5,000	U	5,000	U	5,000	U	5,000	U	20,000	U
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	200	U	1000	U	1000	U	1,200	U	1,200	U	2,500	U	2,500	U	5,000	UJ	5,000	U	5,000	U	5,000	U	10,000	U
Methylcyclohexane		108-87-2	NE	400	U	2,000	U	2,000	U	2,500	U	2,500	U	5,000	U	5,000	U	10,000	U	10,000	U	10,000	U	10,000	U	2,000	U
Methylene chloride		75-09-2	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
Styrene		100-42-5	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
1,1,2,2-Tetrachloroethane		79-34-5	5	20	U	100	U	100	U	120	U	120	U	250	U	250	U	500	U	500	U	500	U	500	U	2,000	U
Tetrachloroethene (PCE)		127-18-4	5	**40,000**	**J**	**21,000**		**20,000**		**34,000**		**24,000**		**63,000**		**41,000**		**1,400**		500	U	500	U	500	U	2,000	U
Tetrahydrofuran		109-99-9	50*	--		--		--		--		--		--		--		5,000	U	--		--		--		--	
Toluene		108-88-3	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113		76-13-1	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
1,2,3-Trichlorobenzene		87-61-6	5	2.5	U	100	U	500	U	500	U	620	U	620	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	--	
1,2,4-Trichlorobenzene		120-82-1	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
1,1,2-Trichloroethane		79-00-5	1	60	U	300	U	300	U	380	U	380	U	750	U	750	U	1,500	U	1,500	U	1,500	U	1,500	U	2,000	U
Trichloroethene (TCE)		79-01-6	5	**6,500**		**6,100**		**4,300**		**8,100**		**4,900**		**5,000**		**30,000**		**650**		500	U	500	U	500	U	2,000	U
Trichlorofluoromethane (Freon 11)		75-69-4	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
Vinyl chloride		75-01-4	2	**2,200**		**2,400**	**J**	**970**		**2,100**		**1,500**		**2,800**		**12,000**		**38,000**		**99,000**		**66,000**		**71,000**		**120,000**	
o-Xylene		95-47-6	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	--	
m/p-Xylene		179601-23-1	5	100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	--	
Total Xylene				100	U	500	U	500	U	620	U	620	U	1,200	U	1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	4,000	U
Total VOCs		--	NE	**64,850**	**J**	**40,670**	**J**	**33,670**		**58,200**		**40,400**		**127,900**	**J**	**173,200**	**J**	**230,270**	**J**	**191,000**		**166,000**		**157,000**		**280,000**	

Table 4A. Historical Groundwater VOC Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW3I_G		MW3I		MW3I		MW3I		MW3I		MW3I		MW3I	
			Sample Date	2/8/2023		7/6/2023		11/30/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
Analyte	**Units**	**CAS No.**	**NYS AWQS**														
Volatile Organic Compounds	ug/L																
Acetone		67-64-1	50*	500	U	50	U	5	U	5	U	5	U	5	U	10	U
Benzene		71-43-2	1	50	U	5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Bromochloromethane		74-97-5	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	--	
Bromodichloromethane		75-27-4	50*	50	U	5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Bromoform		75-25-2	50*	200	U	20	UJ	2	U	2	U	2	U	2	U	1	U
Bromomethane		74-83-9	5	250	UJ	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Carbon disulfide		75-15-0	60*	500	UJ	50	U	5	U	5	U	5	U	5	U	1	U
Carbon tetrachloride		56-23-5	5	50	U	5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Chlorobenzene		108-90-7	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloroethane		75-00-3	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloroform (Trichloromethane)		67-66-3	7	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloromethane		74-87-3	5	250	U	25	U	2.5	UJ	2.5	U	2.5	U	2.5	U	1	U
Cyclohexane		110-82-7	NE	1000	U	100	U	10	U	10	U	10	U	10	U	1	U
1,2-Dibromo-3-chloropropane		96-12-8	0.04	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Dibromochloromethane		124-48-1	50*	50	U	5	UJ	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	200	U	20	U	2	U	2	U	2	U	2	U	1	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	500	U	50	U	5	U	5	U	5	U	5	U	1	U
1,1-Dichloroethane		75-34-3	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2-Dichloroethane		107-06-2	0.6	50	U	5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,1-Dichloroethene		75-35-4	5	50	U	5	U	**0.26**	**J**	**0.64**		**0.4**	**J**	0.5	U	1	U
cis-1,2-Dichloroethene		156-59-2	5	**4,000**		**970**		**140**		**42**		**56**		**5**		**2.9**	
trans-1,2-Dichloroethene		156-60-5	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2-Dichloropropane		78-87-5	1	100	U	10	U	1	U	1	U	1	U	1	U	1	U
cis-1,3-Dichloropropene		10061-01-5	0.4	50	U	5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
trans-1,3-Dichloropropene		10061-02-6	0.4	50	UJ	5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,4-Dioxane		123-91-1	NE	25,000	U	2,500	U	250	U	250	U	250	U	250	U	--	
Ethylbenzene		100-41-4	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
2-Hexanone		591-78-6	50*	500	U	50	U	5	U	5	U	5	U	5	U	5	U
Isopropylbenzene		98-82-8	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Methyl acetate		79-20-9	NE	200	U	20	U	2	UJ	2	U	2	U	2	U	2.5	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	500	U	50	U	5	U	5	U	5	U	5	U	10	U
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	500	U	50	U	5	UJ	5	U	5	U	5	U	5	U
Methylcyclohexane		108-87-2	NE	1000	U	100	U	10	U	10	U	10	U	10	U	1	U
Methylene chloride		75-09-2	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Styrene		100-42-5	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2,2-Tetrachloroethane		79-34-5	5	50	U	5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Tetrachloroethene (PCE)		127-18-4	5	**13,000**		**19**		**2.3**		**0.64**		**1.1**		**0.41**	**J**	1	U
Tetrahydrofuran		109-99-9	50*	--		--		5	U	--		--		--		--	
Toluene		108-88-3	5	**150**	**J**	**26**		**1.8**	**J**	2.5	U	2.5	U	2.5	U	1	U
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113		76-13-1	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2,3-Trichlorobenzene		87-61-6	5	2,500	U	250	U	25	U	2.5	U	2.5	U	2.5	U	--	
1,2,4-Trichlorobenzene		120-82-1	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2-Trichloroethane		79-00-5	1	150	U	15	U	1.5	U	1.5	U	1.5	U	1.5	U	1	U
Trichloroethene (TCE)		79-01-6	5	**690**		**18**		**1.4**		**0.6**		**1.6**		**0.75**		**0.53**	**J**
Trichlorofluoromethane (Freon 11)		75-69-4	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride		75-01-4	2	**120**		**8.6**	**J**	**7.3**		**39**		**24**		**6.1**		**8.4**	
o-Xylene		95-47-6	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	--	
m/p-Xylene		179601-23-1	5	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	--	
Total Xylene				250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U
Total VOCs		--	NE	**17,960**	**J**	**1,041.60**	**J**	**153.06**	**J**	**82.88**		**83.1**	**J**	**12.26**	**J**	**11.83**	**J**

Table 4A. Historical Groundwater VOC Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	TW9		TW9		TW9		TW9		MW4S_G		MW4S_G (DUI		MW4S		MW4S		MW4S		MW4S		MW4S		MW4S (DUP)		MW4S	
			Sample Date	3/21/2021		7/11/2021		11/20/2021		2/26/2022		2/8/2023		2/8/2023		7/6/2023		12/1/2023		12/10/2024		3/12/2025		6/17/2025		6/17/2025		9/24/2025	
Analyte	**Units**	**CAS No.**	**NYS AWQS**																										
Volatile Organic Compounds	ug/L																												
Acetone		67-64-1	50*	100	U	10	U	20	U	20	U	2,500	U	2,500	U	5,000	U	1000	U	5,000	U	500	U	500	U	1000	U	20,000	U
Benzene		71-43-2	1	10	U	1	U	2	U	2	U	250	U	250	U	500	U	100	U	500	U	50	U	50	U	100	U	2,000	U
Bromochloromethane		74-97-5	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	--	
Bromodichloromethane		75-27-4	50*	10	U	1	U	2	U	2	U	250	U	250	U	500	U	100	U	500	U	50	U	50	U	100	U	2,000	U
Bromoform		75-25-2	50*	40	U	4	U	8	U	8	U	1000	U	1000	U	2,000	UJ	400	U	2,000	U	200	U	200	U	400	U	2,000	U
Bromomethane		74-83-9	5	50	UJ	5	UJ	10	U	10	U	1,200	UJ	1,200	UJ	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
Carbon disulfide		75-15-0	60*	100	U	10	U	20	U	20	U	2,500	UJ	2,500	UJ	5,000	U	1000	U	5,000	U	500	U	500	U	1000	U	2,000	U
Carbon tetrachloride		56-23-5	5	10	U	1	U	2	U	2	U	250	U	250	U	500	U	100	U	500	U	50	U	50	U	100	U	2,000	U
Chlorobenzene		108-90-7	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
Chloroethane		75-00-3	5	50	UJ	5	UJ	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
Chloroform (Trichloromethane)		67-66-3	7	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
Chloromethane		74-87-3	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	UJ	2,500	U	250	U	250	U	500	U	2,000	U
Cyclohexane		110-82-7	NE	200	U	20	U	40	U	40	U	5,000	U	5,000	U	10,000	U	2,000	U	10,000	U	1000	U	1000	U	2,000	U	2,000	U
1,2-Dibromo-3-chloropropane		96-12-8	0.04	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
Dibromochloromethane		124-48-1	50*	10	U	1	U	2	U	2	U	250	U	250	U	500	UJ	100	U	500	U	50	U	50	U	100	U	2,000	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	40	U	4	U	8	U	8	U	1000	U	1000	U	2,000	U	400	U	2,000	U	200	U	200	U	400	U	2,000	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	100	U	10	UJ	20	U	20	U	2,500	U	2,500	U	5,000	U	1000	U	5,000	U	500	U	500	U	1000	U	2,000	U
1,1-Dichloroethane		75-34-3	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
1,2-Dichloroethane		107-06-2	0.6	10	U	1	U	2	U	2	U	250	U	250	U	500	U	100	U	500	U	50	U	50	U	100	U	2,000	U
1,1-Dichloroethene		75-35-4	5	**3.4**	**J**	1	U	2	U	2	U	**100**	**J**	**94**	**J**	500	U	**35**	**J**	500	U	50	U	50	U	100	U	2,000	U
cis-1,2-Dichloroethene		156-59-2	5	**1,800**		**290**		**470**		**460**		**76,000**		**70,000**		**16,000**		**17,000**		**9,600**		**1000**		**1,400**		**1,400**		**8,100**	
trans-1,2-Dichloroethene		156-60-5	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
1,2-Dichloropropane		78-87-5	1	20	U	2	U	4	U	4	U	500	U	500	U	1000	U	200	U	1000	U	100	U	100	U	200	U	2,000	U
cis-1,3-Dichloropropene		10061-01-5	0.4	10	U	1	U	2	U	2	U	250	U	250	U	500	U	100	U	500	U	50	U	50	U	100	U	2,000	U
trans-1,3-Dichloropropene		10061-02-6	0.4	10	U	1	U	2	U	2	U	250	UJ	250	UJ	500	U	100	U	500	U	50	U	50	U	100	U	2,000	U
1,4-Dioxane		123-91-1	NE	5,000	UJ	500	U	1000	U	1000	U	120,000	U	120,000	U	250,000	U	50,000	U	250,000	U	25,000	U	25,000	U	50,000	U	--	
Ethylbenzene		100-41-4	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
2-Hexanone		591-78-6	50*	100	U	10	U	20	U	20	U	2,500	U	2,500	U	5,000	U	1000	U	5,000	U	500	U	500	U	1000	U	10,000	U
Isopropylbenzene		98-82-8	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
Methyl acetate		79-20-9	NE	40	UJ	4	U	8	U	8	U	1000	U	1000	U	2,000	U	400	UJ	2,000	U	200	U	200	U	400	U	5,000	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	100	U	10	U	20	U	20	U	2,500	U	2,500	U	5,000	U	1000	U	5,000	U	500	U	500	U	1000	U	20,000	U
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	100	U	10	U	20	U	20	U	2,500	U	2,500	U	5,000	U	1000	UJ	5,000	U	500	U	500	U	1000	U	10,000	U
Methylcyclohexane		108-87-2	NE	200	U	20	U	40	U	40	U	5,000	U	5,000	U	10,000	U	2,000	U	10,000	U	1000	U	1000	U	2,000	U	2,000	U
Methylene chloride		75-09-2	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
Styrene		100-42-5	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
1,1,2,2-Tetrachloroethane		79-34-5	5	10	U	1	U	2	U	2	U	250	U	250	U	500	U	100	U	500	U	50	U	50	U	100	U	2,000	U
Tetrachloroethene (PCE)		127-18-4	5	**390**		**53**		**140**		**61**		**43,000**		**41,000**		**120,000**		**140,000**	**J**	**82,000**		**19,000**		**19,000**		**20,000**		**85,000**	
Tetrahydrofuran		109-99-9	50*	--		--		--		--		--		--		--		1000	U	--		--		--		--		--	
Toluene		108-88-3	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113		76-13-1	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
1,2,3-Trichlorobenzene		87-61-6	5	2.5	U	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	--	
1,2,4-Trichlorobenzene		120-82-1	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
1,1,2-Trichloroethane		79-00-5	1	30	U	3	U	6	U	6	U	750	U	750	U	1,500	U	300	U	1,500	U	150	U	150	U	300	U	2,000	U
Trichloroethene (TCE)		79-01-6	5	**400**		**27**		**40**		**47**		**7,200**		**6,800**		**7,400**		**5,900**		**4,500**		**590**		**690**		**730**		**4,100**	
Trichlorofluoromethane (Freon 11)		75-69-4	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	2,000	U
Vinyl chloride		75-01-4	2	**120**		**10**	**J**	**7.8**		**24**		**8,500**		**8,100**		**1,400**		**870**		**860**	**J**	**84**	**J**	**96**	**J**	**100**	**J**	2,000	U
o-Xylene		95-47-6	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	--	
m/p-Xylene		179601-23-1	5	50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	--	
Total Xylene				50	U	5	U	10	U	10	U	1,200	U	1,200	U	2,500	U	500	U	2,500	U	250	U	250	U	500	U	4,000	U
Total VOCs		--	NE	**2,713.40**	**J**	**380**	**J**	**657.8**		**592**		**134,800**	**J**	**125,994**	**J**	**144,800**		**163,805**	**J**	**96,960**	**J**	**20,674**	**J**	**21,186**	**J**	**22,230**	**J**	**97,200**	

Table 4A. Historical Groundwater VOC Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW4I_G		MW4I		MW4I		MW4I (DUP)		MW4I		MW 4I (DUP)		MW4I		MW4I		MW4I	
			Sample Date	2/8/2023		7/6/2023		12/1/2023		12/1/2023		12/10/2024		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
Analyte	**Units**	**CAS No.**	**NYS AWQS**																		
Volatile Organic Compounds	ug/L																				
Acetone		67-64-1	50*	500	U	250	U	250	U	250	U	120	U	120	U	100	U	120	U	400	U
Benzene		71-43-2	1	50	U	25	U	25	U	25	U	12	U	12	U	10	U	12	U	40	U
Bromochloromethane		74-97-5	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	--	
Bromodichloromethane		75-27-4	50*	50	U	25	U	25	U	25	U	12	U	12	U	10	U	12	U	40	U
Bromoform		75-25-2	50*	200	U	100	U	100	U	100	U	50	U	50	U	40	U	50	U	40	U
Bromomethane		74-83-9	5	250	UJ	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
Carbon disulfide		75-15-0	60*	500	UJ	250	U	250	U	250	U	120	U	120	U	100	U	120	U	40	U
Carbon tetrachloride		56-23-5	5	50	U	25	U	25	U	25	U	12	U	12	U	10	U	12	U	40	U
Chlorobenzene		108-90-7	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
Chloroethane		75-00-3	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
Chloroform (Trichloromethane)		67-66-3	7	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
Chloromethane		74-87-3	5	250	U	120	U	120	UJ	120	UJ	62	U	62	U	50	U	62	U	40	U
Cyclohexane		110-82-7	NE	1000	U	500	U	500	U	500	U	250	U	250	U	200	U	250	U	40	U
1,2-Dibromo-3-chloropropane		96-12-8	0.04	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
Dibromochloromethane		124-48-1	50*	50	U	25	U	25	U	25	U	12	U	12	U	10	U	12	U	40	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	200	U	100	U	100	U	100	U	50	U	50	U	40	U	50	U	40	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	500	U	250	U	250	U	250	U	120	U	120	U	100	U	120	U	40	U
1,1-Dichloroethane		75-34-3	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
1,2-Dichloroethane		107-06-2	0.6	50	U	25	U	25	U	25	U	12	U	12	U	10	U	12	U	40	U
1,1-Dichloroethene		75-35-4	5	50	U	**25**		**29**		**26**		**23**		**26**		**24**		**15**		**36**	**J**
cis-1,2-Dichloroethene		156-59-2	5	**8,500**		**3,600**		**6,700**		**6,000**		**3,200**		**3,200**		**2,900**		**2,000**		**3,800**	**F1**
trans-1,2-Dichloroethene		156-60-5	5	250	U	120	U	120	U	120	U	**25**	**J**	**23**	**J**	50	U	62	U	40	U
1,2-Dichloropropane		78-87-5	1	100	U	50	U	50	U	50	U	25	U	25	U	20	U	25	U	40	U
cis-1,3-Dichloropropene		10061-01-5	0.4	50	U	25	U	25	U	25	U	12	U	12	U	10	U	12	U	40	U
trans-1,3-Dichloropropene		10061-02-6	0.4	50	UJ	25	U	25	U	25	U	12	U	12	U	10	U	12	U	40	U
1,4-Dioxane		123-91-1	NE	25,000	U	12,000	U	12,000	U	12,000	U	6,200	U	6,200	U	5,000	U	6,200	U	--	
Ethylbenzene		100-41-4	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
2-Hexanone		591-78-6	50*	500	U	250	U	250	U	250	U	120	U	120	U	100	U	120	U	200	U
Isopropylbenzene		98-82-8	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
Methyl acetate		79-20-9	NE	200	U	100	U	100	UJ	100	UJ	50	U	50	U	40	U	50	U	100	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	500	U	250	U	250	U	250	U	120	U	120	U	100	U	120	U	400	U
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	500	U	250	U	250	UJ	250	UJ	120	U	120	U	100	U	120	U	200	U
Methylcyclohexane		108-87-2	NE	1000	U	500	U	500	U	500	U	250	U	250	U	200	U	250	U	40	U
Methylene chloride		75-09-2	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
Styrene		100-42-5	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
1,1,2,2-Tetrachloroethane		79-34-5	5	50	U	25	U	25	U	25	U	12	U	12	U	10	U	12	U	40	U
Tetrachloroethene (PCE)		127-18-4	5	**5,800**		**7,700**	**J**	**720**		**670**		**36**		**34**		**31**		**16**		**20**	**J**
Tetrahydrofuran		109-99-9	50*	--		--		250	U	250	U	--		--		--		--		**--**	
Toluene		108-88-3	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113		76-13-1	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
1,2,3-Trichlorobenzene		87-61-6	5	500	U	250	U	120	U	120	U	120	U	62	U	62	U	50	U	--	
1,2,4-Trichlorobenzene		120-82-1	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
1,1,2-Trichloroethane		79-00-5	1	150	U	75	U	75	U	75	U	38	U	38	U	30	U	38	U	40	U
Trichloroethene (TCE)		79-01-6	5	**720**		**2,000**		**2,200**		**1,900**		**70**		**78**		**28**		**12**		40	U
Trichlorofluoromethane (Freon 11)		75-69-4	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	40	U
Vinyl chloride		75-01-4	2	**540**		**360**		**330**		**330**		**390**		**380**		**300**		**280**		**550**	
o-Xylene		95-47-6	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	**--**	
m/p-Xylene		179601-23-1	5	250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	**--**	
Total Xylene				250	U	120	U	120	U	120	U	62	U	62	U	50	U	62	U	80	U
Total VOCs		--	NE	**15,560**		**13,685**	**J**	**9,979**		**8,926**		**3,744**	**J**	**3,741**	**J**	**3,283**		**2,323**		**4,406**	**J**

Table 4A. Historical Groundwater VOC Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW5S		MW5S		MW5S		MW5S		MW5S		MW5S		MW5S	
			Sample Date	2/21/2023		7/5/2023		11/29/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
Analyte	**Units**	**CAS No.**	**NYS AWQS**														
Volatile Organic Compounds	ug/L																
Acetone		67-64-1	50*	20	U	25	U	25	U	25	U	**94**		**61**		**86**	**J**
Benzene		71-43-2	1	**72**		**41**		**19**		**14**		**22**		**22**		**18**	
Bromochloromethane		74-97-5	5	10	U	12	U	12	U	12	U	12	U	10	U	--	
Bromodichloromethane		75-27-4	50*	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U	10	U
Bromoform		75-25-2	50*	8	UJ	10	U	10	U	10	U	10	U	8	U	10	U
Bromomethane		74-83-9	5	10	U	12	U	12	U	12	U	12	U	10	U	10	U
Carbon disulfide		75-15-0	60*	20	U	25	U	25	U	25	U	25	U	20	U	10	U
Carbon tetrachloride		56-23-5	5	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U	10	U
Chlorobenzene		108-90-7	5	10	U	12	U	12	U	12	U	12	U	10	U	10	U
Chloroethane		75-00-3	5	10	U	12	U	12	U	12	U	12	U	10	U	10	U
Chloroform (Trichloromethane)		67-66-3	7	10	U	12	U	12	U	12	U	12	U	10	U	10	U
Chloromethane		74-87-3	5	10	U	12	U	12	U	12	U	12	U	10	U	10	U
Cyclohexane		110-82-7	NE	**230**		**270**		**200**		**120**		**200**		**200**		**370**	
1,2-Dibromo-3-chloropropane		96-12-8	0.04	10	U	12	U	12	U	12	U	12	U	10	U	10	U
Dibromochloromethane		124-48-1	50*	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U	10	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	8	U	10	U	10	U	10	U	10	U	8	U	10	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	10	U	12	U	12	U	12	U	12	U	10	U	10	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	10	U	12	U	12	U	12	U	12	U	10	U	10	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	10	U	12	U	12	U	12	U	12	U	10	U	10	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	20	U	25	U	25	U	25	U	25	U	20	U	10	U
1,1-Dichloroethane		75-34-3	5	10	U	12	U	12	U	12	U	12	U	10	U	10	U
1,2-Dichloroethane		107-06-2	0.6	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U	10	U
1,1-Dichloroethene		75-35-4	5	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U	10	U
cis-1,2-Dichloroethene		156-59-2	5	10	U	12	U	12	U	12	U	12	U	10	U	10	U
trans-1,2-Dichloroethene		156-60-5	5	10	U	12	U	12	U	12	U	12	U	10	U	10	U
1,2-Dichloropropane		78-87-5	1	4	U	5	U	5	U	5	U	5	U	4	U	10	U
cis-1,3-Dichloropropene		10061-01-5	0.4	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U	10	U
trans-1,3-Dichloropropene		10061-02-6	0.4	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U	10	U
1,4-Dioxane		123-91-1	NE	1000	U	1,200	U	1,200	U	1,200	U	1,200	U	1000	U	--	
Ethylbenzene		100-41-4	5	**500**		**780**		**600**		**380**		**560**		**370**		**260**	
2-Hexanone		591-78-6	50*	20	U	25	U	25	U	25	U	25	U	20	U	50	U
Isopropylbenzene		98-82-8	5	**54**		**100**		**96**		**85**		**120**		**89**		**110**	
Methyl acetate		79-20-9	NE	8	U	10	U	10	U	10	U	10	U	8	U	25	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	20	U	25	U	25	U	25	U	25	U	20	U	**34**	**J**
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	10	U	12	U	12	U	12	U	12	U	10	U	10	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	20	U	25	U	25	U	25	U	25	U	20	U	50	U
Methylcyclohexane		108-87-2	NE	**160**		**180**		**120**		**76**		**110**		**98**		**150**	
Methylene chloride		75-09-2	5	10	U	12	U	12	U	12	U	**4.9**	**J**	**3**	**J**	10	U
Styrene		100-42-5	5	10	U	12	U	12	U	12	U	12	U	10	U	10	U
1,1,2,2-Tetrachloroethane		79-34-5	5	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U	10	U
Tetrachloroethene (PCE)		127-18-4	5	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U	10	U
Tetrahydrofuran		109-99-9	50*	--		--		25	U	--		--		--		--	
Toluene		108-88-3	5	**11**		**11**	**J**	**7.4**	**J**	**5.8**	**J**	**7**	**J**	**5.5**	**J**	**6.4**	**J**
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113		76-13-1	5	10	U	12	U	12	U	12	U	12	U	10	U	10	U
1,2,3-Trichlorobenzene		87-61-6	5	62	U	10	U	12	U	12	U	12	U	12	U	--	
1,2,4-Trichlorobenzene		120-82-1	5	10	U	12	U	12	U	12	U	12	U	10	U	10	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	10	U	12	U	12	U	12	U	12	U	10	U	10	U
1,1,2-Trichloroethane		79-00-5	1	6	U	7.5	U	7.5	U	7.5	U	7.5	U	6	U	10	U
Trichloroethene (TCE)		79-01-6	5	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U	10	U
Trichlorofluoromethane (Freon 11)		75-69-4	5	10	U	12	U	12	U	12	U	12	U	10	U	10	U
Vinyl chloride		75-01-4	2	**2**	**J**	**0.58**	**J**	5	U	5	U	5	U	**2.4**	**J**	10	U
o-Xylene		95-47-6	5	**11**		**8.3**	**J**	**6.9**	**J**	**5.2**	**J**	**4.6**	**J**	**3.4**	**J**	--	
m/p-Xylene		179601-23-1	5	**40**		**42**		**30**		**19**		**24**		**17**		--	
Total Xylene				**51**		**50.3**		**36.9**		**24.2**		**28.6**		**20.4**		**23**	
Total VOCs		--	NE	**1,080**		**1,432.88**	**J**	**1,079.30**	**J**	**705**	**J**	**1,146.50**	**J**	**871.3**	**J**	**1,057.40**	**J**

Table 4A. Historical Groundwater VOC Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW5I		MW5I		MW5I		MW5I		MW5I		MW5I		MW5I	
			Sample Date	2/21/2023		7/5/2023		11/29/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
Analyte	**Units**	**CAS No.**	**NYS AWQS**														
Volatile Organic Compounds	ug/L																
Acetone		67-64-1	50*	5	U	5	U	5	U	5	U	5	U	**2.2**	**J**	10	U
Benzene		71-43-2	1	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Bromochloromethane		74-97-5	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	--	
Bromodichloromethane		75-27-4	50*	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Bromoform		75-25-2	50*	2	UJ	2	UJ	2	U	2	U	2	U	2	U	1	U
Bromomethane		74-83-9	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Carbon disulfide		75-15-0	60*	5	U	5	U	5	U	5	U	5	U	5	U	1	U
Carbon tetrachloride		56-23-5	5	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Chlorobenzene		108-90-7	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloroethane		75-00-3	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloroform (Trichloromethane)		67-66-3	7	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloromethane		74-87-3	5	2.5	U	2.5	U	2.5	UJ	2.5	U	2.5	U	2.5	U	1	U
Cyclohexane		110-82-7	NE	**1.4**	**J**	**0.38**	**J**	10	U	10	U	10	U	10	U	1	U
1,2-Dibromo-3-chloropropane		96-12-8	0.04	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Dibromochloromethane		124-48-1	50*	0.5	U	0.5	UJ	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	2	U	2	U	2	U	2	U	2	U	2	U	1	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	5	U	5	U	5	U	5	U	5	U	5	U	1	U
1,1-Dichloroethane		75-34-3	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2-Dichloroethane		107-06-2	0.6	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,1-Dichloroethene		75-35-4	5	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
cis-1,2-Dichloroethene		156-59-2	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
trans-1,2-Dichloroethene		156-60-5	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2-Dichloropropane		78-87-5	1	1	U	1	U	1	U	1	U	1	U	1	U	1	U
cis-1,3-Dichloropropene		10061-01-5	0.4	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
trans-1,3-Dichloropropene		10061-02-6	0.4	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,4-Dioxane		123-91-1	NE	250	U	250	U	250	U	250	U	250	U	250	U	--	
Ethylbenzene		100-41-4	5	**1.4**	**J**	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
2-Hexanone		591-78-6	50*	5	U	5	U	5	UJ	5	U	5	U	5	U	5	U
Isopropylbenzene		98-82-8	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Methyl acetate		79-20-9	NE	2	U	2	U	2	UJ	2	U	2	U	2	U	2.5	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	5	U	5	U	5	U	5	U	5	U	5	U	10	U
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	5	U	5	U	5	UJ	5	U	5	U	5	U	5	U
Methylcyclohexane		108-87-2	NE	**2.2**	**J**	**0.65**	**J**	**0.4**	**J**	10	U	10	U	10	U	1	U
Methylene chloride		75-09-2	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Styrene		100-42-5	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2,2-Tetrachloroethane		79-34-5	5	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Tetrachloroethene (PCE)		127-18-4	5	**0.3**	**J**	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Tetrahydrofuran		109-99-9	50*	--		--		5	U	--		--		--		--	
Toluene		108-88-3	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113		76-13-1	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2,3-Trichlorobenzene		87-61-6	5	10	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	--	
1,2,4-Trichlorobenzene		120-82-1	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2-Trichloroethane		79-00-5	1	1.5	U	1.5	U	1.5	U	1.5	U	1.5	U	1.5	U	1	U
Trichloroethene (TCE)		79-01-6	5	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Trichlorofluoromethane (Freon 11)		75-69-4	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride		75-01-4	2	1	U	1	U	1	U	1	U	1	U	**1.1**		1	U
o-Xylene		95-47-6	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	--	
m/p-Xylene		179601-23-1	5	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	--	
Total Xylene				2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U
Total VOCs		--	NE	**5.3**	**J**	**1.03**	**J**	**0.4**	**J**	ND		ND		**3.3**	**J**	ND	

Table 4A. Historical Groundwater VOC Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW6		MW6		MW6		MW6		MW6		MW7		MW7		MW7		MW7		MW7	
			Sample Date	11/28/2023		12/9/2024		3/11/2025		6/16/2025		9/23/2025		11/28/2023		12/9/2024		3/11/2025		6/16/2025		9/23/2025	
Analyte	**Units**	**CAS No.**	**NYS AWQS**																				
Volatile Organic Compounds	ug/L																						
Acetone		67-64-1	50*	5	U	5	U	5	U	5	U	10	U	5	U	5	U	5	U	**4.9**	**J**	10	U
Benzene		71-43-2	1	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Bromochloromethane		74-97-5	5	2.5	U	2.5	U	2.5	U	2.5	U	--		2.5	U	2.5	U	2.5	U	2.5	U	--	
Bromodichloromethane		75-27-4	50*	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Bromoform		75-25-2	50*	2	U	2	U	2	U	2	U	1	U	2	U	2	U	2	U	2	U	1	U
Bromomethane		74-83-9	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Carbon disulfide		75-15-0	60*	5	U	5	U	5	U	5	U	1	U	5	U	5	U	5	U	5	U	1	U
Carbon tetrachloride		56-23-5	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Chlorobenzene		108-90-7	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloroethane		75-00-3	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloroform (Trichloromethane)		67-66-3	7	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloromethane		74-87-3	5	2.5	UJ	2.5	U	2.5	U	2.5	U	1	U	2.5	UJ	2.5	U	2.5	U	2.5	U	1	U
Cyclohexane		110-82-7	NE	10	U	10	U	10	U	10	U	1	U	10	U	10	U	10	U	10	U	1	U
1,2-Dibromo-3-chloropropane		96-12-8	0.04	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Dibromochloromethane		124-48-1	50*	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	2	U	2	U	2	U	2	U	1	U	2	U	2	U	2	U	2	U	1	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	5	U	5	U	5	U	5	U	1	U	5	U	5	U	5	U	5	U	1	U
1,1-Dichloroethane		75-34-3	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2-Dichloroethane		107-06-2	0.6	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,1-Dichloroethene		75-35-4	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
cis-1,2-Dichloroethene		156-59-2	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
trans-1,2-Dichloroethene		156-60-5	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2-Dichloropropane		78-87-5	1	1	U	1	U	1	U	1	U	1	U	1	U	1	U	1	U	1	U	1	U
cis-1,3-Dichloropropene		10061-01-5	0.4	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
trans-1,3-Dichloropropene		10061-02-6	0.4	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,4-Dioxane		123-91-1	NE	250	U	250	U	250	U	250	U	--		250	U	250	U	250	U	250	U	--	
Ethylbenzene		100-41-4	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
2-Hexanone		591-78-6	50*	5	U	5	U	5	U	5	U	5	U	5	U	5	U	5	U	5	U	5	U
Isopropylbenzene		98-82-8	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Methyl acetate		79-20-9	NE	2	UJ	2	U	2	U	2	U	2.5	U	2	UJ	2	U	2	U	2	U	2.5	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	5	U	5	U	5	U	5	U	10	U	5	U	5	U	5	U	5	U	10	U
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	5	UJ	5	U	5	U	5	U	5	U	5	UJ	5	U	5	U	5	U	5	U
Methylcyclohexane		108-87-2	NE	10	U	10	U	10	U	10	U	1	U	10	U	10	U	10	U	10	U	1	U
Methylene chloride		75-09-2	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Styrene		100-42-5	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2,2-Tetrachloroethane		79-34-5	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Tetrachloroethene (PCE)		127-18-4	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Tetrahydrofuran		109-99-9	50*	5	U	--		--		--		--		5	U	--		--		--		--	
Toluene		108-88-3	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113		76-13-1	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2,3-Trichlorobenzene		87-61-6	5	2.5	U	2.5	U	2.5	U	2.5	U	--		2.5	U	2.5	U	2.5	U	2.5	U	--	
1,2,4-Trichlorobenzene		120-82-1	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2-Trichloroethane		79-00-5	1	1.5	U	1.5	U	1.5	U	1.5	U	1	U	1.5	U	1.5	U	1.5	U	1.5	U	1	U
Trichloroethene (TCE)		79-01-6	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Trichlorofluoromethane (Freon 11)		75-69-4	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride		75-01-4	2	1	U	1	U	1	U	**4.1**		1	U	**4.7**		**15**		**12**		**15**		**13**	
o-Xylene		95-47-6	5	2.5	U	2.5	U	2.5	U	2.5	U	--		2.5	U	2.5	U	2.5	U	2.5	U	**--**	
m/p-Xylene		179601-23-1	5	2.5	U	2.5	U	2.5	U	2.5	U	--		2.5	U	2.5	U	2.5	U	2.5	U	**--**	
Total Xylene				2.5	U	2.5	U	2.5	U	2.5	U	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U
Total VOCs		--	NE	ND		ND		ND		**4.1**		ND		**4.7**		**15**		**12**		**19.9**	**J**	**13**	

Table 4A. Historical Groundwater VOC Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW8		MW8		MW8		MW8		MW8		MW11		MW11		MW11		MW11		MW11	
			Sample Date	11/29/2023		12/9/2024		3/11/2025		6/16/2025		9/24/2025		11/29/2023		12/11/2024		3/11/2025		6/16/2025		9/24/2025	
Analyte	**Units**	**CAS No.**	**NYS AWQS**																				
Volatile Organic Compounds	ug/L																						
Acetone		67-64-1	50*	5	U	**2.2**	**J**	**3.9**	**J**	**8.1**		**3.5**	**J**	5	U	5	U	5	U	**1.9**	**J**	10	U
Benzene		71-43-2	1	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Bromochloromethane		74-97-5	5	2.5	U	2.5	U	2.5	U	2.5	U	--		2.5	U	2.5	U	2.5	U	2.5	U	--	
Bromodichloromethane		75-27-4	50*	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Bromoform		75-25-2	50*	2	U	2	U	2	U	2	U	1	U	2	U	2	U	2	U	2	U	1	U
Bromomethane		74-83-9	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Carbon disulfide		75-15-0	60*	5	U	5	U	5	U	5	U	1	U	5	U	5	U	5	U	5	U	1	U
Carbon tetrachloride		56-23-5	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Chlorobenzene		108-90-7	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloroethane		75-00-3	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloroform (Trichloromethane)		67-66-3	7	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloromethane		74-87-3	5	2.5	UJ	2.5	U	2.5	U	2.5	U	1	U	2.5	UJ	2.5	U	2.5	U	2.5	U	1	U
Cyclohexane		110-82-7	NE	10	U	10	U	10	U	10	U	1	U	10	U	10	U	10	U	10	U	1	U
1,2-Dibromo-3-chloropropane		96-12-8	0.04	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Dibromochloromethane		124-48-1	50*	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	2	U	2	U	2	U	2	U	1	U	2	U	2	U	2	U	2	U	1	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	5	U	5	U	5	U	5	U	1	U	5	U	5	U	5	U	5	U	1	U
1,1-Dichloroethane		75-34-3	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2-Dichloroethane		107-06-2	0.6	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,1-Dichloroethene		75-35-4	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
cis-1,2-Dichloroethene		156-59-2	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
trans-1,2-Dichloroethene		156-60-5	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2-Dichloropropane		78-87-5	1	1	U	1	U	1	U	1	U	1	U	1	U	1	U	1	U	1	U	1	U
cis-1,3-Dichloropropene		10061-01-5	0.4	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
trans-1,3-Dichloropropene		10061-02-6	0.4	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,4-Dioxane		123-91-1	NE	250	U	250	U	250	U	250	U	--		250	U	250	U	250	U	250	U	--	
Ethylbenzene		100-41-4	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
2-Hexanone		591-78-6	50*	5	U	5	U	5	U	5	U	5	U	5	U	5	U	5	U	5	U	5	U
Isopropylbenzene		98-82-8	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Methyl acetate		79-20-9	NE	2	UJ	2	U	2	U	2	U	2.5	U	2	UJ	2	U	2	U	2	U	2.5	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	5	U	5	U	5	U	5	U	10	U	5	U	5	U	5	U	5	U	10	U
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	2.5	U	**0.34**	**J**	**0.2**	**J**	**0.31**	**J**	1	U	2.5	U	**0.28**	**J**	2.5	U	2.5	U	1	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	5	UJ	5	U	5	U	5	U	5	U	5	UJ	5	U	5	U	5	U	5	U
Methylcyclohexane		108-87-2	NE	10	U	10	U	10	U	10	U	1	U	10	U	10	U	10	U	10	U	1	U
Methylene chloride		75-09-2	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Styrene		100-42-5	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2,2-Tetrachloroethane		79-34-5	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Tetrachloroethene (PCE)		127-18-4	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Tetrahydrofuran		109-99-9	50*	5	U	--		--		--		--		--		--		--		--		--	
Toluene		108-88-3	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113		76-13-1	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2,3-Trichlorobenzene		87-61-6	5	2.5	U	2.5	U	2.5	U	2.5	U	--		2.5	U	2.5	U	2.5	U	2.5	U	--	
1,2,4-Trichlorobenzene		120-82-1	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2-Trichloroethane		79-00-5	1	1.5	U	1.5	U	1.5	U	1.5	U	1	U	1.5	U	1.5	U	1.5	U	1.5	U	1	U
Trichloroethene (TCE)		79-01-6	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U	**0.27**	**J**	0.5	U	0.5	U	0.5	U	1	U
Trichlorofluoromethane (Freon 11)		75-69-4	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride		75-01-4	2	1	U	**0.21**	**J**	**0.11**	**J**	**0.43**	**J**	1	U	**0.34**	**J**	**0.36**	**J**	**0.14**	**J**	**2.6**		1	U
o-Xylene		95-47-6	5	2.5	U	2.5	U	2.5	U	2.5	U	--		2.5	U	2.5	U	2.5	U	2.5	U	--	
m/p-Xylene		179601-23-1	5	2.5	U	2.5	U	2.5	U	2.5	U	--		2.5	U	2.5	U	2.5	U	2.5	U	--	
Total Xylene				2.5	U	2.5	U	2.5	U	2.5	U	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U
Total VOCs		--	NE	ND		**2.75**	**J**	**4.21**	**J**	**8.84**	**J**	**3.5**		**0.61**	**J**	**0.64**	**J**	**0.14**	**J**	**4.5**	**J**	ND	

Table 4A. Historical Groundwater VOC Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW12S		MW12S		MW12S		MW12S		MW12S	
			Sample Date	11/30/2023		12/11/2024		3/11/2025		6/18/2025		9/24/2025	
Analyte	**Units**	**CAS No.**	**NYS AWQS**										
Volatile Organic Compounds	ug/L												
Acetone		67-64-1	50*	5	U	5	U	5	U	**2.7**	**J**	10	U
Benzene		71-43-2	1	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Bromochloromethane		74-97-5	5	2.5	U	2.5	U	2.5	U	2.5	U	--	
Bromodichloromethane		75-27-4	50*	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Bromoform		75-25-2	50*	2	U	2	U	2	U	2	U	1	U
Bromomethane		74-83-9	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Carbon disulfide		75-15-0	60*	5	U	5	U	5	U	5	U	1	U
Carbon tetrachloride		56-23-5	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Chlorobenzene		108-90-7	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloroethane		75-00-3	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloroform (Trichloromethane)		67-66-3	7	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Chloromethane		74-87-3	5	2.5	UJ	2.5	U	2.5	U	2.5	U	1	U
Cyclohexane		110-82-7	NE	10	U	10	U	10	U	10	U	1	U
1,2-Dibromo-3-chloropropane		96-12-8	0.04	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Dibromochloromethane		124-48-1	50*	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,2-Dibromoethane (EDB)		106-93-4	0.0006	2	U	2	U	2	U	2	U	1	U
1,2-Dichlorobenzene (o-DCB)		95-50-1	3	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,3-Dichlorobenzene (m-DCB)		541-73-1	3	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,4-Dichlorobenzene (p-DCB)		106-46-7	3	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Dichlorodifluoromethane (Freon 12)		75-71-8	5	5	U	5	U	5	U	5	U	1	U
1,1-Dichloroethane		75-34-3	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2-Dichloroethane		107-06-2	0.6	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,1-Dichloroethene		75-35-4	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U
cis-1,2-Dichloroethene		156-59-2	5	**7.2**		**3.4**		**2.3**	**J**	**1.7**	**J**	**2.3**	
trans-1,2-Dichloroethene		156-60-5	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2-Dichloropropane		78-87-5	1	1	U	1	U	1	U	1	U	1	U
cis-1,3-Dichloropropene		10061-01-5	0.4	0.5	U	0.5	U	0.5	U	0.5	U	1	U
trans-1,3-Dichloropropene		10061-02-6	0.4	0.5	U	0.5	U	0.5	U	0.5	U	1	U
1,4-Dioxane		123-91-1	NE	250	U	250	U	250	U	250	U	--	
Ethylbenzene		100-41-4	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
2-Hexanone		591-78-6	50*	5	U	5	U	5	U	5	U	5	U
Isopropylbenzene		98-82-8	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Methyl acetate		79-20-9	NE	2	UJ	2	U	2	U	2	U	2.5	U
Methyl ethyl ketone (2-Butanone)		78-93-3	50*	5	U	5	U	5	U	5	U	10	U
Methyl tert-butyl ether (MTBE)		1634-04-4	10*	2.5	U	2.5	U	2.5	U	2.5	U	1	U
4-Methyl-2-pentanone (MIBK)		108-10-1	NE	5	UJ	5	U	5	U	5	U	5	U
Methylcyclohexane		108-87-2	NE	10	U	10	U	10	U	10	U	1	U
Methylene chloride		75-09-2	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Styrene		100-42-5	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2,2-Tetrachloroethane		79-34-5	5	0.5	U	0.5	U	0.5	U	0.5	U	1	U
Tetrachloroethene (PCE)		127-18-4	5	**0.5**		0.5	U	0.5	U	0.5	U	1	U
Tetrahydrofuran		109-99-9	50*	5	U	--		--		--		--	
Toluene		108-88-3	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113		76-13-1	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,2,3-Trichlorobenzene		87-61-6	5	2.5	U	2.5	U	2.5	U	2.5	U	--	
1,2,4-Trichlorobenzene		120-82-1	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,1-Trichloroethane (TCA)		71-55-6	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
1,1,2-Trichloroethane		79-00-5	1	1.5	U	1.5	U	1.5	U	1.5	U	1	U
Trichloroethene (TCE)		79-01-6	5	**0.61**		**0.32**	**J**	**0.24**	**J**	**0.45**	**J**	1	U
Trichlorofluoromethane (Freon 11)		75-69-4	5	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride		75-01-4	2	**0.47**	**J**	**3.7**		**1.4**		**1.3**		**1.5**	
o-Xylene		95-47-6	5	2.5	U	2.5	U	2.5	U	2.5	U	--	
m/p-Xylene		179601-23-1	5	2.5	U	2.5	U	2.5	U	2.5	U	--	
Total Xylene				2.5	U	2.5	U	2.5	U	2.5	U	2	U
Total VOCs		--	NE	**8.78**	**J**	**7.42**	**J**	**3.94**	**J**	**6.15**	**J**	**3.8**	

Table 4B. Historical Groundwater PFAS Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW1		MW1R		MW2		MW2R	
			Sample Date	2/28/2022		11/30/2023		2/28/2022		12/1/2023	
			Parent Sample								
Analyte	**Units**	**CAS No.**	**NYS AWQS**								
PFAS	ng/L										
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)		763051-92-9	NE	--		--		--		--	
2H,2H,3H,3H-Perfluorooctanoic Acid (5:3FTCA)		914637-49-3	NE	--		--		--		--	
3-Perfluoroheptyl Propanoic Acid (7:3FTCA)		812-70-4	NE	--		--		--		--	
3-Perfluoropropyl Propanoic Acid (3:3FTCA)		356-02-5	NE	--		--		--		--	
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)		919005-14-4	NE	--		--		--		--	
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)		756426-58-1	NE	--		--		--		--	
N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA)		4151-50-2	NE	--		--		--		--	
N-Ethyl Perfluorooctanesulfonamido Ethanol (NEtFOSE)		1691-99-2	NE	--		--		--		--	
N-ethyl perfluorooctanesulfonamidoacetic acid (NEtFOSAA)		2991-50-6	NE	1.71	U	1.88	U	1.84	U	1.88	U
N-Methyl Perfluorooctane Sulfonamide (NMeFOSA)		31506-32-8	NE	--		--		--		--	
N-Methyl Perfluorooctanesulfonamido Ethanol (NMeFOSE)		24448-09-7	NE	--		--		--		--	
N-methyl perfluorooctanesulfonamidoacetic acid (NMeFOSAA)		2355-31-9	NE	1.71	U	1.88	U	1.84	U	1.88	U
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)		151772-58-6	NE	--		--		--		--	
Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA)		113507-82-7	NE	--		--		--		--	
Perfluoro(2-Propoxypropanoic) Acid		13252-13-6	NE	--		--		--		--	
Perfluoro-3-methoxypropanoic acid (PFMPA)		377-73-1	NE	--		--		--		--	
Perfluoro-4-methoxybutanoic acid (PFMBA)		863090-89-5	NE	--		--		--		--	
Perfluorobutanesulfonic acid (PFBS)		375-73-5	NE	**7.01**		**1.17**	**J**	**0.646**	**J**	**1.14**	**J**
Perfluorobutanoic acid (PFBA)		375-22-4	NE	**15.8**		**10.7**		1.84	U	**4.26**	
Perfluorodecanesulfonic acid (PFDS)		335-77-3	NE	1.71	U	1.88	U	1.84	U	1.88	U
Perfluorodecanoic acid (PFDA)		335-76-2	NE	1.71	U	1.88	U	**0.41**	**J**	1.88	U
Perfluorododecanesulfonic Acid (PFDoS)		79780-39-5	NE	--		--		--		--	
Perfluorododecanoic acid (PFDoA)		307-55-1	NE	1.71	U	1.88	U	1.84	U	1.88	U
Perfluoroheptanesulfonic acid (PFHpS)		375-92-8	NE	**1.72**		**0.773**	**J**	1.84	U	1.88	U
Perfluoroheptanoic acid (PFHpA)		375-85-9	NE	**18.8**		**3.34**		**0.933**	**J**	**3.03**	
Perfluorohexanoic acid (PFHxA)		307-24-4	NE	**21.4**	**J**	**10.2**		**2.26**		**4.66**	
Perfluorononanesulfonic Acid (Pfns)		68259-12-1	NE	--		--		--		--	
Perfluorooctanesulfonamide (FOSA)		754-91-6	NE	1.71	U	1.88	U	1.84	U	1.88	U
Perfluoropentanesulfonic Acid (Pfpes)		2706-91-4	NE	--		--		--		--	
Perfluoropentanoic Acid (PFPeA)		2706-90-3	NE	**44.6**	**J**	**12.5**		**2.25**		**6.14**	
Perfluorotetradecanoic acid (PFTA/PFTeDA)		376-06-7	NE	1.71	U	1.88	U	1.84	U	1.88	U
Perfluorotridecanoic acid (PFTriA/PFTrDA)		72629-94-8	NE	1.71	U	1.88	U	1.84	U	1.88	U
Perfluoroundecanoic acid (PFUnA)		2058-94-8	NE	1.71	U	1.88	U	1.84	U	1.88	U
Sodium 1H,1H,2H,2H-Perfluorodecane Sulfonate (8:2)		39108-34-4	NE	1.71	U	1.88	U	1.84	U	1.88	U
Sodium 1H,1H,2H,2H-Perfluorohexane Sulfonate (4:2)		757124-72-4	NE	--		--		--		--	
Sodium 1H,1H,2H,2H-Perfluorooctane Sulfonate (6:2)		27619-97-2	NE	1.71	U	1.88	U	1.84	U	**1.25**	**J**
Perfluorohexanesulfonic acid (PFHxS)		355-46-4	NE	**12.6**		**1.29**	**J**	1.84	U	**2.8**	
Perfluorononanoic Acid (PFNA)		375-95-1	NE	**2.16**		**2.52**		**0.376**	**J**	**1.08**	**J**
Perfluorooctanesulfonic acid (PFOS)		1763-23-1	2.7	**61.8**	**J**	**22.4**		**2.21**	**J**	**17.9**	
Perfluorooctanoic Acid (PFOA)		335-67-1	6.7	**30.7**	**J**	**6.32**		**0.886**	**J**	**13**	
PFOA/PFOS, Total		N/A	NE	**92.5**		**28.7**		**3.1**	**J**	**30.9**	

Table 4B. Historical Groundwater PFAS Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	TW6		DUP		MW3S		MW3S		MW3S		MW3S		MW3S	
			Sample Date	2/27/2022		2/27/2022		12/1/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
			Parent Sample			TW6											
Analyte	**Units**	**CAS No.**	**NYS AWQS**														
PFAS	ng/L																
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)		763051-92-9	NE	--		--		--		6.26	U	32	U	32	U	1.7	U
2H,2H,3H,3H-Perfluorooctanoic Acid (5:3FTCA)		914637-49-3	NE	--		--		--		39.1	U	200	U	200	U	8.3	U
3-Perfluoroheptyl Propanoic Acid (7:3FTCA)		812-70-4	NE	--		--		--		39.1	U	200	U	200	U	8.3	U
3-Perfluoropropyl Propanoic Acid (3:3FTCA)		356-02-5	NE	--		--		--		7.83	U	40	U	40	U	3.3	U
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)		919005-14-4	NE	--		--		--		6.26	U	32	U	32	U	1.7	U
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)		756426-58-1	NE	--		--		--		6.26	U	32	U	32	U	1.7	U
N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA)		4151-50-2	NE	--		--		--		1.56	U	8	U	8	U	1.7	U
N-Ethyl Perfluorooctanesulfonamido Ethanol (NEtFOSE)		1691-99-2	NE	--		--		--		15.6	U	80	U	80	U	8.2	U
N-ethyl perfluorooctanesulfonamidoacetic acid (NEtFOSAA)		2991-50-6	NE	1.83	U	1.82	U	2	U	1.56	U	8	U	8	U	1.7	U
N-Methyl Perfluorooctane Sulfonamide (NMeFOSA)		31506-32-8	NE	--		--		--		1.56	U	8	U	8	U	1.7	U
N-Methyl Perfluorooctanesulfonamido Ethanol (NMeFOSE)		24448-09-7	NE	--		--		--		15.6	U	80	U	80	U	8.2	U
N-methyl perfluorooctanesulfonamidoacetic acid (NMeFOSAA)		2355-31-9	NE	**6.07**	**J**	**6.14**	**J**	2	U	**1.27**	**J**	8	U	8	U	1.7	U
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)		151772-58-6	NE	--		--		--		3.13	U	16	U	16	U	1.7	U
Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA)		113507-82-7	NE	--		--		--		3.13	U	16	U	16	U	1.7	U
Perfluoro(2-Propoxypropanoic) Acid		13252-13-6	NE	--		--		--		6.26	U	32	U	32	U	1.7	U
Perfluoro-3-methoxypropanoic acid (PFMPA)		377-73-1	NE	--		--		--		3.13	U	16	U	**5.64**	**J**	1.7	U
Perfluoro-4-methoxybutanoic acid (PFMBA)		863090-89-5	NE	--		--		--		3.13	U	16	U	16	U	1.7	U
Perfluorobutanesulfonic acid (PFBS)		375-73-5	NE	**6.12**		**5.77**		**4.58**		**11.8**	**F**	**13.3**	**F**	**6.72**	**J**	**4.5**	
Perfluorobutanoic acid (PFBA)		375-22-4	NE	**13**		**13.3**		**14.6**		6.26	U	**18.4**	**J**	**19.2**	**J**	**5.1**	
Perfluorodecanesulfonic acid (PFDS)		335-77-3	NE	1.83	U	1.82	U	2	U	1.56	U	8	U	8	U	1.7	U
Perfluorodecanoic acid (PFDA)		335-76-2	NE	**2.9**		**3.2**		2	U	1.56	U	8	U	8	U	1.7	U
Perfluorododecanesulfonic Acid (PFDoS)		79780-39-5	NE	--		--		--		1.56	U	8	U	8	U	1.7	U
Perfluorododecanoic acid (PFDoA)		307-55-1	NE	**0.539**	**J**	**0.505**	**J**	2	U	1.56	U	8	U	8	U	1.7	U
Perfluoroheptanesulfonic acid (PFHpS)		375-92-8	NE	**1.95**		**2.22**		**3.05**	**J**	**0.689**	**J**	8	U	8	U	1.7	U
Perfluoroheptanoic acid (PFHpA)		375-85-9	NE	**16.2**		**16.2**		**8.5**		**10.3**	**J**	**11**		**7.64**	**J**	**5.7**	
Perfluorohexanoic acid (PFHxA)		307-24-4	NE	**30.9**		**30.9**		**13.6**		**14**		**21.1**		**14.2**		**12**	
Perfluorononanesulfonic Acid (Pfns)		68259-12-1	NE	--		--		--		1.56	U	8	U	8	U	1.7	U
Perfluorooctanesulfonamide (FOSA)		754-91-6	NE	**1.24**	**J**	**1.76**	**J**	2	U	1.56	U	8	U	8	U	1.7	U
Perfluoropentanesulfonic Acid (Pfpes)		2706-91-4	NE	--		--		--		**2.14**		**2.56**	**J**	**2.08**	**J**	**1.7**	
Perfluoropentanoic Acid (PFPeA)		2706-90-3	NE	**49.4**		**48.8**		**20.5**		**19**		**24**		**14.7**	**J**	**7.3**	
Perfluorotetradecanoic acid (PFTA/PFTeDA)		376-06-7	NE	1.83	U	1.82	U	2	U	1.56	U	8	U	8	U	1.6	U
Perfluorotridecanoic acid (PFTriA/PFTrDA)		72629-94-8	NE	1.83	U	1.82	U	2	U	1.56	U	8	U	8	U	1.6	U
Perfluoroundecanoic acid (PFUnA)		2058-94-8	NE	**1.04**	**J**	**0.978**	**J**	2	U	1.56	U	8	U	8	U	1.7	U
Sodium 1H,1H,2H,2H-Perfluorodecane Sulfonate (8:2)		39108-34-4	NE	1.83	U	1.82	U	2	U	6.26	U	32	U	32	U	3.3	U
Sodium 1H,1H,2H,2H-Perfluorohexane Sulfonate (4:2)		757124-72-4	NE	--		--		--		6.26	U	32	U	32	U	3.3	U
Sodium 1H,1H,2H,2H-Perfluorooctane Sulfonate (6:2)		27619-97-2	NE	1.83	U	1.82	U	**1.57**	**J**	6.26	U	32	U	32	U	3.3	U
Perfluorohexanesulfonic acid (PFHxS)		355-46-4	NE	**10.1**		**10.3**		**12.4**	**J**	**6.48**		**7.08**	**J**	**3.92**	**J**	**3.4**	
Perfluorononanoic Acid (PFNA)		375-95-1	NE	**4.64**		**4.98**		2	U	**0.697**	**J**	**1.52**	**J**	8	U	1.7	U
Perfluorooctanesulfonic acid (PFOS)		1763-23-1	2.7	**118**		**122**		**10.1**		**19.1**		**24.4**		**10.4**		**6.6**	**I**
Perfluorooctanoic Acid (PFOA)		335-67-1	6.7	**34.1**		**35**		**17.2**		**32.2**		**30.4**		**18.3**		**15**	
PFOA/PFOS, Total		N/A	NE	**152**		**157**		**27.3**		**51.3**		**54.8**		**28.7**		**21.6**	**I**

Table 4B. Historical Groundwater PFAS Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW3I		MW3I		MW3I		MW3I		MW3I	
			Sample Date	11/30/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
			Parent Sample										
Analyte	**Units**	**CAS No.**	**NYS AWQS**										
PFAS	ng/L												
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)		763051-92-9	NE	--		6.53	U	6.24	U	6.5	U	1.6	U
2H,2H,3H,3H-Perfluorooctanoic Acid (5:3FTCA)		914637-49-3	NE	--		40.8	U	39	U	40.6	U	7.8	U
3-Perfluoroheptyl Propanoic Acid (7:3FTCA)		812-70-4	NE	--		40.8	U	39	U	40.6	U	7.8	U
3-Perfluoropropyl Propanoic Acid (3:3FTCA)		356-02-5	NE	--		8.16	U	7.8	U	8.13	U	3.1	U
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)		919005-14-4	NE	--		6.53	U	6.24	U	6.5	U	1.6	U
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)		756426-58-1	NE	--		6.53	U	6.24	U	6.5	U	1.6	U
N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA)		4151-50-2	NE	--		1.63	U	1.56	U	1.62	U	1.6	U
N-Ethyl Perfluorooctanesulfonamido Ethanol (NEtFOSE)		1691-99-2	NE	--		16.3	U	15.6	U	16.2	U	7.8	U
N-ethyl perfluorooctanesulfonamidoacetic acid (NEtFOSAA)		2991-50-6	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
N-Methyl Perfluorooctane Sulfonamide (NMeFOSA)		31506-32-8	NE	--		1.63	U	1.56	U	1.62	U	1.6	U
N-Methyl Perfluorooctanesulfonamido Ethanol (NMeFOSE)		24448-09-7	NE	--		16.3	U	15.6	U	16.2	U	7.8	U
N-methyl perfluorooctanesulfonamidoacetic acid (NMeFOSAA)		2355-31-9	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)		151772-58-6	NE	--		3.26	U	3.12	U	3.25	U	1.6	U
Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA)		113507-82-7	NE	--		3.26	U	3.12	U	3.25	U	1.6	U
Perfluoro(2-Propoxypropanoic) Acid		13252-13-6	NE	--		6.53	U	6.24	U	6.5	U	1.6	U
Perfluoro-3-methoxypropanoic acid (PFMPA)		377-73-1	NE	--		3.26	U	3.12	U	3.25	U	1.6	U
Perfluoro-4-methoxybutanoic acid (PFMBA)		863090-89-5	NE	--		3.26	U	3.12	U	3.25	U	1.6	U
Perfluorobutanesulfonic acid (PFBS)		375-73-5	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Perfluorobutanoic acid (PFBA)		375-22-4	NE	**1.59**	**J**	**2.23**	**J**	**2.28**	**J**	6.5	U	**0.97**	**J**
Perfluorodecanesulfonic acid (PFDS)		335-77-3	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Perfluorodecanoic acid (PFDA)		335-76-2	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Perfluorododecanesulfonic Acid (PFDoS)		79780-39-5	NE	--		1.63	U	1.56	U	1.62	U	1.6	U
Perfluorododecanoic acid (PFDoA)		307-55-1	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Perfluoroheptanesulfonic acid (PFHpS)		375-92-8	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Perfluoroheptanoic acid (PFHpA)		375-85-9	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Perfluorohexanoic acid (PFHxA)		307-24-4	NE	1.87	U	1.63	U	**0.39**	**JF**	1.62	U	1.6	U
Perfluorononanesulfonic Acid (Pfns)		68259-12-1	NE	--		1.63	U	1.56	U	1.62	U	1.6	U
Perfluorooctanesulfonamide (FOSA)		754-91-6	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Perfluoropentanesulfonic Acid (Pfpes)		2706-91-4	NE	--		1.63	U	1.56	U	1.62	U	1.6	U
Perfluoropentanoic Acid (PFPeA)		2706-90-3	NE	1.87	U	**0.53**	**J**	**0.764**	**J**	3.25	U	1.6	U
Perfluorotetradecanoic acid (PFTA/PFTeDA)		376-06-7	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Perfluorotridecanoic acid (PFTriA/PFTrDA)		72629-94-8	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Perfluoroundecanoic acid (PFUnA)		2058-94-8	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Sodium 1H,1H,2H,2H-Perfluorodecane Sulfonate (8:2)		39108-34-4	NE	1.87	U	6.53	U	6.24	U	6.5	U	3.1	U
Sodium 1H,1H,2H,2H-Perfluorohexane Sulfonate (4:2)		757124-72-4	NE	--		6.53	U	6.24	U	6.5	U	3.1	U
Sodium 1H,1H,2H,2H-Perfluorooctane Sulfonate (6:2)		27619-97-2	NE	**1.48**	**J**	6.53	U	6.24	U	6.5	U	3.1	U
Perfluorohexanesulfonic acid (PFHxS)		355-46-4	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Perfluorononanoic Acid (PFNA)		375-95-1	NE	1.87	U	1.63	U	1.56	U	1.62	U	1.6	U
Perfluorooctanesulfonic acid (PFOS)		1763-23-1	2.7	1.87	U	1.63	U	**3.49**		**0.626**	**J**	1.6	U
Perfluorooctanoic Acid (PFOA)		335-67-1	6.7	1.87	U	1.63	U	**0.288**	**J**	1.62	U	1.6	U
PFOA/PFOS, Total		N/A	NE	1.87	U	1.63	U	**3.778**	J	**0.626**	**J**	1.6	U

Table 4B. Historical Groundwater PFAS Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW4S		MW4S		MW4S		MW4S		DUP-061725		MW4S	
			Sample Date	12/1/2023		12/10/2024		3/12/2025		6/17/2025		6/17/2025		9/24/2025	
			Parent Sample									MW4S			
Analyte	**Units**	**CAS No.**	**NYS AWQS**												
PFAS	ng/L														
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)		763051-92-9	NE	--		5.76	U	6.02	U	6.03	U	6.41	U	1.5	U
2H,2H,3H,3H-Perfluorooctanoic Acid (5:3FTCA)		914637-49-3	NE	--		36	U	37.6	U	37.7	U	40	U	7.7	U
3-Perfluoroheptyl Propanoic Acid (7:3FTCA)		812-70-4	NE	--		36	U	37.6	U	37.7	U	40	U	7.7	U
3-Perfluoropropyl Propanoic Acid (3:3FTCA)		356-02-5	NE	--		7.2	U	7.52	U	7.54	U	8.01	U	3.1	U
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)		919005-14-4	NE	--		5.76	U	6.02	U	6.03	U	6.41	U	1.5	U
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)		756426-58-1	NE	--		5.76	U	6.02	U	6.03	U	6.41	U	1.5	U
N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA)		4151-50-2	NE	--		1.44	U	1.5	U	1.51	U	1.6	U	1.5	U
N-Ethyl Perfluorooctanesulfonamido Ethanol (NEtFOSE)		1691-99-2	NE	--		14.4	U	15	U	15.1	U	16	U	7.7	U
N-ethyl perfluorooctanesulfonamidoacetic acid (NEtFOSAA)		2991-50-6	NE	1.94	U	1.44	U	1.5	U	1.51	U	1.6	U	1.5	U
N-Methyl Perfluorooctane Sulfonamide (NMeFOSA)		31506-32-8	NE	--		1.44	U	1.5	U	1.51	U	1.6	U	1.5	U
N-Methyl Perfluorooctanesulfonamido Ethanol (NMeFOSE)		24448-09-7	NE	--		14.4	U	15	U	15.1	U	16	U	7.7	U
N-methyl perfluorooctanesulfonamidoacetic acid (NMeFOSAA)		2355-31-9	NE	1.94	U	1.44	U	1.5	U	1.51	U	1.6	U	1.5	U
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)		151772-58-6	NE	--		2.88	U	3.01	U	3.01	U	3.2	U	1.5	U
Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA)		113507-82-7	NE	--		2.88	U	3.01	U	3.01	U	3.2	U	1.5	U
Perfluoro(2-Propoxypropanoic) Acid		13252-13-6	NE	--		5.76	U	6.02	U	6.03	U	6.41	U	1.5	U
Perfluoro-3-methoxypropanoic acid (PFMPA)		377-73-1	NE	--		2.88	U	3.01	U	3.01	U	3.2	U	1.5	U
Perfluoro-4-methoxybutanoic acid (PFMBA)		863090-89-5	NE	--		2.88	U	3.01	U	3.01	U	3.2	U	1.5	U
Perfluorobutanesulfonic acid (PFBS)		375-73-5	NE	**4.52**		**4.43**		**5.12**		**2.53**		**3.32**		**4.4**	
Perfluorobutanoic acid (PFBA)		375-22-4	NE	**12.6**		**11.3**		**29.7**		**8.3**		**9.64**		**12**	
Perfluorodecanesulfonic acid (PFDS)		335-77-3	NE	1.94	U	1.44	U	1.5	U	1.51	U	1.6	U	1.5	U
Perfluorodecanoic acid (PFDA)		335-76-2	NE	1.94	U	**3.07**		**7.81**		**6.54**		**6.02**		**1.5**	
Perfluorododecanesulfonic Acid (PFDoS)		79780-39-5	NE	--		1.44	U	1.5	U	1.51	U	1.6	U	1.5	U
Perfluorododecanoic acid (PFDoA)		307-55-1	NE	1.94	U	1.44	U	1.5	U	1.51	U	1.6	U	1.5	U
Perfluoroheptanesulfonic acid (PFHpS)		375-92-8	NE	1.94	U	**0.777**	**J**	**0.459**	**J**	**1.08**	**J**	**1.05**	**J**	1.5	U
Perfluoroheptanoic acid (PFHpA)		375-85-9	NE	**15.9**		**9.42**		**17.5**		**8.07**		**8.62**		**11**	
Perfluorohexanoic acid (PFHxA)		307-24-4	NE	**24.2**		**15.4**		**37.3**		**10.5**		**11.5**		**18**	
Perfluorononanesulfonic Acid (Pfns)		68259-12-1	NE	--		1.44	U	1.5	U	1.51	U	1.6	U	1.5	U
Perfluorooctanesulfonamide (FOSA)		754-91-6	NE	1.94	U	**0.389**	**JF**	**0.293**	**J**	**0.505**	**JF**	**0.521**	**JF**	1.5	U
Perfluoropentanesulfonic Acid (Pfpes)		2706-91-4	NE	--		**0.936**	**J**	**0.579**	**J**	**0.829**	**J**	**0.913**	**J**	**0.8**	**J**
Perfluoropentanoic Acid (PFPeA)		2706-90-3	NE	**36.4**		**26.6**		**67.4**		**12.4**		**15.4**		**25**	
Perfluorotetradecanoic acid (PFTA/PFTeDA)		376-06-7	NE	1.94	U	1.44	U	1.5	U	1.51	U	1.6	U	1.5	U
Perfluorotridecanoic acid (PFTriA/PFTrDA)		72629-94-8	NE	1.94	U	1.44	U	1.5	U	1.51	U	1.6	U	1.5	U
Perfluoroundecanoic acid (PFUnA)		2058-94-8	NE	1.94	U	1.44	U	**0.323**	**J**	1.51	U	**0.216**	**J**	1.5	U
Sodium 1H,1H,2H,2H-Perfluorodecane Sulfonate (8:2)		39108-34-4	NE	1.94	U	5.76	U	6.02	U	6.03	U	6.41	U	3.1	U
Sodium 1H,1H,2H,2H-Perfluorohexane Sulfonate (4:2)		757124-72-4	NE	--		5.76	U	6.02	U	6.03	U	6.41	U	3.1	U
Sodium 1H,1H,2H,2H-Perfluorooctane Sulfonate (6:2)		27619-97-2	NE	1.94	U	5.76	U	6.02	U	6.03	U	6.41	U	3.1	U
Perfluorohexanesulfonic acid (PFHxS)		355-46-4	NE	**4.55**		**4.21**		**2.38**		**4.26**		**3.94**		**3.7**	
Perfluorononanoic Acid (PFNA)		375-95-1	NE	**3.61**		**6.46**		**7.66**		**10.4**		**10.6**		**4.1**	
Perfluorooctanesulfonic acid (PFOS)		1763-23-1	2.7	**12.9**		**45.4**		**31.2**		**75**		**72.1**		**17**	
Perfluorooctanoic Acid (PFOA)		335-67-1	6.7	**37.5**		**30.9**		**26.8**		**21.2**		**23.1**		**24**	
PFOA/PFOS, Total		N/A	NE	**50.4**		**76.3**		**58**		**96.2**		**95.2**		**41**	

Table 4B. Historical Groundwater PFAS Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW4I		DUP-12.1.23		MW4I		DUP-1210202		MW4I		MW4I		MW4I	
			Sample Date	12/1/2023		12/1/2023		12/10/2024		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
			Parent Sample			MW4I				MW 4I							
Analyte	**Units**	**CAS No.**	**NYS AWQS**														
PFAS	ng/L																
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)		763051-92-9	NE	--		--		6.38	U	6.36	U	6.32	U	6.03	U	1.5	U
2H,2H,3H,3H-Perfluorooctanoic Acid (5:3FTCA)		914637-49-3	NE	--		--		39.9	U	39.8	U	39.5	U	37.7	U	7.6	U
3-Perfluoroheptyl Propanoic Acid (7:3FTCA)		812-70-4	NE	--		--		39.9	U	39.8	U	39.5	U	37.7	U	7.6	U
3-Perfluoropropyl Propanoic Acid (3:3FTCA)		356-02-5	NE	--		--		7.97	U	7.96	U	7.9	U	7.54	U	3	U
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)		919005-14-4	NE	--		--		6.38	U	6.36	U	6.32	U	6.03	U	1.5	U
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)		756426-58-1	NE	--		--		6.38	U	6.36	U	6.32	U	6.03	U	1.5	U
N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA)		4151-50-2	NE	--		--		1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
N-Ethyl Perfluorooctanesulfonamido Ethanol (NEtFOSE)		1691-99-2	NE	--		--		15.9	U	15.9	U	15.8	U	15.1	U	7.6	U
N-ethyl perfluorooctanesulfonamidoacetic acid (NEtFOSAA)		2991-50-6	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
N-Methyl Perfluorooctane Sulfonamide (NMeFOSA)		31506-32-8	NE	--		--		1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
N-Methyl Perfluorooctanesulfonamido Ethanol (NMeFOSE)		24448-09-7	NE	--		--		15.9	U	15.9	U	15.8	U	15.1	U	7.6	U
N-methyl perfluorooctanesulfonamidoacetic acid (NMeFOSAA)		2355-31-9	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)		151772-58-6	NE	--		--		3.19	U	3.18	U	3.16	U	3.02	U	1.5	U
Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA)		113507-82-7	NE	--		--		3.19	U	3.18	U	3.16	U	3.02	U	1.5	U
Perfluoro(2-Propoxypropanoic) Acid		13252-13-6	NE	--		--		6.38	U	6.36	U	6.32	U	6.03	U	1.5	U
Perfluoro-3-methoxypropanoic acid (PFMPA)		377-73-1	NE	--		--		3.19	U	3.18	U	3.16	U	3.02	U	1.5	U
Perfluoro-4-methoxybutanoic acid (PFMBA)		863090-89-5	NE	--		--		3.19	U	3.18	U	3.16	U	3.02	U	1.5	U
Perfluorobutanesulfonic acid (PFBS)		375-73-5	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluorobutanoic acid (PFBA)		375-22-4	NE	**1.61**	**J**	**1.63**	**J**	**1.32**	**J**	**1.5**	**J**	**1.86**	**J**	6.03	U	**1.1**	**J**
Perfluorodecanesulfonic acid (PFDS)		335-77-3	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluorodecanoic acid (PFDA)		335-76-2	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluorododecanesulfonic Acid (PFDoS)		79780-39-5	NE	--		--		1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluorododecanoic acid (PFDoA)		307-55-1	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluoroheptanesulfonic acid (PFHpS)		375-92-8	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluoroheptanoic acid (PFHpA)		375-85-9	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluorohexanoic acid (PFHxA)		307-24-4	NE	**0.906**	**J**	**0.916**	**J**	**0.542**	**J**	**0.493**	**J**	**0.41**	**J**	**0.302**	**J**	1.5	U
Perfluorononanesulfonic Acid (Pfns)		68259-12-1	NE	--		--		1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluorooctanesulfonamide (FOSA)		754-91-6	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluoropentanesulfonic Acid (Pfpes)		2706-91-4	NE	--		--		1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluoropentanoic Acid (PFPeA)		2706-90-3	NE	**3.19**		**3.69**		**0.51**	**J**	**0.573**	**J**	**0.41**	**J**	**0.528**	**J**	1.5	U
Perfluorotetradecanoic acid (PFTA/PFTeDA)		376-06-7	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluorotridecanoic acid (PFTriA/PFTrDA)		72629-94-8	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluoroundecanoic acid (PFUnA)		2058-94-8	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Sodium 1H,1H,2H,2H-Perfluorodecane Sulfonate (8:2)		39108-34-4	NE	1.84	U	1.84	U	6.38	U	6.36	U	6.32	U	6.03	U	3	U
Sodium 1H,1H,2H,2H-Perfluorohexane Sulfonate (4:2)		757124-72-4	NE	--		--		6.38	U	6.36	U	6.32	U	6.03	U	3	U
Sodium 1H,1H,2H,2H-Perfluorooctane Sulfonate (6:2)		27619-97-2	NE	**1.82**	**J**	**3.38**		6.38	U	6.36	U	6.32	U	6.03	U	3	U
Perfluorohexanesulfonic acid (PFHxS)		355-46-4	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluorononanoic Acid (PFNA)		375-95-1	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
Perfluorooctanesulfonic acid (PFOS)		1763-23-1	2.7	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	**0.256**	**J**	1.5	U
Perfluorooctanoic Acid (PFOA)		335-67-1	6.7	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	1.51	U	1.5	U
PFOA/PFOS, Total		N/A	NE	1.84	U	1.84	U	1.59	U	1.59	U	1.58	U	**0.256**	**J**	1.5	U

Table 4B. Historical Groundwater PFAS Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW5S		MW5S		MW5S		MW5S		MW5S		MW5I		MW5I	
			Sample Date	11/29/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025		11/29/2023		12/10/2024	
			Parent Sample														
Analyte	**Units**	**CAS No.**	**NYS AWQS**														
PFAS	ng/L																
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)		763051-92-9	NE	--		6.06	U	5.95	U	6.07	U	1.8	U	--		6.54	U
2H,2H,3H,3H-Perfluorooctanoic Acid (5:3FTCA)		914637-49-3	NE	--		37.9	U	37.2	U	37.9	U	8.8	U	--		40.8	U
3-Perfluoroheptyl Propanoic Acid (7:3FTCA)		812-70-4	NE	--		37.9	U	37.2	U	37.9	U	8.8	U	--		40.8	U
3-Perfluoropropyl Propanoic Acid (3:3FTCA)		356-02-5	NE	--		7.58	U	7.44	U	7.58	U	3.5	U	--		8.17	U
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)		919005-14-4	NE	--		6.06	U	5.95	U	6.07	U	1.8	U	--		6.54	U
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)		756426-58-1	NE	--		6.06	U	5.95	U	6.07	U	1.8	U	--		6.54	U
N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA)		4151-50-2	NE	--		1.52	U	1.49	U	1.52	U	1.8	U	--		1.63	U
N-Ethyl Perfluorooctanesulfonamido Ethanol (NEtFOSE)		1691-99-2	NE	--		15.2	U	14.9	U	15.2	U	8.8	U	--		16.3	U
N-ethyl perfluorooctanesulfonamidoacetic acid (NEtFOSAA)		2991-50-6	NE	2.11	U	1.52	U	1.49	U	1.52	U	1.8	U	1.89	U	1.63	U
N-Methyl Perfluorooctane Sulfonamide (NMeFOSA)		31506-32-8	NE	--		1.52	U	1.49	U	1.52	U	1.8	U	--		1.63	U
N-Methyl Perfluorooctanesulfonamido Ethanol (NMeFOSE)		24448-09-7	NE	--		15.2	U	14.9	U	15.2	U	8.8	U	--		16.3	U
N-methyl perfluorooctanesulfonamidoacetic acid (NMeFOSAA)		2355-31-9	NE	2.11	U	1.52	U	1.49	U	1.52	U	1.8	U	1.89	U	1.63	U
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)		151772-58-6	NE	--		3.03	U	2.98	U	3.03	U	1.8	U	--		3.27	U
Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA)		113507-82-7	NE	--		3.03	U	2.98	U	3.03	U	1.8	U	--		3.27	U
Perfluoro(2-Propoxypropanoic) Acid		13252-13-6	NE	--		6.06	U	5.95	U	6.07	U	1.8	U	--		6.54	U
Perfluoro-3-methoxypropanoic acid (PFMPA)		377-73-1	NE	--		3.03	U	2.98	U	6.07	U	1.8	U	--		3.27	U
Perfluoro-4-methoxybutanoic acid (PFMBA)		863090-89-5	NE	--		3.03	U	2.98	U	3.03	U	1.8	U	--		3.27	U
Perfluorobutanesulfonic acid (PFBS)		375-73-5	NE	**4.84**		**5.82**		**3.36**		**2.8**		**0.85**	**J**	1.89	U	1.63	U
Perfluorobutanoic acid (PFBA)		375-22-4	NE	**18.1**		**12**		**9.82**		**10.7**	**J**	**9.1**		**1.46**	**J**	**2.01**	**J**
Perfluorodecanesulfonic acid (PFDS)		335-77-3	NE	2.11	U	1.52	U	1.49	U	1.52	U	1.8	U	1.89	U	1.63	U
Perfluorodecanoic acid (PFDA)		335-76-2	NE	**0.468**	**J**	**0.833**	**J**	**0.439**	**J**	**0.531**	**J**	1.8	U	1.89	U	1.63	U
Perfluorododecanesulfonic Acid (PFDoS)		79780-39-5	NE	--		1.52	U	1.49	U	1.52	U	1.8	U	--		1.63	U
Perfluorododecanoic acid (PFDoA)		307-55-1	NE	2.11	U	1.52	U	1.49	U	1.52	U	1.8	U	1.89	U	1.63	U
Perfluoroheptanesulfonic acid (PFHpS)		375-92-8	NE	**1.3**	**J**	**1.28**	**J**	**1.49**		**0.872**	**J**	1.8	U	1.89	U	1.63	U
Perfluoroheptanoic acid (PFHpA)		375-85-9	NE	**13.3**		**7.98**		**7.88**		**8.73**		**2.6**		1.89	U	**0.31**	**J**
Perfluorohexanoic acid (PFHxA)		307-24-4	NE	**25.2**		**12.8**		**13.1**		**12.2**		**5**		**0.43**	**J**	**0.899**	**J**
Perfluorononanesulfonic Acid (Pfns)		68259-12-1	NE	--		1.52	U	1.49	U	1.52	U	1.8	U	--		1.63	U
Perfluorooctanesulfonamide (FOSA)		754-91-6	NE	2.11	U	**0.788**	**JF**	**0.566**	**J**	**0.508**	**J**	**0.48**	**J**	1.89	U	1.63	U
Perfluoropentanesulfonic Acid (Pfpes)		2706-91-4	NE	--		**1.1**	**J**	**0.618**	**J**	**0.516**	**J**	1.8	U	--		1.63	U
Perfluoropentanoic Acid (PFPeA)		2706-90-3	NE	**37.4**		**22.2**		**22.5**		**17.8**		**8.1**		**0.472**	**J**	**0.596**	**J**
Perfluorotetradecanoic acid (PFTA/PFTeDA)		376-06-7	NE	2.11	U	1.52	U	1.49	U	1.52	U	1.8	U	1.89	U	1.63	U
Perfluorotridecanoic acid (PFTriA/PFTrDA)		72629-94-8	NE	2.11	U	1.52	U	1.49	U	1.52	U	1.8	U	1.89	U	1.63	U
Perfluoroundecanoic acid (PFUnA)		2058-94-8	NE	2.11	U	1.52	U	1.49	U	1.52	U	1.8	U	1.89	U	1.63	U
Sodium 1H,1H,2H,2H-Perfluorodecane Sulfonate (8:2)		39108-34-4	NE	2.11	U	6.06	U	5.95	U	6.07	U	3.5	U	1.89	U	6.54	U
Sodium 1H,1H,2H,2H-Perfluorohexane Sulfonate (4:2)		757124-72-4	NE	--		6.06	U	5.95	U	6.07	U	3.5	U	--		6.54	U
Sodium 1H,1H,2H,2H-Perfluorooctane Sulfonate (6:2)		27619-97-2	NE	**1.95**	**J**	6.06	U	5.95	U	6.07	U	3.5	U	**42.7**	**J**	**101**	
Perfluorohexanesulfonic acid (PFHxS)		355-46-4	NE	**7.38**		**7.64**		**6.13**		**4.48**		**1.4**	**J**	**0.483**	**J**	**0.784**	**J**
Perfluorononanoic Acid (PFNA)		375-95-1	NE	**4.32**		**2.17**		**1.47**	**J**	**1.75**		1.8	U	1.89	U	1.63	U
Perfluorooctanesulfonic acid (PFOS)		1763-23-1	2.7	**32.2**		**41.4**		**37.9**		**36.3**		**20**		1.89	U	**0.531**	**J**
Perfluorooctanoic Acid (PFOA)		335-67-1	6.7	**91.5**		**46.4**		**37.4**		**38.8**		**14**		**0.257**	**J**	**0.343**	**J**
PFOA/PFOS, Total		N/A	NE	**124**		**87.8**		**75.3**		**75.1**		**34**		**0.257**	**J**	**0.874**	**J**

Table 4B. Historical Groundwater PFAS Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

			Sample Name	MW5I		MW5I		MW5I		MW6		MW7		MW8		MW11		MW12S	
			Sample Date	3/12/2025		6/17/2025		9/24/2025		11/28/2023		11/28/2023		11/29/2023		11/29/2023		11/30/2023	
			Parent Sample																
Analyte	**Units**	**CAS No.**	**NYS AWQS**																
PFAS	ng/L																		
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)		763051-92-9	NE	6.27	U	5.98	U	1.6	U	--		--		--		--		--	
2H,2H,3H,3H-Perfluorooctanoic Acid (5:3FTCA)		914637-49-3	NE	39.2	U	37.4	U	8.2	U	--		--		--		--		--	
3-Perfluoroheptyl Propanoic Acid (7:3FTCA)		812-70-4	NE	39.2	U	37.4	U	8.2	U	--		--		--		--		--	
3-Perfluoropropyl Propanoic Acid (3:3FTCA)		356-02-5	NE	7.83	U	7.47	U	3.3	U	--		--		--		--		--	
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)		919005-14-4	NE	6.27	U	5.98	U	1.6	U	--		--		--		--		--	
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)		756426-58-1	NE	6.27	U	5.98	U	1.6	U	--		--		--		--		--	
N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA)		4151-50-2	NE	1.57	U	1.49	U	1.6	U	--		--		--		--		--	
N-Ethyl Perfluorooctanesulfonamido Ethanol (NEtFOSE)		1691-99-2	NE	15.7	U	14.9	U	8.2	U	--		--		--		--		--	
N-ethyl perfluorooctanesulfonamidoacetic acid (NEtFOSAA)		2991-50-6	NE	1.57	U	1.49	U	1.6	U	1.93	U	1.84	U	1.89	U	1.86	U	1.87	U
N-Methyl Perfluorooctane Sulfonamide (NMeFOSA)		31506-32-8	NE	1.57	U	1.49	U	1.6	U	--		--		--		--		--	
N-Methyl Perfluorooctanesulfonamido Ethanol (NMeFOSE)		24448-09-7	NE	15.7	U	14.9	U	8.2	U	--		--		--		--		--	
N-methyl perfluorooctanesulfonamidoacetic acid (NMeFOSAA)		2355-31-9	NE	1.57	U	1.49	U	1.6	U	1.93	U	1.84	U	1.89	U	1.86	U	1.87	U
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)		151772-58-6	NE	3.13	U	2.99	U	1.6	U	--		--		--		--		--	
Perfluoro(2-ethoxyethane)sulfonic acid (PFEESA)		113507-82-7	NE	3.13	U	2.99	U	1.6	U	--		--		--		--		--	
Perfluoro(2-Propoxypropanoic) Acid		13252-13-6	NE	6.27	U	5.98	U	1.6	U	--		--		--		--		--	
Perfluoro-3-methoxypropanoic acid (PFMPA)		377-73-1	NE	3.13	U	2.99	U	1.6	U	--		--		--		--		--	
Perfluoro-4-methoxybutanoic acid (PFMBA)		863090-89-5	NE	3.13	U	2.99	U	1.6	U	--		--		--		--		--	
Perfluorobutanesulfonic acid (PFBS)		375-73-5	NE	1.57	U	1.49	U	1.6	U	1.93	U	1.84	U	**2.91**		**1.16**	**J**	**2.48**	
Perfluorobutanoic acid (PFBA)		375-22-4	NE	**2.19**	**J**	**2.03**	**J**	**1.7**	**J**	**1.31**	**J**	**2.16**		**10.8**		**3.06**		**5.65**	
Perfluorodecanesulfonic acid (PFDS)		335-77-3	NE	1.57	U	1.49	U	1.6	U	1.93	U	1.84	U	1.89	U	1.86	U	1.87	U
Perfluorodecanoic acid (PFDA)		335-76-2	NE	1.57	U	1.49	U	1.6	U	1.93	U	1.84	U	1.89	U	1.86	U	1.87	U
Perfluorododecanesulfonic Acid (PFDoS)		79780-39-5	NE	1.57	U	1.49	U	1.6	U	--		--		--		--		--	
Perfluorododecanoic acid (PFDoA)		307-55-1	NE	1.57	U	1.49	U	1.6	U	1.93	U	1.84	U	1.89	U	1.86	U	1.87	U
Perfluoroheptanesulfonic acid (PFHpS)		375-92-8	NE	1.57	U	1.49	U	1.6	U	1.93	U	1.84	U	1.89	U	1.86	U	1.87	U
Perfluoroheptanoic acid (PFHpA)		375-85-9	NE	1.57	U	**0.471**	**J**	**1.5**	**J**	1.93	U	1.84	U	**8.44**		**1.86**	**J**	**3.02**	
Perfluorohexanoic acid (PFHxA)		307-24-4	NE	**0.407**	**J**	**1.41**	**J**	**4.4**		1.93	U	**1**	**J**	**12.7**		**2.69**		**5.11**	
Perfluorononanesulfonic Acid (Pfns)		68259-12-1	NE	1.57	U	1.49	U	1.6	U	--		--		--		--		--	
Perfluorooctanesulfonamide (FOSA)		754-91-6	NE	1.57	U	1.49	U	1.6	U	1.93	U	1.84	U	1.89	U	1.86	U	1.87	U
Perfluoropentanesulfonic Acid (Pfpes)		2706-91-4	NE	1.57	U	1.49	U	1.6	U	--		--		--		--		--	
Perfluoropentanoic Acid (PFPeA)		2706-90-3	NE	**0.47**	**J**	**1.33**	**J**	**4.9**		**0.717**	**J**	**1.73**	**J**	**21**		**3.47**		**6.97**	
Perfluorotetradecanoic acid (PFTA/PFTeDA)		376-06-7	NE	1.57	U	1.49	U	1.6	U	1.93	U	1.84	U	1.89	U	1.86	U	1.87	U
Perfluorotridecanoic acid (PFTriA/PFTrDA)		72629-94-8	NE	1.57	U	1.49	U	1.6	U	1.93	U	1.84	U	1.89	U	1.86	U	1.87	U
Perfluoroundecanoic acid (PFUnA)		2058-94-8	NE	1.57	U	1.49	U	1.6	U	1.93	U	1.84	U	1.89	U	1.86	U	1.87	U
Sodium 1H,1H,2H,2H-Perfluorodecane Sulfonate (8:2)		39108-34-4	NE	6.27	U	5.98	U	3.3	U	1.93	U	1.84	U	1.89	U	1.86	U	1.87	U
Sodium 1H,1H,2H,2H-Perfluorohexane Sulfonate (4:2)		757124-72-4	NE	6.27	U	5.98	U	3.3	U	--		--		--		--		--	
Sodium 1H,1H,2H,2H-Perfluorooctane Sulfonate (6:2)		27619-97-2	NE	**39.2**		**25.8**		**39**		1.93	U	1.84	U	1.89	U	1.86	U	1.87	U
Perfluorohexanesulfonic acid (PFHxS)		355-46-4	NE	**0.462**	**J**	**0.508**	**J**	**1.1**	**J**	1.93	U	1.84	U	1.89	U	**0.63**	**J**	**0.998**	**J**
Perfluorononanoic Acid (PFNA)		375-95-1	NE	1.57	U	1.49	U	1.6	U	1.93	U	1.84	U	**0.662**	**J**	**1.24**	**J**	**1.2**	**J**
Perfluorooctanesulfonic acid (PFOS)		1763-23-1	2.7	**0.509**	**J**	**0.448**	**J**	1.6	U	1.93	U	1.84	U	1.89	U	**9.53**		**10.3**	
Perfluorooctanoic Acid (PFOA)		335-67-1	6.7	1.57	U	**0.247**	**J**	1.6	U	1.93	U	1.84	U	**8.09**		**3.95**		**5.35**	**J**
PFOA/PFOS, Total		N/A	NE	**0.509**	**J**	**0.695**	J	1.6	U	1.93	U	1.84	U	**8.09**		**13.5**		**15.7**	

Table 4C. Historical Groundwater Natural Attenuation Parameter Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

Analyte	Units	CAS No.	MW1 11/22/2021		MW1 2/28/2022		MW1_G 2/7/2023		MW1R 7/6/2023		MW1R 11/30/2023		MW1R 12/11/2024		MW1R 3/12/2025		MW1R 9/24/2025	
Dissolved Gases	ug/L																	
Ethane		74-84-0	--		--		**4.38**		**3.69**		**0.89**		**5**	**J**	**3.9**	**J**	**5.2**	**J**
Ethene		74-85-1	--		--		0.5	U	**3.82**		**4.05**		17	U	17	U	7	U
Methane		74-82-8	--		--		**76.5**		**182**		**203**		**970**		**400**		**1,800**	
General Chemistry																		
DO	mg/L	(field parameter)	0.54		0.08		0.98		--		0.81		0.2		0.00		2.51	
Oxidation Reduction Potential	millivolts	(field parameter)	28		101		-65		--		-103.9		-121		-131		-68	
pH	s.u.	(field parameter)	7.6		7.49		7.2		--		7.43		7.52		7.3		7.86	
Sulfate	ug/L	14808-79-8	--		--		**129,000**		**140,000**		**99,000**		**32,400**		--		**30,600**	
Metals	ug/L																	
Iron, Dissolved		7439-89-6	--		--		**21.1**	**J**	**294**	**J**	**418**		**182**		**3,490**		**210**	
Iron, Total		7439-89-6	--		**96.5**		**70,900**		**44,600**		**465**		**1,290**		**4,000**		**610**	
Manganese		7439-96-5	--		**146.2**		--		--		**609.9**		--		--		--	
Volatile Fatty Acids	ug/L																	
Butyric Acid		107-92-6	--		--		4,000	U	500	U	500	U	100,000	U	200,000	U	--	
Lactic Acid		50-21-5	--		--		**3,900**	**J**	**3,100**		**3,000**		--		--		--	
Pyruvic Acid		127-17-3	--		--		4,000	U	500	U	500	U	--		--		--	
Acetic Acid		64-19-7	--		--		**6,700**	**J**	500	U	500	U	100,000	U	200,000	U	--	
Formic acid		64-18-6	--		--		**96,000**		500	U	500	U	--		--		--	
Pentanoic acid		109-52-4	--		--		4,000	U	500	U	500	U	--		--		--	
Propionic Acid		79-09-4	--		--		4,000	U	500	U	500	U	100,000	U	200,000	U	--	
Hexanoic acid		142-62-1	--		--		**1,700**	**J**	500	U	500	U	--		300,000	U	--	

Table 4C. Historical Groundwater Natural Attenuation Parameter Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

Analyte	Units	CAS No.	MW2 2/28/2022		MW2_G 2/7/2023		MW2R 7/6/2023		MW2R 12/1/2023		MW2R 12/11/2024		MW2R 3/12/2025		MW2R 6/17/2025		MW2R 9/24/2025	
Dissolved Gases	ug/L																	
Ethane		74-84-0	--		**1.17**		**0.555**		0.5	U	14	U	14	U	**5.2**		7.5	U
Ethene		74-85-1	--		0.5	U	**0.67**		0.5	U	17	U	17	U	**6.5**		7	U
Methane		74-82-8	--		**30.2**		**72.3**		**51.6**		**230**		**560**		**790**		**890**	
General Chemistry																		
DO	mg/L	(field parameter)	1.08		1.29		--		1.01		0.00		0.39		3.49		0.00	
Oxidation Reduction Potential	millivolts	(field parameter)	-81		-77		--		-51.3		45		118		-134		-85	
pH	s.u.	(field parameter)	10.91		7.25		--		7.19		7.23		7.18		7.37		8.41	
Sulfate	ug/L	14808-79-8	--		**42,300**		**28,000**		**20,000**		**14,500**		--		**16,600**		**27,100**	
Metals	ug/L																	
Iron, Dissolved		7439-89-6	--		50	U	**842**	**J**	**106**		**186**		**51.5**		**2,270**		**1,200**	
Iron, Total		7439-89-6	**1,090**		**130,000**		**1,090**		50	U	**468**		**255**		**4,000**		**1000**	
Manganese		7439-96-5	**27.82**		--		--		**969.1**		--		--		--		--	
Volatile Fatty Acids	ug/L																	
Butyric Acid		107-92-6	--		2,000	U	500	U	5,000	U	100,000	U	200,000	U	--		--	
Lactic Acid		50-21-5	--		**4,900**	**J**	**720**		5,000	U	--		--		--		--	
Pyruvic Acid		127-17-3	--		2,000	U	500	U	5,000	U	--		--		--		--	
Acetic Acid		64-19-7	--		**8,300**		500	U	5,000	U	100,000	U	200,000	U	--		--	
Formic acid		64-18-6	--		**52,000**		500	U	5,000	U	--		--		--		--	
Pentanoic acid		109-52-4	--		2,000	U	500	U	5,000	U	--		--		--		--	
Propionic Acid		79-09-4	--		**850**	**J**	500	U	5,000	U	100,000	U	200,000	U	--		--	
Hexanoic acid		142-62-1	--		**2,300**	**J**	500	U	5,000	U	--		300,000	U	--		--	

Table 4C. Historical Groundwater Natural Attenuation Parameter Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

Analyte	Units	CAS No.	MW3S_G 2/7/2023		MW3S 7/6/2023		MW3S 12/1/2023		MW3S 12/10/2024		MW3S 3/12/2025		MW3S 6/17/2025		MW3S 9/24/2025	
Dissolved Gases	ug/L															
Ethane		74-84-0	**188**		**1,050**		**944**		**1,500**		**1,200**		**560**		**1,100**	
Ethene		74-85-1	**497**		**3,110**		**4,250**		**18,000**		**15,000**		**14,000**		**29,000**	
Methane		74-82-8	**1,820**		**4,630**		**4,510**		**6,900**		**6,300**		**3,500**		**6,800**	
General Chemistry																
DO	mg/L	(field parameter)	--		0.46		0.51		0.00		0.00		0.35		0.05	
Oxidation Reduction Potential	millivolts	(field parameter)	--		-74		-69.7		-39		-91		-129		-104	
pH	s.u.	(field parameter)	--		6.37		6.45		6.48		6.5		6.55		6.54	
Sulfate	ug/L	14808-79-8	**109,000**		**46,000**		**2,000**	**J**	**1,010**		--		**1,160**		**25,000**	
Metals	ug/L															
Iron, Dissolved		7439-89-6	**2,690**		**39,200**	**J**	**106,000**		**78,200**		**83,100**		**120,000**		**88,000**	
Iron, Total		7439-89-6	**104,000**		**44,400**		**98,500**		**72,000**		**84,100**		**91,800**		**91,600**	
Manganese		7439-96-5	--		--		**3,190**		--		--		--		--	
Volatile Fatty Acids	ug/L															
Butyric Acid		107-92-6	**36,000**	**J**	50,000	U	50,000	U	100,000	U	200,000	U	--		--	
Lactic Acid		50-21-5	**6,700**	**J**	50,000	U	50,000	U	--		--		--		--	
Pyruvic Acid		127-17-3	20,000	U	50,000	U	50,000	U	--		--		--		--	
Acetic Acid		64-19-7	**350,000**		**257,000**		**432,000**		**456,000**		**439,000**		--		--	
Formic acid		64-18-6	**100,000**		50,000	U	50,000	U	--		--		--		--	
Pentanoic acid		109-52-4	**7,600**	**J**	50,000	U	50,000	U	--		--		--		--	
Propionic Acid		79-09-4	**730,000**		**641,000**		**540,000**		**315,000**		**319,000**		--		--	
Hexanoic acid		142-62-1	20,000	U	50,000	U	50,000	U	--		300,000	U	--		--	

Table 4C. Historical Groundwater Natural Attenuation Parameter Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

Analyte	Units	CAS No.	MW3I_G 2/8/2023		MW3I 7/6/2023		MW3I 11/30/2023		MW3I 12/10/2024		MW3I 3/12/2025		MW3I 6/17/2025		MW3I 9/24/2025	
Dissolved Gases	ug/L															
Ethane		74-84-0	**15.4**		**4.67**		**1.31**		**1.9**	**J**	**3.3**	**J**	**5.2**		7.5	U
Ethene		74-85-1	**44.1**		**9.45**		**1.11**		**10**	**J**	17	U	**6**		7	U
Methane		74-82-8	**128**		**748**		**277**		**310**		**71**		**31**		**140**	
General Chemistry																
DO	mg/L	(field parameter)	--		0.48		0.27		0.00		0.00		0.29		0.00	
Oxidation Reduction Potential	millivolts	(field parameter)	--		-116		-134		-70		-45		-93		-108	
pH	s.u.	(field parameter)	--		7.11		7.15		7.2		7.18		7.24		7.17	
Sulfate	ug/L	14808-79-8	**73,600**		**120,000**		**78,000**		**79,000**		--		**57,800**		**79,900**	
Metals	ug/L															
Iron, Dissolved		7439-89-6	**21.5**	**J**	**1,910**	**J**	**1,170**		**1,370**		**496**		**554**		**1,000**	
Iron, Total		7439-89-6	**195,000**		**2,500**		**1,230**		**1,370**		**607**		**709**		**1,100**	
Manganese		7439-96-5	--		--		**47.42**		--		--		--		--	
Volatile Fatty Acids	ug/L															
Butyric Acid		107-92-6	**6,600**		500	U	5,000	U	100,000	U	200,000	U	--		--	
Lactic Acid		50-21-5	**3,900**	**J**	**670**		5,000	U	--		--		--		--	
Pyruvic Acid		127-17-3	**750**	**J**	500	U	5,000	U	--		--		--		--	
Acetic Acid		64-19-7	**31,000**		**2,300**		5,000	U	100,000	U	200,000	U	--		--	
Formic acid		64-18-6	**55,000**		500	U	5,000	U	--		--		--		--	
Pentanoic acid		109-52-4	2,000	U	500	U	5,000	U	--		--		--		--	
Propionic Acid		79-09-4	**84,000**		500	U	5,000	U	100,000	U	200,000	U	--		--	
Hexanoic acid		142-62-1	**1,100**	**J**	500	U	5,000	U	--		300,000	U	--		--	

Table 4C. Historical Groundwater Natural Attenuation Parameter Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

Analyte	Units	CAS No.	MW4S_G 2/8/2023		DUP01_020823 2/8/2023 MW4S_G		MW4S 7/6/2023		MW4S 12/1/2023		MW4S 12/10/2024		MW4S 3/12/2025		MW4S 6/17/2025		MW4S 9/24/2025	
Dissolved Gases	ug/L																	
Ethane		74-84-0	**520**		--		**578**		**638**		**330**		**42**		**60**		**390**	
Ethene		74-85-1	**1,720**		--		**60.4**		**65.3**		**22**		17	U	**2.1**		**13**	
Methane		74-82-8	**3,910**		--		**5,610**		**6,100**		**3,900**		**390**		**790**		**4,600**	
General Chemistry																		
DO	mg/L	(field parameter)	--		--		0.24		0.88		0.00		3.28		0.00		0.15	
Oxidation Reduction Potential	millivolts	(field parameter)	--		--		115		9.8		41		24		85		80	
pH	s.u.	(field parameter)	--		--		6.12		6.83		6.91		7.29		7.24		6.82	
Sulfate	ug/L	14808-79-8	**313,000**		--		**190,000**		**150,000**		**148,000**		--		**143,000**		**131,000**	
Metals	ug/L																	
Iron, Dissolved		7439-89-6	**2,440**	**J**	**25**	**J**	**604**	**J**	**382**		**270**		**543**		**1,270**		**81**	
Iron, Total		7439-89-6	**53,200**	**J**	**151,000**	**J**	**57,100**		**9,000**		**1,900**		**2,360**		**1,340**		**1,500**	
Manganese		7439-96-5	--		--		--		**1,203**		--		--		--		--	
Volatile Fatty Acids	ug/L																	
Butyric Acid		107-92-6	**57,000**		--		500	U	5,000	U	100,000	U	200,000	U	--		--	
Lactic Acid		50-21-5	10,000	U	--		**710**		5,000	U	--		--		--		--	
Pyruvic Acid		127-17-3	**3,300**	**J**	--		500	U	5,000	U	--		--		--		--	
Acetic Acid		64-19-7	**140,000**		--		500	U	5,000	U	100,000	U	200,000	U	--		--	
Formic acid		64-18-6	**53,000**		--		500	U	5,000	U	--		--		--		--	
Pentanoic acid		109-52-4	**14,000**	**J**	--		500	U	5,000	U	--		--		--		--	
Propionic Acid		79-09-4	**380,000**		--		500	U	5,000	U	100,000	U	200,000	U	--		--	
Hexanoic acid		142-62-1	**3,800**	**J**	--		500	U	5,000	U	--		300,000	U	--		--	

Table 4C. Historical Groundwater Natural Attenuation Parameter Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

Analyte	Units	CAS No.	MW4I_G 2/8/2023		MW4I 7/6/2023		MW4I 12/1/2023		DUP-12.1.23 12/1/2023 MW4I		MW4I 12/10/2024		MW4I 3/12/2025		MW4I 6/17/2025		MW4I 9/24/2025	
Dissolved Gases	ug/L																	
Ethane		74-84-0	**16.4**		**44.8**		**32.1**		--		**24**		**16**		**17**		**15**	
Ethene		74-85-1	**59.4**		**104**		**82.5**		--		**70**		**62**		**62**		**54**	
Methane		74-82-8	**162**		**1,770**		**2,050**		--		**2,000**		**1,300**		**1,200**		**1,400**	
General Chemistry																		
DO	mg/L	(field parameter)	**--**		0.49		0.28		--		0.00		0.33		0.00		0.00	
Oxidation Reduction Potential	millivolts	(field parameter)	--		-90		-110.9		--		-62		-48		107		-72	
pH	s.u.	(field parameter)	--		6.78		7.01		--		7.15		6.86		7.38		8.08	
Sulfate	ug/L	14808-79-8	**99,600**		**140,000**		**150,000**		--		**158,000**		--		**161,000**		**128,000**	
Metals	ug/L																	
Iron, Dissolved		7439-89-6	**3,690**		**1,230**	**J**	**1,520**		--		**1,670**		**949**		**23.9**	**J**	**1,300**	
Iron, Total		7439-89-6	**547,000**		**1,500**		**1,590**		**1,600**	**J**	**1,560**		**1,020**		**1,570**		**1,200**	
Manganese		7439-96-5	--		--		**182.6**		**176.4**		--		--		--		--	
Volatile Fatty Acids	ug/L																	
Butyric Acid		107-92-6	**11,000**		500	U	5,000	U	--		100,000	U	200,000	U	--		--	
Lactic Acid		50-21-5	**2,800**	**J**	**710**		5,000	U	--		--		--		--		--	
Pyruvic Acid		127-17-3	**93,000**	**J**	500	U	5,000	U	--		--		--		--		--	
Acetic Acid		64-19-7	**36,000**		**940**		5,000	U	--		100,000	U	200,000	U	--		--	
Formic acid		64-18-6	**57,000**		500	U	5,000	U	--		--		--		--		--	
Pentanoic acid		109-52-4	**3,100**	**J**	500	U	5,000	U	--		--		--		--		--	
Propionic Acid		79-09-4	**93,000**		**1,300**		5,000	U	--		100,000	U	200,000	U	--		--	
Hexanoic acid		142-62-1	**1,200**	**J**	500	U	5,000	U	--		--		300,000	U	--		--	

Table 4. Historical Groundwater Analysis Results
UniFirst Corporation
Liverpool, NY
BCP Site No. 734152

Notes:
Data for these sampling events have not been validated. Qualifiers are Lab Qualifiers.
Analytes in blue are not detected in any sample
ng/L = nanogram per liter (ppt)
s.u. = standard units
ug/L = micrograms per liter or parts per billion (ppb)

PFAS = Per- and polyfluoroalkyl substances
VOC = Volatile Organic Compound

Total VOCs are calculated using detects only

NYS AWQS = New York State Ambient Water Quality Standards and Guidance Values for GA groundwater
* indicates the value is a guidance value and not a standard

CAS No. = Chemical Abstracts Service Number
ND = Not Detected
NE = Not Established

Bolding indicates a detected result concentration
Shading and bolding indicates that the detected concentration is above the NYSDOH guidance to which it was compared

Laboratory Qualifiers:
F = MS and/or MSD recovery exceeds the control limits
J = The result is an estimated value
U = The result was not detected above the reporting limit

Figures

SOURCE:
USGS QUADRANGLE MAP OF SYRACUSE WEST, NY 2023.

Quarterly Groundwater Monitoring Report - Q3 2025 113 7th North Street Liverpool, New York	GEI Consultants	SITE LOCATION MAP
Unifirst Liverpool Liverpool, New York	Project 2001642	February 2026 Fig. 1

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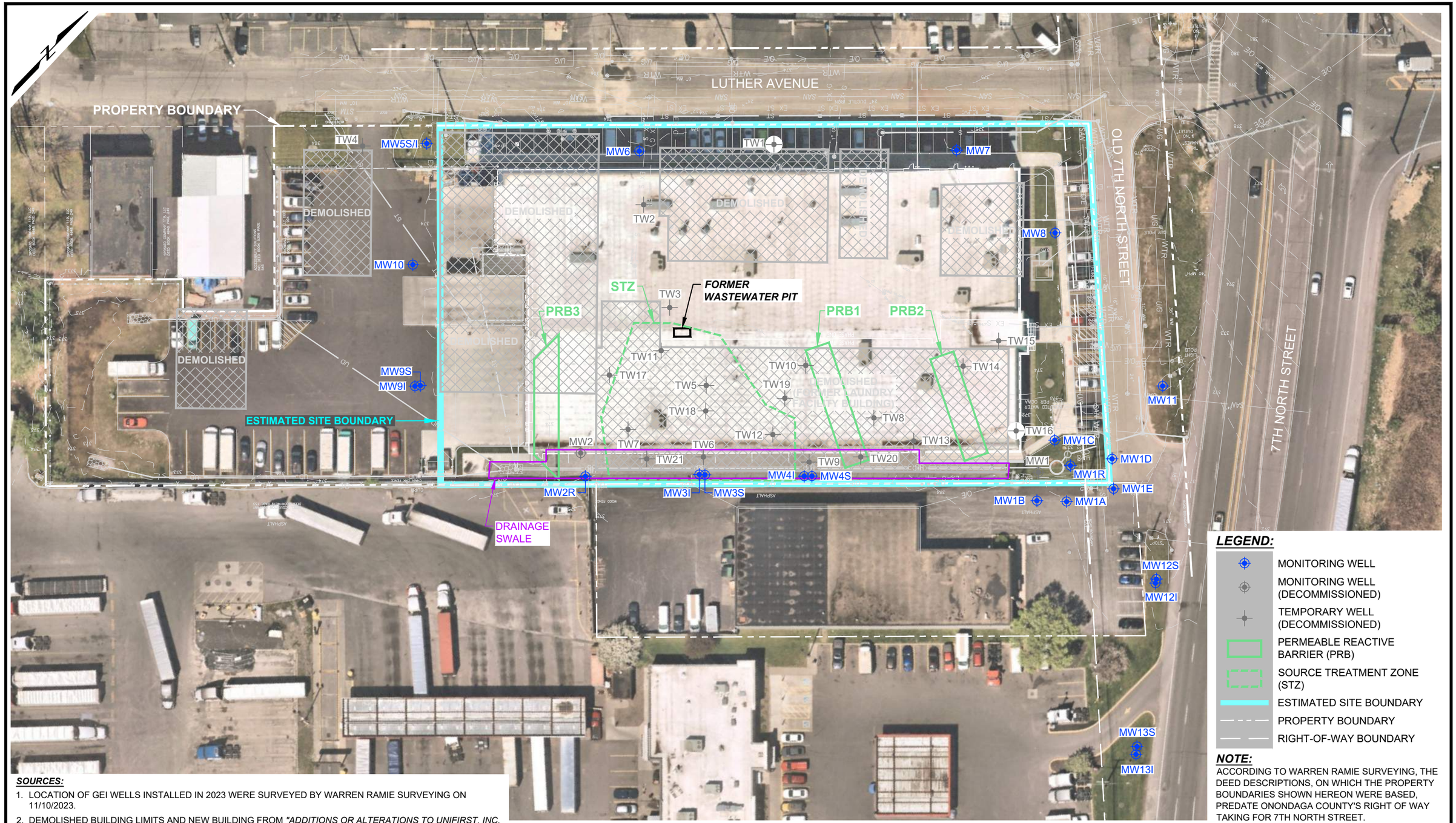

LEGEND:

- MONITORING WELL
- MONITORING WELL (DECOMMISSIONED)
- TEMPORARY WELL (DECOMMISSIONED)
- PERMEABLE REACTIVE BARRIER (PRB)
- SOURCE TREATMENT ZONE (STZ)
- ESTIMATED SITE BOUNDARY
- PROPERTY BOUNDARY
- RIGHT-OF-WAY BOUNDARY

NOTE:

ACCORDING TO WARREN RAMIE SURVEYING, THE DEED DESCRIPTIONS, ON WHICH THE PROPERTY BOUNDARIES SHOWN HEREON WERE BASED, PREDATE ONONDAGA COUNTY'S RIGHT OF WAY TAKING FOR 7TH NORTH STREET.

SOURCES:

1. LOCATION OF GEI WELLS INSTALLED IN 2023 WERE SURVEYED BY WARREN RAMIE SURVEYING ON 11/10/2023.
2. DEMOLISHED BUILDING LIMITS AND NEW BUILDING FROM *"ADDITIONS OR ALTERATIONS TO UNIFIRST, INC. NEW FACILITY, OLD 7th NORTH STREET LIVERPOOL, NEW YORK, SITE PHASING PLAN,"* SHEETS L-2 REV. 12 DATED 07.05.22. AND L2.1 REV 4 DATED 11.05.19 . PREPARED BY KEPLINGER FREEMAN ASSOCIATES.
3. SURVEY OF PROPERTIES SOUTH OF SITE BY WARREN RAMIE, PROVIDED IN A FILE NAMED "WRS 7th North Street Boundary 08.15.23.dwg".
4. WELLS MW-1A, MW-1B, MW-1C, MW-1D & MW-1E FROM FIGURE 2: AREA WIDE DIAGRAM, PREPARED BY SOVEREIGN CONSULTING, INC., DATED 4/24/2024.
5. NEARMAP AERIAL ©2024 NEARMAP, IMAGERY DATED 05/07/2024, ACCESSED ON 08/15/2024.

0 60 120

SCALE: 1" = 60'

Quarterly Groundwater Monitoring Report - Q3 2025
113 7th North Street
Liverpool, New York

Unifirst Liverpool
Liverpool, New York

GEI Consultants

SITE PLAN

Project 2001642 | February 2026 | Fig. 2

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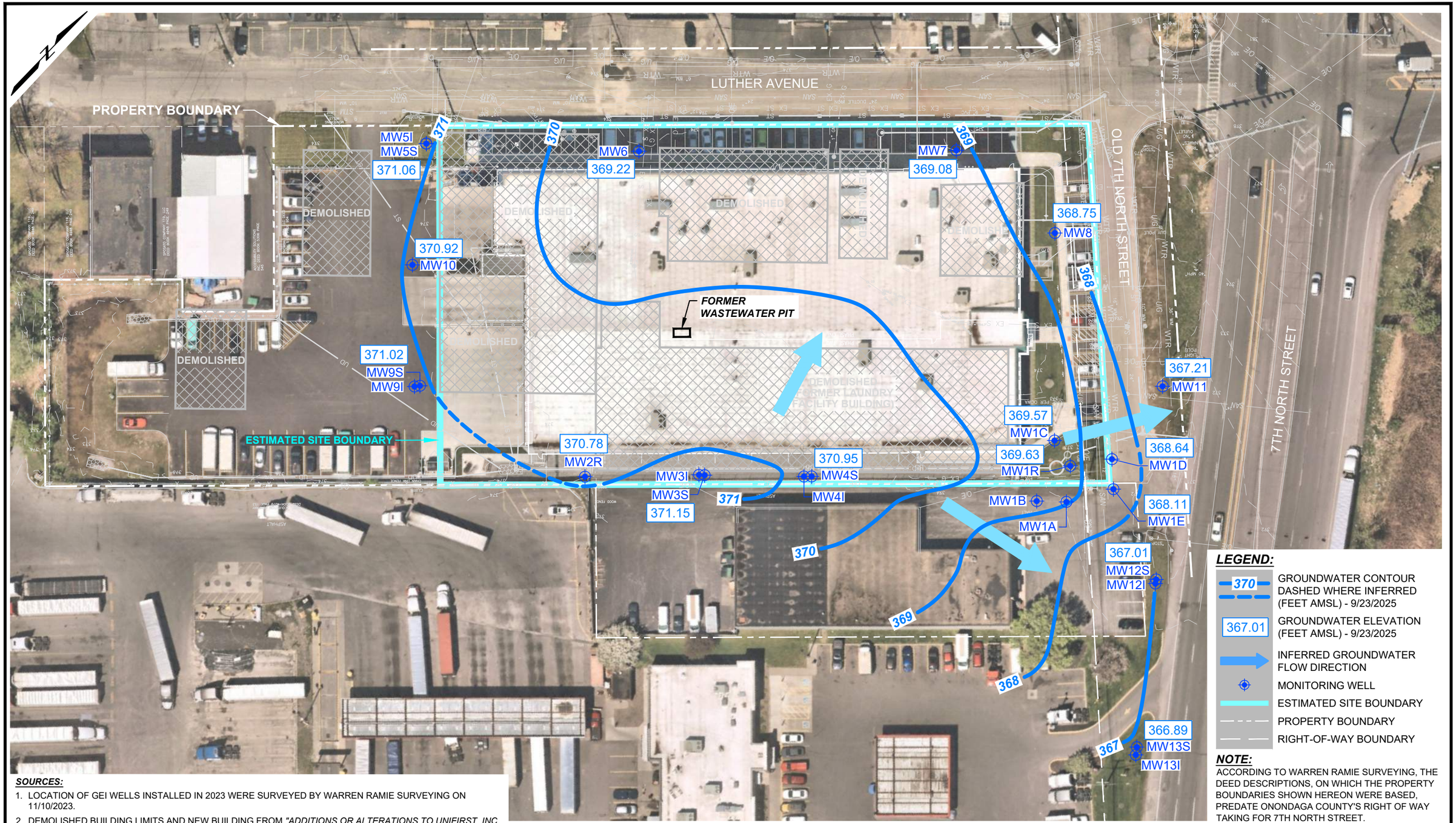

LEGEND:

- 370 — GROUNDWATER CONTOUR DASHED WHERE INFERRED (FEET AMSL) - 9/23/2025
- 367.01 — GROUNDWATER ELEVATION (FEET AMSL) - 9/23/2025
- INFERRED GROUNDWATER FLOW DIRECTION
- MONITORING WELL
- ESTIMATED SITE BOUNDARY
- PROPERTY BOUNDARY
- RIGHT-OF-WAY BOUNDARY

NOTE:

ACCORDING TO WARREN RAMIE SURVEYING, THE DEED DESCRIPTIONS, ON WHICH THE PROPERTY BOUNDARIES SHOWN HEREON WERE BASED, PREDATE ONONDAGA COUNTY'S RIGHT OF WAY TAKING FOR 7TH NORTH STREET.

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0 60 120

SCALE: 1" = 60'

Quarterly Groundwater Monitoring Report - Q3 2025
113 7th North Street
Liverpool, New York

Unifirst Liverpool
Liverpool, New York

GEI Consultants

GROUNDWATER ELEVATION SUMMARY
9/23/2025

Project 2001642 | February 2026 | Fig. 3

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MW5S*	2/21/2023		7/5/2023		11/29/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
PCE	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U	10	U
TCE	2	U	2.5	U	2.5	U	2.5	U	2.5	U	2	U	10	U
cis-1,2-DCE	10	U	12	U	12	U	12	U	12	U	10	U	10	U
trans-1,2-DCE	10	U	12	U	12	U	12	U	12	U	10	U	10	U
Vinyl chloride	**2**	**J**	**0.58**	**J**	5	U	5	U	5	U	**2.4**	**J**	10	U
Total VOCs*	**1,079**	**J**	**1,433**	**J**	**1,080**	**J**	**705**	**J**	**1,147**	**J**	**871**	**J**	**1,057**	**J**

MW6	11/28/2023		12/9/2024		3/11/2025		6/16/2025		9/23/2025	
PCE	1	U	0.5	U	0.5	U	0.5	U	1	U
TCE	0.5	U	0.5	U	0.5	U	0.5	U	1	U
cis-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U	1	U
trans-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride	1	U	1	U	1	U	**4.1**		1	U
Total VOCs	ND		ND		ND		**4.1**		ND	

TW10	3/20/2021		7/11/2021		11/21/2021		2/26/2022	
PCE	**3.7**		**46**		**8.4**		**6.6**	
TCE	**1.3**		**7.8**		**4.2**		**4.2**	
cis-1,2-DCE	**36**		**69**		**86**		**96**	
trans-1,2-DCE	2.5	U	**0.83**	**J**	**1**	**J**	**1.1**	**J**
Vinyl chloride	**4.4**		**9.1**	**J**	**7.8**		**8.6**	
Total VOCs	**45.4**		**133**	**J**	**108**	**J**	**120**	**J**

MW5I*	2/21/2023		7/5/2023		11/29/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
PCE	**0.3**	**J**	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
TCE	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U	1	U
cis-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
trans-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride	1	U	1	U	1	U	**1.1**		1	U	1	U	1	U
Total VOCs*	**5.3**	**J**	**1.03**	**J**	**0.4**	**J**	ND		ND		**3.3**	**J**	ND	

TW1	5/28/2020	
PCE	1	U
TCE	0.5	U
cis-1,2-DCE	2.5	U
trans-1,2-DCE	2.5	U
Vinyl chloride	1	UJ
Total VOCs	ND	

MW7	11/28/2023		12/9/2024		3/11/2025		6/16/2025		9/23/2025	
PCE	0.5	U	0.5	U	0.5	U	0.5	U	1	U
TCE	0.5	U	0.5	U	0.5	U	0.5	U	1	U
cis-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U	1	U
trans-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride	**4.7**		**15**		**12**		**15**		13	
Total VOCs	**4.7**		**15**		**12**		**19.9**	**J**	13	

TW15	3/20/2021		7/10/2021		11/22/2021		2/27/2022	
PCE	**1.9**		**2.5**		**1.9**		**1.5**	
TCE	**0.29**	**J**	**0.36**	**J**	**0.43**	**J**	**0.22**	**J**
cis-1,2-DCE	2.5	U	2.5	U	**1.1**	**J**	**2.6**	
trans-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U
Vinyl chloride	1	U	1	UJ	1	UJ	1	U
Total VOCs	**3.99**	**J**	**7.83**	**J**	**9.73**	**J**	**13.0**	**J**

TW11	3/20/2021		7/11/2021		11/21/2021		2/26/2022	
PCE	**23,000**	**J**	**51,000**		**61,000**		**43,000**	
TCE	**2,400**		**2,600**		**3,800**		**2,200**	
cis-1,2-DCE	**5,700**		**5,100**		**7,700**		**7,800**	
trans-1,2-DCE	**42**	**J**	1,200	U	1,200	U	1,200	U
Vinyl chloride	**1,400**		**1,400**		**1,400**		**1,100**	
Total VOCs	**32,566**	**J**	**60,100**		**73,900**		**54,100**	

TW4	5/28/2020	
PCE	0.5	U
TCE	0.5	U
cis-1,2-DCE	2.5	U
trans-1,2-DCE	2.5	U
Vinyl chloride	1	UJ
Total VOCs	**16.3**	**J**

TW2	5/28/2020	
PCE	**0.62**	
TCE	0.5	U
cis-1,2-DCE	**11**	
trans-1,2-DCE	2.5	U
Vinyl chloride	**2.2**	**J**
Total VOCs	**16.4**	**J**

TW3	5/28/2020	
PCE	**5,300**	
TCE	**810**	
cis-1,2-DCE	**3,400**	
trans-1,2-DCE	120	U
Vinyl chloride	**350**	**J**
Total VOCs	**9,871**	**J**

TW19	11/21/2021		2/26/2022	
PCE	**9.5 [9.8]**		**7.2**	
TCE	**1.2 [1.2]**		**0.82**	
cis-1,2-DCE	**11 [11 J]**		**17**	
trans-1,2-DCE	2.5 [2.5]	U	2.5	U
Vinyl chloride	1 [1]	UJ	1	U
Total VOCs	**23.2 [23.5] J**		**26.4**	**J**

MW8*	11/29/2023		12/9/2024		3/11/2025		6/16/2025		9/24/2025	
PCE	0.5	U	0.5	U	0.5	U	0.5	U	1	U
TCE	0.5	U	0.5	U	0.5	U	0.5	U	1	U
cis-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U	1	U
trans-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride	1	U	**0.21**	**J**	**0.11**	**J**	**0.43**	**J**	1	U
Total VOCs	ND		**2.75**	**J**	**4.21**	**J**	**8.84**	**J**	**3.5**	**J**

TW14	3/21/2021		7/10/2021		11/20/2021		2/26/2022	
PCE	**7.8**		**0.34**	**J**	**3.6**		**7.3**	
TCE	**6.6**		0.5	U	**0.98**		**1.3**	
cis-1,2-DCE	**6.2**		**2.2**	**J**	**19**		**16**	
trans-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U
Vinyl chloride	**0.15**	**J**	1	UJ	1	UJ	1	U
Total VOCs	**27.4**	**J**	**15.7**	**J**	**39.6**	**J**	**51.8**	

MW10	11/29/2023	
PCE	0.5	U
TCE	0.5	U
cis-1,2-DCE	2.5	U
trans-1,2-DCE	2.5	U
Vinyl chloride	1	U
Total VOCs	ND	

TW16	3/21/2021		7/10/2021		11/22/2021		2/28/2022	
PCE	**41**		**18**		**13**		**13**	
TCE	**3.7**		**2**		**1.6**		**1.1**	
cis-1,2-DCE	**22**		**16**		**14**		**15**	
trans-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U
Vinyl chloride	**0.53**	**J**	**0.33**	**J**	**0.2**	**J**	**0.26**	**J**
Total VOCs	**67.2**	**J**	**36.3**	**J**	**28.8**	**J**	**29.4**	**J**

TW17	11/21/2021		2/26/2022	
PCE	**0.43**	**J**	**0.26**	**J**
TCE	**0.71**		**0.26**	**J**
cis-1,2-DCE	**1.7**	**J**	**1.2**	**J**
trans-1,2-DCE	2.5	U	2.5	U
Vinyl chloride	**0.91**	**J**	**0.68**	**J**
Total VOCs	**3.75**	**J**	**2.4**	**J**

MW11	11/29/2023		12/11/2024		3/11/2025		6/16/2025		9/24/2025	
PCE	0.5	U	0.5	U	0.5	U	0.5	U	1	U
TCE	**0.27**	**J**	0.5	U	0.5	U	0.5	U	1	U
cis-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U	1	U
trans-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride	**0.34**	**J**	**0.36**	**J**	**0.14**	**J**	**2.6**		1	U
Total VOCs	**0.61**	**J**	**0.64**	**J**	**0.14**	**J**	**4.5**		ND	

MW9S	11/29/2023	
PCE	**0.41**	**J**
TCE	0.5	U
cis-1,2-DCE	2.5	U
trans-1,2-DCE	2.5	U
Vinyl chloride	1	U
Total VOCs	**0.41**	**J**

MW-1C	6/18/2025	9/25/2025
PCE	**1.1**	1 U
TCE	0.5 U	1 U
cis-1,2-DCE	2.5 U	1 U
trans-1,2-DCE	2.5 U	1 U
Vinyl chloride	**0.13** J	1 U
Total VOCs	**7.13**	**5.6** J

MW9I	11/29/2023	
PCE	0.5	U
TCE	0.5	U
cis-1,2-DCE	2.5	U
trans-1,2-DCE	2.5	U
Vinyl chloride	1	U
Total VOCs	ND	

MW-1D	6/18/2025	9/25/2025
PCE	**1,000**	**540**
TCE	**210**	**210**
cis-1,2-DCE	**230**	**990**
trans-1,2-DCE	25 U	20 U
Vinyl chloride	**16**	**28**
Total VOCs	**1,468**	**1,768**

MW-1E	6/18/2025	9/25/2025
PCE	**3,600**	**3,600**
TCE	**760**	**1,100**
cis-1,2-DCE	**430**	**2,900**
trans-1,2-DCE	62 U	100 U
Vinyl chloride	**12** J	100 U
Total VOCs	**4,802**	**7,600**

TW5	3/20/2021		7/11/2021		11/21/2021		2/26/2022	
PCE	**6,800**		**490**		**450**		**420**	
TCE	**510**		**58**		**50**		**48**	
cis-1,2-DCE	**1,400**		**330**		**270**		**250**	
trans-1,2-DCE	120	U	10	U	**1.4**	**J**	10	U
Vinyl chloride	**77**		**3.1**	**J**	**2.1**		**1.5**	**J**
Total VOCs	**8,787**		**881**	**J**	**774**	**J**	**720**	**J**

MW12S	11/30/2023		12/11/2024		3/11/2025		6/18/2025		9/24/2025	
PCE	**0.5**		0.5	U	0.5	U	0.5	U	1	U
TCE	**0.61**		**0.32**	**J**	**0.24**	**J**	**0.45**	**J**	1	U
cis-1,2-DCE	**7.2**		**3.4**		**2.3**	**J**	**1.7**	**J**	**2.3**	
trans-1,2-DCE	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride	**0.47**	**J**	**3.7**		**1.4**		**1.3**		**1.5**	
Total VOCs	**8.78**	**J**	**7.42**	**J**	**3.94**	**J**	**6.15**		**3.8**	

TW18	11/21/2021		2/26/2022	
PCE	**690**		**590**	
TCE	**78**		**69**	
cis-1,2-DCE	**350**		**320**	
trans-1,2-DCE	12	U	10	U
Vinyl chloride	**1.9**	**J**	**0.75**	**J**
Total VOCs	**1,120**	**J**	**980**	**J**

MW12I	11/30/2023	
PCE	0.5	U
TCE	0.5	U
cis-1,2-DCE	2.5	U
trans-1,2-DCE	2.5	U
Vinyl chloride	1	U
Total VOCs	ND	

MW13S	11/30/2023	
PCE	0.5	U
TCE	0.5	U
cis-1,2-DCE	2.5	U
trans-1,2-DCE	2.5	U
Vinyl chloride	**0.3**	**J**
Total VOCs	**1.6**	**J**

TW7	3/20/2021		7/11/2021		11/21/2021		2/26/2022	
PCE	**4,100**		**13**		**0.98**		**1.5**	
TCE	**560**		**14**		**1.2**		**1.1**	
cis-1,2-DCE	**760**		**75**		**12**		**6.8**	
trans-1,2-DCE	62	U	4.5		**0.98**	**J**	2.5	U
Vinyl chloride	**22**	**J**	**12**		**2.9**		**1.7**	
Total VOCs	**5,442**	**J**	**119**	**J**	**18.1**	**J**	**11.1**	

MW13I	11/30/2023	
PCE	0.5	U
TCE	0.5	U
cis-1,2-DCE	2.5	U
trans-1,2-DCE	2.5	U
Vinyl chloride	1	U
Total VOCs	ND	

MW2/MW2G/MW2R	5/28/2020		3/20/2021		7/11/2021		11/22/2021		2/28/2022		2/7/2023		7/6/2023		12/1/2023		12/11/2024		3/12/2025		6/17/2025		9/24/2025	
PCE	**4**		**64 [61]**		**0.76**		**2.1**		**0.43**	**J**	**470**		**150**	**J**	**42**		**1.7**		**16**		**0.27**	**J**	6.7	
TCE	**0.91**		**7.1 [6.7]**		**0.34**	**J**	0.35	**J**	0.5	U	**120**		**18**		**17**		**2.8**		**7.2**		**3.7**		4.4	
cis-1,2-DCE	**3.4**		**8.5 [7.8]**		**0.92**	**J**	2.5	U	2.5	U	**130**		**38**		**51**		**28**		**21**		**37**		39	
trans-1,2-DCE	2.5	U	2.5 [2.5]	U	2.5	U	2.5	U	2.5	U	12	U	2.5	U	2.5	U	2.5	U	2.5	U	**0.85**	**J**	1.7	
Vinyl chloride	**0.44**	**J**	**0.66 [0.64]**	**J**	1	U	**0.09**	**J**	1	U	**1**	**J**	**0.79**	**J**	**1.2**		**1.4**		**0.4**	**J**	**1.7**		3.9	
Total VOCs	**31.9**	**J**	**134 [130]**	**J**	**31.2**	**J**	**11.8**	**J**	**12.1**	**J**	**1,267**	**J**	**225**	**J**	**111**	**J**	**36**	**J**	**44.6**	**J**	**49.9**	**J**	55.7	

TW12	3/21/2021		7/11/2021		11/21/2021		2/26/2022	
PCE	**25,000**	**J**	**1,200**		**480**		**630**	
TCE	**10,000**		**440**		**110**		**170**	
cis-1,2-DCE	**22,000**		**1,600**		**710**		**1,100**	
trans-1,2-DCE	120		14	**J**	6.2	**J**	25	U
Vinyl chloride	**3,200**		**120**	**J**	**21**		**39**	
Total VOCs	**60,385**	**J**	**3,376**	**J**	**1,327**	**J**	**1,939**	

MW1/MW1G/MW1R*	5/28/2020		3/21/2021		7/10/2021		11/22/2021		2/28/2022		2/7/2023		11/30/2023		12/11/2024		3/12/2025		9/24/2025	
PCE	**1,400**		**29**		**8.7**		**22**		**25**		**52**		**91**		0.5	U	0.5	U	**1.5**	**J**	10	U
TCE	**100**		**3.8**		**1.9**		**3.7**		**4**	**J**	**8.7**		**27**		**0.29**	**J**	0.5	U	2.5	U	10	U
cis-1,2-DCE	**80**		**17**		**20**		**29**		**24**	**J**	**37**		**130**		**14**		**0.78**	**J**	12	U	10	U
trans-1,2-DCE	25	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	**0.71**	**J**	2.5	U	2.5	U	12	U	10	U
Vinyl chloride	**3.7**	**J**	**0.73**	**J**	**1.4**	**J**	**2**	**J**	2	U	2	U	**54**		**14**		**0.57**	**J**	**0.61**	**J**	10	U
Total VOCs	**1,620**	**J**	**50.5**	**J**	**32**	**J**	**56.7**	**J**	**55**	**J**	**99.9**	**J**	**303**	**J**	**48.6**	**J**	**72**	**J**	**22**	**J**	5.1	**J**

TW21	11/20/2021		2/26/2022	
PCE	**3,400**		**2,000**	
TCE	**370**		**200**	
cis-1,2-DCE	**310**		**330**	
trans-1,2-DCE	62	U	50	U
Vinyl chloride	**69**	**J**	**95**	
Total VOCs	**4,149**	**J**	**2,625**	

MW4IG/MW4I	2/8/2023		7/6/2023		12/1/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
PCE	**5,800**		**7,700**	**J**	**720 [670]**		**36 [34]**		**31**		**16**		20	**J**
TCE	**720**		**2,000**		**2,200 [1,900]**		**70 [78]**		**28**		**12**		40	U
cis-1,2-DCE	**8,500**		**3,600**		**6,700 [6,000]**		**3,200 [3,200]**		**2,900**		**2,000**		3,800	
trans-1,2-DCE	250	U	120	U	120 [120]	U	**25 [23]**	**J**	50	U	62	U	40	U
Vinyl chloride	**540**		**360**		**330 [330]**		**390 [380]**		**300**		**280**		550	
Total VOCs	**15,560**		**13,685**	**J**	**9,979 [8,926]**		**3,744 [3,741]**	**J**	**3,283**		**2,323**		4,406	**J**

TW13	3/21/2021		7/10/2021		11/21/2021		2/27/2022	
PCE	**40**		**72 [59]**		**70**		**57**	
TCE	**7**		**13 [11]**		**14**		**10**	
cis-1,2-DCE	**110**		**180 [180]**		**220**		**140**	
trans-1,2-DCE	2.5	U	0.82 J [2.5 U]		5	U	2.5	U
Vinyl chloride	**4.6**		**2.2 [2.3 J]**		**1.3**	**J**	**1.1**	
Total VOCs	**162**		**268 [252]**	**J**	**305**	**J**	**210**	**J**

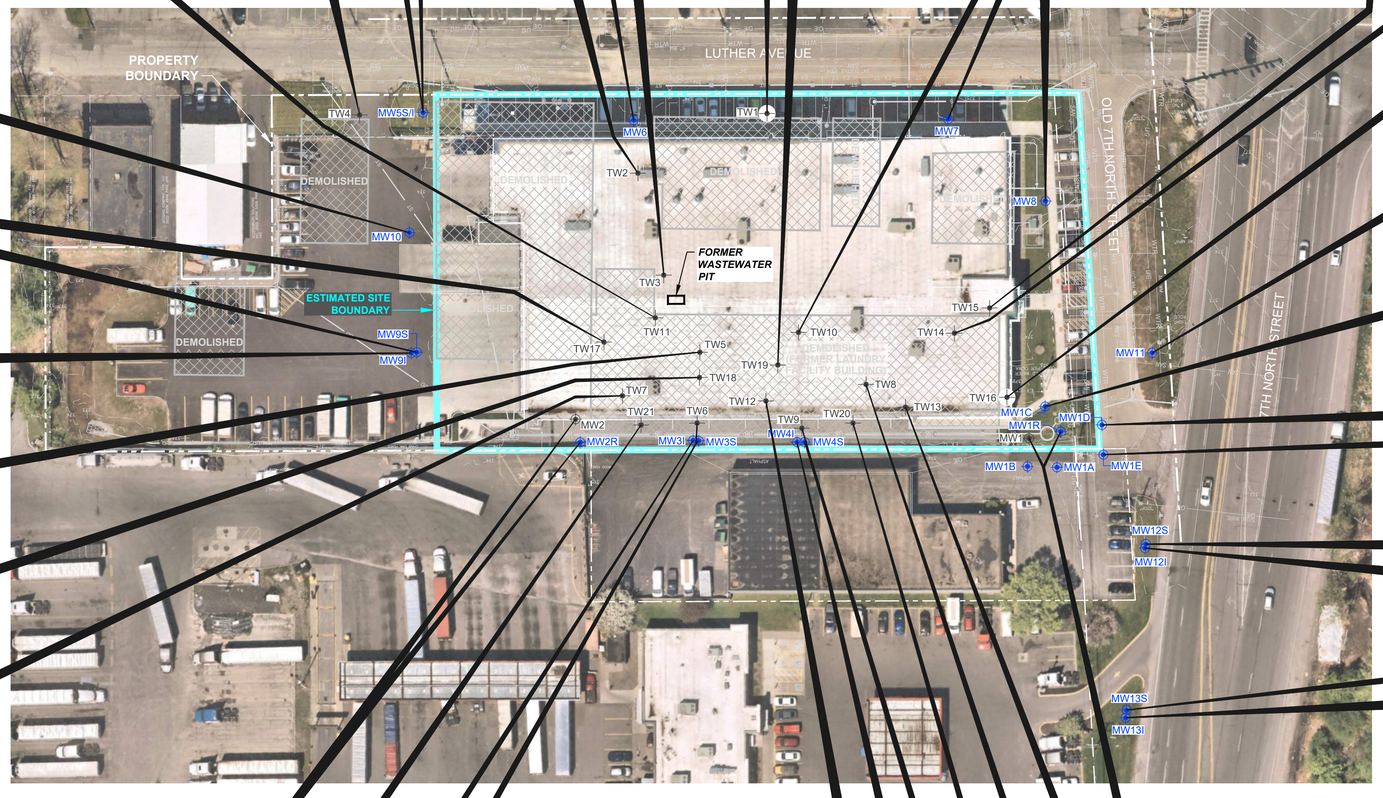
MW3IG/MW3I	2/8/2023		7/6/2023		11/30/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
PCE	**13,000**		**19**		**2.3**		**0.64**		**1.1**		**0.41**	**J**	1	U
TCE	**690**		**18**		**1.4**		**0.6**		**1.6**		**0.75**		0.53	**J**
cis-1,2-DCE	**4,000**		**970**		**140**		**42**		**56**		**5**		2.9	
trans-1,2-DCE	250	U	25	U	2.5	U	2.5	U	2.5	U	2.5	U	1	U
Vinyl chloride	**120**		**8.6**	**J**	**7.3**		**39**		**24**		**6.1**		8.4	
Total VOCs	**17,960**	**J**	**1,042**	**J**	**153**		**83**		**83.1**	**J**	**12.3**	**J**	11.8	**J**

TW6/MW3SG/MW3S**	3/20/2021		7/11/2021		11/21/2021		2/27/2022		2/7/2023		7/6/2023		12/1/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
PCE	**40,000**	**J**	**21,000**		**20,000**		**34,000 [24,000]**		**63,000**		**41,000**		**1,400**		500	U	500	U	500	U	2,000	U
TCE	**6,500**		**6,100**		**4,300**		**8,100 [4,900]**		**5,000**		**30,000**		**650**		500	U	500	U	500	U	2,000	U
cis-1,2-DCE	**16,000**		**11,000**		**8,400**		**14,000 [10,000]**		**100**	**J**	**90000**		**190,000**		**92,000**		**100,000**		**86,000**		160,000	
trans-1,2-DCE	**110**		**170**	**J**	500	U	620 [620]	U	**57,000**		1,200	U	2,500	U	2,500	U	2,500	U	2,500	U	2,000	U
Vinyl chloride	**2,200**		**2,400**	**J**	**970**		**2,100 [1,500]**		**2,800**		**12,000**		**38,000**		**99,000**		**66,000**		**71,000**		120,000	
Total VOCs	**64,850**	**J**	**40,670**	**J**	**33,670**		**58,200 [40,400]**		**127,900**	**J**	**173,200**	**J**	**230,270**	**J**	**191,000**		**166,000**		**157,000**		280,000	

TW8	3/21/2021		7/10/2021		11/20/2021		2/26/2022	
PCE	**97**		**88**		**100**		**110**	
TCE	**12**		**13**		**13**		**15**	
cis-1,2-DCE	**160**		**160**		**170**		**180**	
trans-1,2-DCE	**0.94**	**J**	**0.95**	**J**	**0.91**	**J**	**1.1**	**J**
Vinyl chloride	**2**		**0.2**	**J**	**0.14**	**J**	**0.23**	**J**
Total VOCs	**273**	**J**	**264**	**J**	**286**	**J**	**308**	**J**

TW9/MW4SG/MW4S**	3/21/2021		7/11/2021		11/20/2021		2/26/2022		2/8/2023		7/6/2023		12/1/2023		12/10/2024		3/12/2025		6/17/2025		9/24/2025	
PCE	**390**		**53**		**140**		**61**		**43,000 [41,000]**		**120,000**		**140,000**	**J**	**82,000**		**19,000**		**19,000**		85,000	
TCE	**400**		**27**		**40**		**47**		**7,200 [6,800]**		**7,400**		**5,900**		**4,500**		**590**		**690**		4,100	
cis-1,2-DCE	**1,800**		**290**		**470**		**460**		**76,000 [70,000]**	**J**	**16,000**		**17,000**		**9,600**		**1,000**		**1,400**		8,100	
trans-1,2-DCE	50	U	5	U	10	U	10	U	1,200 [1,200]	U	2,500	U	500	U	2,500	U	250	U	250	U	2,000	U
Vinyl chloride	**120**		**10**	**J**	**7.8**		**24**		**8,500 [8,100]**		**1,400**		**870**		**860**	**J**	**84**	**J**	**96**	**J**	2,000	U
Total VOCs	**2,713**	**J**	**380**	**J**	**658**		**592**		**134,800 [125,994]**	**J**	**144,800**		**163,805**	**J**	**96,960**	**J**	**20,674**	**J**	**21,286 J [22,230 J]**		97,200	

TW20	11/20/2021		2/26/2022	
PCE	**88**		**38**	
TCE	**20**		**7.8**	
cis-1,2-DCE	**270**		**130**	
trans-1,2-DCE	5	U	2.5	U
Vinyl chloride	2	UJ	1	U
Total VOCs	**378**		**176**	



LEGEND:

- MONITORING WELL
- MONITORING WELL (DECOMMISSIONED)
- TEMPORARY WELL (DECOMMISSIONED)
- ESTIMATED SITE BOUNDARY
- PROPERTY BOUNDARY
- RIGHT-OF-WAY BOUNDARY

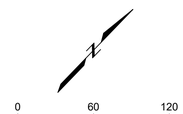
ANALYTICAL BOXES

µg/L	MICROGRAMS PER LITER (ALL RESULTS ARE PRESENTED IN µg/L)
NE	NO EXCEEDANCES
J	RESULT IS AN ESTIMATED VALUE
U	RESULT WAS NOT DETECTED ABOVE THE REPORTING LIMIT
UJ	ANALYTE WAS NOT DETECTED ABOVE THE REPORTED SAMPLE QUANTITATION LIMIT. DATA MAY BE BIASED LOW.
7.8	COMPOUND EXCEEDS NYS AMBIENT WATER QUALITY STANDARDS
BLUE	CURRENT DATA

CVOC NYS AWQS	
Analyte	Standard (µg/L)
Tetrachloroethene (PCE)	5
Trichloroethene (TCE)	5
cis-1,2-Dichloroethene	5
trans-1,2-Dichloroethene	5
Vinyl chloride	2

NOTES:

1. CVOC RESULTS ARE SUMMARIZED ON THIS FIGURE BECAUSE THEY ARE THE PRIMARY SITE CONSTITUENTS OF CONCERN. IN CERTAIN WELLS, OTHER VOCS ALSO EXCEED AWQS. REFER TO THE GROUNDWATER DATA TABLES.
2. * NON-CVOC COMPOUNDS CONTRIBUTE TO THE TOTAL VOC CONCENTRATION AT THESE WELLS.
3. ** PERMANENT WELLS (MW3S AND MW4S) WERE INSTALLED AT THE APPROX. LOCATION OF FORMER TEMPORARY WELLS (TW6 AND TW9), AS SUCH DATA FOR THESE CO-LOCATED WELLS ARE PRESENTED TOGETHER. DATA ASSOCIATED WITH TW-DESIGNATED WELLS WERE COLLECTED IN 2021 AND 2022, WHILE DATA ASSOCIATED WITH MW-DESIGNATED WELLS WERE COLLECTED IN 2023 THROUGH PRESENT.
4. *** HYDROPUNCH SAMPLERS WERE USED TO COLLECT GROUNDWATER FROM TEMPORARY BORINGS AT THE APPROXIMATE LOCATIONS OF MW1R, MW2R, MW3S, MW3I, MW4S, AND MW4I IN FEBRUARY 2023 BECAUSE PERMANENT WELLS COULD NOT BE INSTALLED AT THAT TIME DUE TO ONGOING FACILITY CONSTRUCTION. THESE SAMPLES ARE DESIGNATED AS "MWxxG" IN FIGURE DATA BOXES AND IN REPORT DATA TABLES.
5. ACCORDING TO WARREN RAMIE SURVEYING, THE DEED DESCRIPTIONS, ON WHICH THE PROPERTY BOUNDARIES SHOWN HEREON WERE BASED, PREDATE ONONDAGA COUNTY'S RIGHT OF WAY TAKING FOR 7TH NORTH STREET.



SOURCES:

1. LOCATION OF GEI WELLS INSTALLED IN 2023 WERE SURVEYED BY WARREN RAMIE SURVEYING ON 11/10/2023.
2. DEMOLISHED BUILDING LIMITS AND NEW BUILDING FROM *"ADDITIONS OR ALTERATIONS TO UNIFIRST, INC. NEW FACILITY, OLD 7th NORTH STREET LIVERPOOL, NEW YORK, SITE PHASING PLAN,"* SHEETS L-2 REV. 12 DATED 07.05.22. AND L2.1 REV 4 DATED 11.05.19. PREPARED BY KEPLINGER FREEMAN ASSOCIATES.
3. SURVEY OF PROPERTIES SOUTH OF SITE BY WARREN RAMIE, PROVIDED IN A FILE NAMED "WRS 7th North Street Boundary 08.15.23.dwg".
4. WELLS MW-1A, MW-1B, MW-1C, MW-1D & MW-1E FROM *FIGURE 2: AREA WIDE DIAGRAM*, PREPARED BY SOVEREIGN CONSULTING, INC., DATED 4/24/2024.
5. NEARMAP AERIAL ©2024 NEARMAP, IMAGERY DATED 05/07/2024, ACCESSED ON 08/15/2024.

Quarterly Groundwater Monitoring Report - Q3 2025
113 7th North Street
Liverpool, New York

Unifirst Liverpool
Liverpool, New York

GEI Consultants

SUMMARIZED GROUNDWATER ANALYTICAL RESULTS - CVOCs

Project 2001642 | February 2026 | Fig. 4a

MW5S					
	11/29/2023	12/10/2024	3/12/2025	6/17/2025	9/24/2025
PFOS	32.2	41.4	37.9	36.3	20
PFOA	91.5	46.4	37.4	38.8	14
Total PFAS	238 J	162 J	142 J	125.3 J	41.5 J

MW5I					
	11/29/2023	12/10/2024	3/12/2025	6/17/2025	9/24/2025
PFOS	1.89 U	0.531 J	0.509 J	0.448 J	1.6 U
PFOA	0.257 J	0.343 J	1.57 U	0.247 J	1.6 U
Total PFAS	46 J	106 J	43.2 J	32.2 J	52.6 J

MW10	
	11/29/2023
PFOS	6.43
PFOA	5.56
Total PFAS	38.76 J

TW7	
	3/20/2021
PFOS	13.9
PFOA	169
Total PFAS	236.51 J

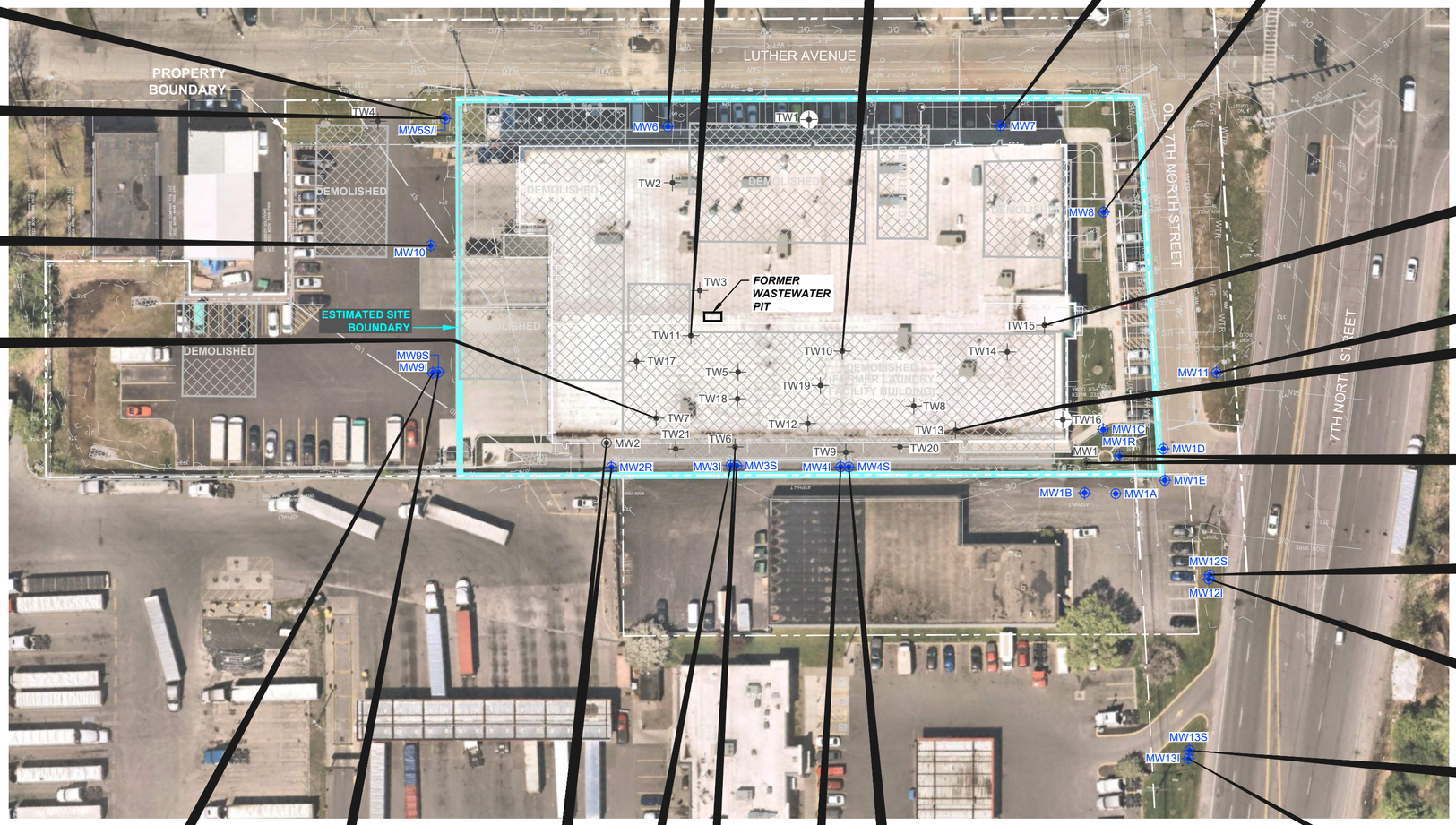
MW6	
	11/28/2023
PFOS	1.93 U
PFOA	1.93 U
Total PFAS	2.03 J

TW11	
	2/26/2022
PFOS	1,230
PFOA	237
Total PFAS	2,413 J

TW10		
	3/20/2021	2/26/2022
PFOS	89 [86.9]	56.6
PFOA	57.9 [56.6]	31.2
Total PFAS	526.2 [522.49] J	211.91 J

MW7	
	11/28/2023
PFOS	1.84 U
PFOA	1.84 U
Total PFAS	4.89 J

MW8	
	11/29/2023
PFOS	1.89 U
PFOA	8.09
Total PFAS	64.6 J



TW15		
	3/20/2021	2/27/2022
PFOS	42.2	11.3
PFOA	20.8	3.8
Total PFAS	154.52 J	217.5 J

MW11	
	11/29/2023
PFOS	9.53
PFOA	3.95
Total PFAS	27.59 J

TW13	
	2/27/2022
PFOS	106
PFOA	27.6
Total PFAS	328.06 J

MW1/MW1G/MW1R***		
	2/28/2022	11/30/2023
PFOS	61.8 J	22.4
PFOA	30.7 J	6.32
Total PFAS	216.59 J	71.21 J

MW12S	
	11/30/2023
PFOS	10.3
PFOA	5.35 J
Total PFAS	41.08 J

MW12I	
	11/30/2023
PFOS	1.9 U
PFOA	1.9 U
Total PFAS	0.913 J

MW13S	
	11/30/2023
PFOS	3.41
PFOA	5.86
Total PFAS	56.34 J

MW9I	
	11/29/2023
PFOS	1.94 U
PFOA	1.94 U
Total PFAS	3.43 J

MW9S	
	11/29/2023
PFOS	8.14
PFOA	5.54
Total PFAS	35.64 J

MW2/MW2G/MW2R***		
	2/28/2022	12/1/2023
PFOS	2.21 J	17.9
PFOA	0.886 J	13
Total PFAS	9.97 J	55.26 J

MW4S					
	12/1/2023	12/10/2024	3/12/2025	6/17/2025	9/24/2025
PFOS	12.9	45.4	31.2	75.0	17
PFOA	37.5	30.9	26.8	21.2	24
Total PFAS	152	159 J	235 J	161.6 J	105 J

MW13I	
	11/30/2023
PFOS	1.89 U
PFOA	1.89 U
Total PFAS	1.02 J

MW3IG/MW3I					
	11/30/2023	12/10/2024	3/12/2025	6/17/2025	9/24/2025
PFOS	1.87 U	1.63 U	3.49	0.626 J	1.6 U
PFOA	1.87 U	1.63 U	0.288 J	1.62 U	1.6 U
Total PFAS	3.1 J	2.8 J	7.21 J	0.626 J	0.97 J

MW4I					
	12/1/2023	12/10/2024	3/12/2025	6/17/2025	9/24/2025
PFOS	1.84 [1.84] U	1.59 [1.59] U	1.58 U	0.256 J	1.5 U
PFOA	1.84 [1.84] U	1.59 [1.59] U	1.58 U	1.51 U	1.5 U
Total PFAS	7.5 [9.6] J	2.4 [2.6] J	2.68 J	1.1 J	1.1 J

TW6/MW3SG/MW3S**						
	2/27/2022	12/1/2023	12/10/2024	3/12/2025	6/17/2025	9/24/2025
PFOS	118 [122]	10.1	19.1	24.4	10.4	6.6
PFOA	34.1 [35]	17.2	32.2	30.4	18.3	15
Total PFAS	296 [302]	106 J	118 J	141 J	102.8 J	54.7

LEGEND:

- MONITORING WELL
- MONITORING WELL (DECOMMISSIONED)
- TEMPORARY WELL (DECOMMISSIONED)
- ESTIMATED SITE BOUNDARY
- PROPERTY BOUNDARY
- RIGHT-OF-WAY BOUNDARY

ANALYTICAL BOXES

ng/L	NANOGRAMS PER LITER
NE	NO EXCEEDANCES
J	RESULT IS AN ESTIMATED VALUE
U	RESULT WAS NOT DETECTED ABOVE THE REPORTING LIMIT
--	NOT ANALYZED
8.14	COMPOUND EXCEEDS NYS AMBIENT WATER QUALITY STANDARDS
BLUE	CURRENT DATA

PFAS NYS AWQS	
Analyte	Standard (ng/L)
PFOS	2.7*
PFOA	6.7*

NOTES:

1. * INDICATES THE VALUE IS A GUIDANCE VALUE NOT A STANDARD.
2. ** PERMANENT WELL (MW3S) WAS INSTALLED AT THE APPROX. LOCATION OF FORMER TEMPORARY WELL (TW6), AS SUCH DATA FOR THIS CO-LOCATED WELL IS PRESENTED TOGETHER. DATA ASSOCIATED WITH TW6 ARE DATED 2021 AND 2022, WHILE DATA ASSOCIATED WITH MW3S ARE DATED 2023 THROUGH PRESENT.
3. ACCORDING TO WARREN RAMIE SURVEYING, THE DEED DESCRIPTIONS, ON WHICH THE PROPERTY BOUNDARIES SHOWN HEREON WERE BASED, PREDATE ONONDAGA COUNTY'S RIGHT OF WAY TAKING FOR 7TH NORTH STREET.

SOURCES:

1. LOCATION OF GEI WELLS INSTALLED IN 2023 WERE SURVEYED BY WARREN RAMIE SURVEYING ON 11/10/2023.
2. DEMOLISHED BUILDING LIMITS AND NEW BUILDING FROM *"ADDITIONS OR ALTERATIONS TO UNIFIRST, INC. NEW FACILITY, OLD 7th NORTH STREET LIVERPOOL, NEW YORK, SITE PHASING PLAN,"* SHEETS L-2 REV. 12 DATED 07.05.22. AND L2.1 REV 4 DATED 11.05.19 . PREPARED BY KEPLINGER FREEMAN ASSOCIATES.
3. SURVEY OF PROPERTIES SOUTH OF SITE BY WARREN RAMIE, PROVIDED IN A FILE NAMED "WRS 7th North Street Boundary 08.15.23.dwg".
4. NEARMAP AERIAL ©2024 NEARMAP, IMAGERY DATED 05/07/2024, ACCESSED ON 08/15/2024.

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SCALE: 1" = 60'

Quarterly Groundwater Monitoring Report - Q3 2025
113 7th North Street
Liverpool, New York

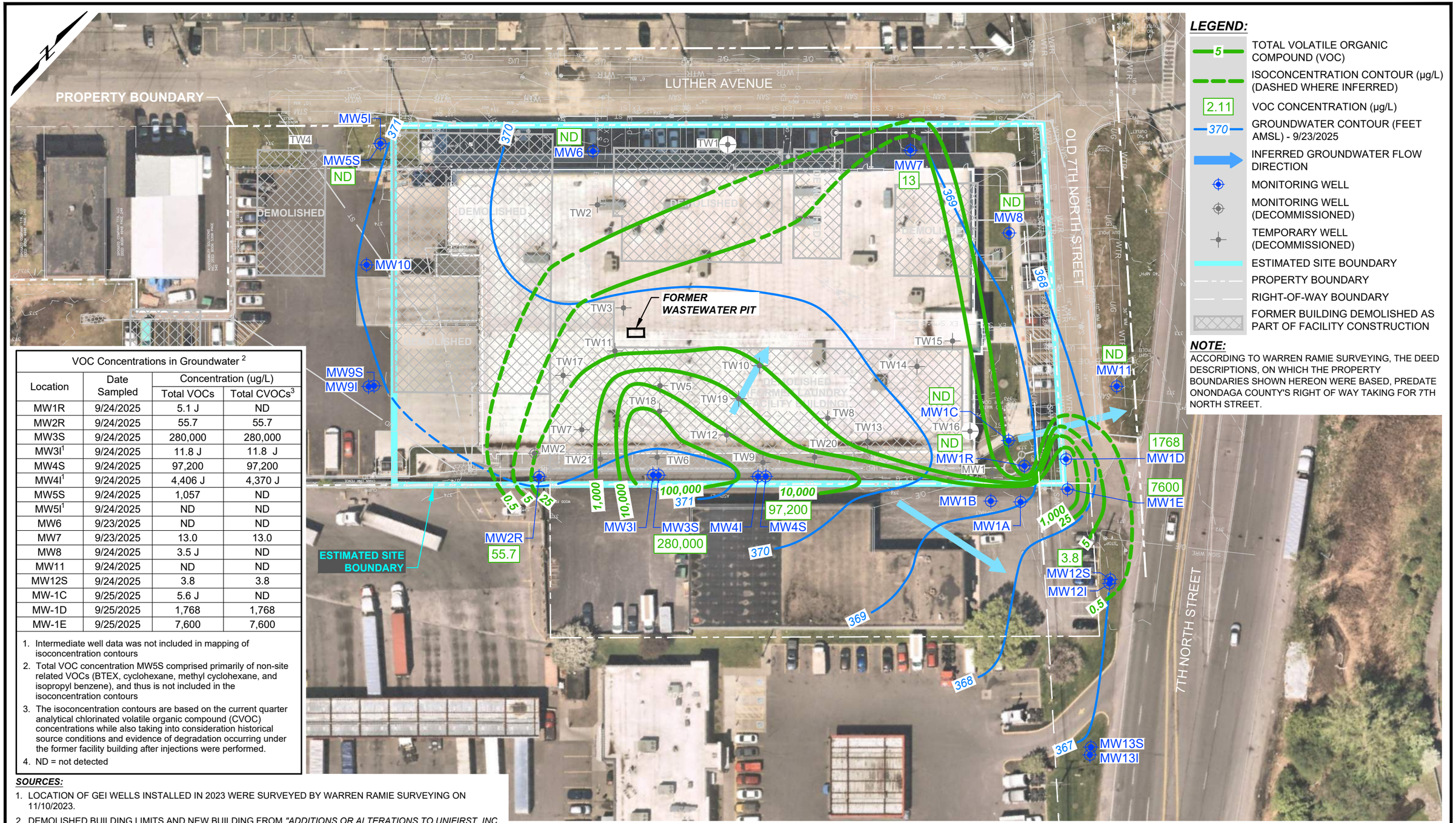
Unifirst Liverpool
Liverpool, New York

GEI Consultants

SUMMARIZED GROUNDWATER ANALYTICAL RESULTS - PFAS

Project 2001642 | February 2026 | Fig. 4b

HERIOT, PAULA B:\Working\UNIFIRST\2001642 Unifirst-Liverpool NY\00_CAD\Figures\QGWR\2025-Q3\4b_2001642_Q325GWR_GWPFS.dwg - 12/4/2025



VOC Concentrations in Groundwater [2]			
Location	Date Sampled	Concentration (ug/L)	
		Total VOCs	Total CVOCs[3]
MW1R	9/24/2025	5.1 J	ND
MW2R	9/24/2025	55.7	55.7
MW3S	9/24/2025	280,000	280,000
MW3I[1]	9/24/2025	11.8 J	11.8 J
MW4S	9/24/2025	97,200	97,200
MW4I[1]	9/24/2025	4,406 J	4,370 J
MW5S	9/24/2025	1,057	ND
MW5I[1]	9/24/2025	ND	ND
MW6	9/23/2025	ND	ND
MW7	9/23/2025	13.0	13.0
MW8	9/24/2025	3.5 J	ND
MW11	9/24/2025	ND	ND
MW12S	9/24/2025	3.8	3.8
MW-1C	9/25/2025	5.6 J	ND
MW-1D	9/25/2025	1,768	1,768
MW-1E	9/25/2025	7,600	7,600

1. Intermediate well data was not included in mapping of isoconcentration contours
2. Total VOC concentration MW5S comprised primarily of non-site related VOCs (BTEX, cyclohexane, methyl cyclohexane, and isopropyl benzene), and thus is not included in the isoconcentration contours
3. The isoconcentration contours are based on the current quarter analytical chlorinated volatile organic compound (CVOC) concentrations while also taking into consideration historical source conditions and evidence of degradation occurring under the former facility building after injections were performed.
4. ND = not detected

LEGEND:

- 5 — TOTAL VOLATILE ORGANIC COMPOUND (VOC)
- ISOCONCENTRATION CONTOUR (µg/L) (DASHED WHERE INFERRED)
- 2.11 — VOC CONCENTRATION (µg/L)
- 370 — GROUNDWATER CONTOUR (FEET AMSL) - 9/23/2025
- INFERRED GROUNDWATER FLOW DIRECTION
- MONITORING WELL
- MONITORING WELL (DECOMMISSIONED)
- TEMPORARY WELL (DECOMMISSIONED)
- ESTIMATED SITE BOUNDARY
- PROPERTY BOUNDARY
- RIGHT-OF-WAY BOUNDARY
- FORMER BUILDING DEMOLISHED AS PART OF FACILITY CONSTRUCTION

NOTE:

ACCORDING TO WARREN RAMIE SURVEYING, THE DEED DESCRIPTIONS, ON WHICH THE PROPERTY BOUNDARIES SHOWN HEREON WERE BASED, PREDATE ONONDAGA COUNTY'S RIGHT OF WAY TAKING FOR 7TH NORTH STREET.

SOURCES:

1. LOCATION OF GEI WELLS INSTALLED IN 2023 WERE SURVEYED BY WARREN RAMIE SURVEYING ON 11/10/2023.
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5. NEARMAP AERIAL ©2024 NEARMAP, IMAGERY DATED 05/07/2024, ACCESSED ON 08/15/2024.

0 60 120

SCALE: 1" = 60'

Quarterly Groundwater Monitoring Report - Q3 2025
113 7th North Street
Liverpool, New York

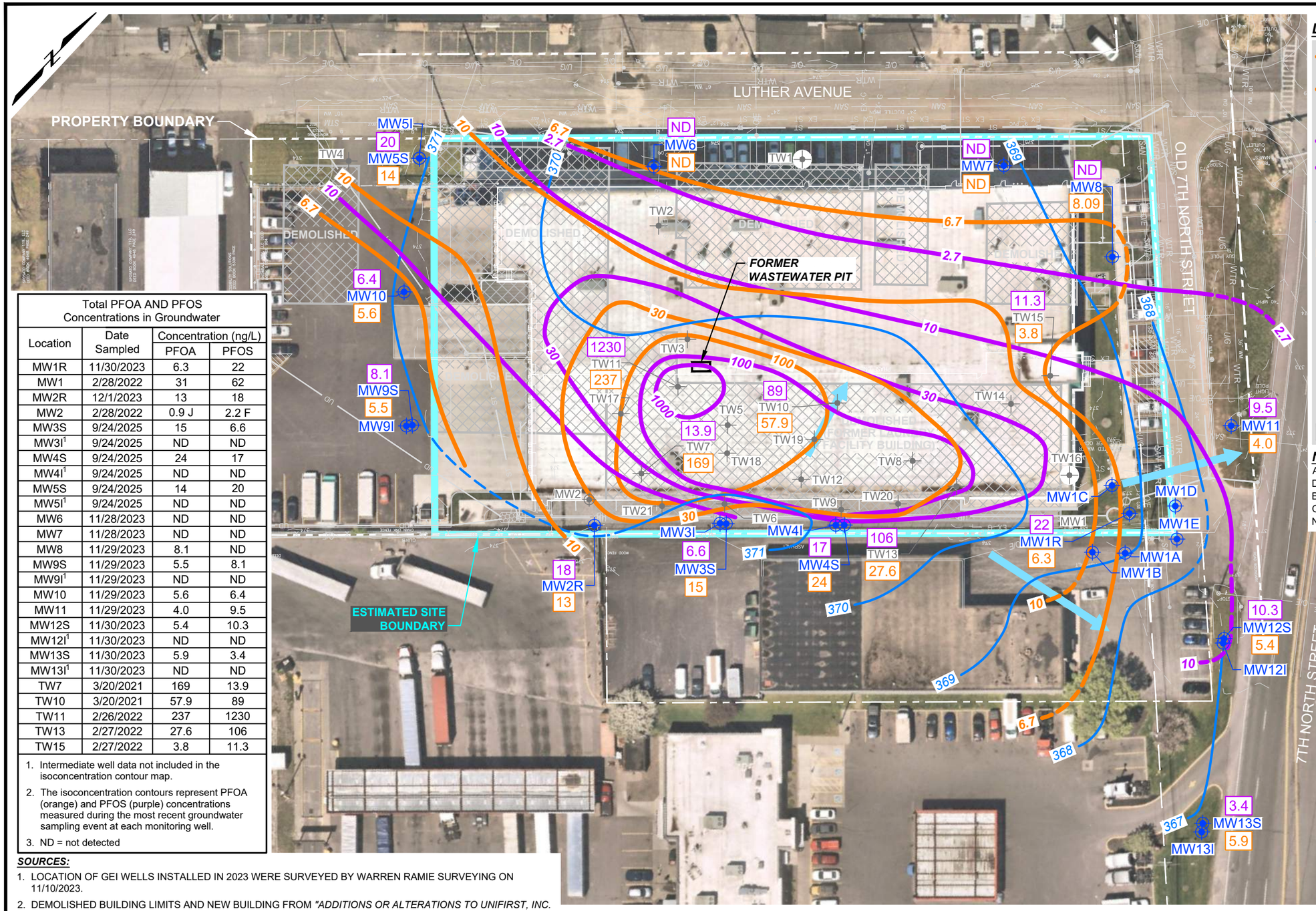
Unifirst Liverpool
Liverpool, New York

GEI Consultants

CVOC ISOCONCENTRATION MAP

Project 2001642 | February 2026 | Fig. 5a

---- B:\Working\UNIFIRST\2001642 Unifirst-Liverpool NY\00_CAD\Figures\QGWR\2025-Q3\5a_2001642_Q325GWR_TVOC.dwg - 11/11/2025



Total PFOA AND PFOS Concentrations in Groundwater			
Location	Date Sampled	Concentration (ng/L)	
		PFOA	PFOS
MW1R	11/30/2023	6.3	22
MW1	2/28/2022	31	62
MW2R	12/1/2023	13	18
MW2	2/28/2022	0.9 J	2.2 F
MW3S	9/24/2025	15	6.6
MW3I[1]	9/24/2025	ND	ND
MW4S	9/24/2025	24	17
MW4I[1]	9/24/2025	ND	ND
MW5S	9/24/2025	14	20
MW5I[1]	9/24/2025	ND	ND
MW6	11/28/2023	ND	ND
MW7	11/28/2023	ND	ND
MW8	11/29/2023	8.1	ND
MW9S	11/29/2023	5.5	8.1
MW9I[1]	11/29/2023	ND	ND
MW10	11/29/2023	5.6	6.4
MW11	11/29/2023	4.0	9.5
MW12S	11/30/2023	5.4	10.3
MW12I[1]	11/30/2023	ND	ND
MW13S	11/30/2023	5.9	3.4
MW13I[1]	11/30/2023	ND	ND
TW7	3/20/2021	169	13.9
TW10	3/20/2021	57.9	89
TW11	2/26/2022	237	1230
TW13	2/27/2022	27.6	106
TW15	2/27/2022	3.8	11.3

1. Intermediate well data not included in the isoconcentration contour map.
2. The isoconcentration contours represent PFOA (orange) and PFOS (purple) concentrations measured during the most recent groundwater sampling event at each monitoring well.
3. ND = not detected

LEGEND:

- 10 — PERFLUOROOCTANOIC ACID (PFOA) ISOCONCENTRATION CONTOUR (ng/L) (NYSDEC criterion = 6.7 ng/l) (DASHED WHERE INFERRED)
- 14 — PFOA CONCENTRATION (ng/L)
- 10 — PERFLUOROOCTANE SULFONATE (PFOS) ISOCONCENTRATION CONTOUR (ng/L) (DASHED WHERE INFERRED) (NYSDEC CRITERION = 2.7 ng/l)
- 20 — PFOS CONCENTRATION (ng/L)
- 370 — GROUNDWATER CONTOUR (FEET AMSL) - 9/23/2025
- INFERRED GROUNDWATER FLOW DIRECTION
- MONITORING WELL
- MONITORING WELL (DECOMMISSIONED)
- TEMPORARY WELL (DECOMMISSIONED)
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0 60 120

SCALE: 1" = 60'

Quarterly Groundwater Monitoring Report - Q3 2025
113 7th North Street
Liverpool, New York

Unifirst Liverpool
Liverpool, New York

GEI Consultants

PFOA AND PFOS ISOCONCENTRATION MAP

Project 2001642 | December 2025 | Fig. 5b

---- B:\Working\UNIFIRST\2001642 Unifirst-Liverpool NY\00_CAD\Figures\QGWR\2025-Q3\5b_2001642_Q325GWR_PFAS.dwg - 11/11/2025

Appendix A Waste Disposal Documentation

Please print or type. (Form designed for use on elite (12-pitch) typewriter.)

Form Approved. OMB No. 2050-0039

UNIFORM HAZARDOUS WASTE MANIFEST

1. Generator ID Number: NYR 000 250 514
2. Page 1 of: 1
3. Emergency Response Phone: (800) 888-4672
4. Manifest Tracking Number: 003465649 GBF

5. Generator's Name and Mailing Address
UNIFIRST LIVERPOOL
113 7TH NORTH STREET
LIVERPOOL, NY 13088
Generator's Phone: (315) 221-3521

Generator's Site Address (if different than mailing address)
103 LUTHER AVE
LIVERPOOL, NY 13088

6. Transporter 1 Company Name: ACV ENVIRONMENTAL SERVICES INC.
U.S. EPA ID Number: NJD 003 812 047

7. Transporter 2 Company Name: Environmental Soil Systems (New York) (NY) 716.685.6720
U.S. EPA ID Number: NYD 986 963 904

8. Designated Facility Name and Site Address: CHEMICAL SOLVENTS, INC.
1010 OLD DENISON AVE
CLEVELAND, OH 44109
U.S. EPA ID Number: OHD 980 897 656

Facility's Phone:

	9a. HM	9b. U.S. DOT Description (including Proper Shipping Name, Hazard Class, ID Number, and Packing Group (if any))	10. Containers No.	10. Containers Type	11. Total Quantity	12. Unit Wt./Vol.	13. Waste Codes		
GENERATOR	X	RQ, NA3082, Hazardous waste, liquid, n.o.s. (Trichloroethylene, Vinyl Chloride), 9, PGIII, ERG #171	003	DM	01200	P	D040	D043	
		2.							
		3.							
		4.							

14. Special Handling Instructions and Additional Information
9.156 / Hazardous Purge Water / PO: P2213.15503 (W:63.03.264100)

15. **GENERATOR'S/OFFEROR'S CERTIFICATION:** I hereby declare that the contents of this consignment are fully and accurately described above by the proper shipping name, and are classified, packaged, marked and labeled/placarded, and are in all respects in proper condition for transport according to applicable international and national governmental regulations. If export shipment and I am the Primary Exporter, I certify that the contents of this consignment conform to the terms of the attached EPA Acknowledgment of Consent.
I certify that the waste minimization statement identified in 40 CFR 262.27(a) (if I am a large quantity generator) or (b) (if I am a small quantity generator) is true.

Generator's/Offeror's Printed/Typed Name: CATHERINE MALAGRIDA
Signature: C Malagrida
Month / Day / Year:

INT'L
16. International Shipments ☐ Import to U.S. ☐ Export from U.S. Port of entry/exit:
Transporter signature (for exports only): Date leaving U.S.:

TRANSPORTER
17. Transporter Acknowledgment of Receipt of Materials
Transporter 1 Printed/Typed Name: [illegible] Signature: [illegible] Month / Day / Year:
Transporter 2 Printed/Typed Name: [illegible] Signature: [illegible] Month: 11 Day: 25 Year: 13

DESIGNATED FACILITY
18. Discrepancy
18a. Discrepancy Indication Space ☐ Quantity ☐ Type ☐ Residue ☐ Partial Rejection ☐ Full Rejection
Manifest Reference Number:
18b. Alternate Facility (or Generator) U.S. EPA ID Number
Facility's Phone:
18c. Signature of Alternate Facility (or Generator) Month / Day / Year
19. Hazardous Waste Report Management Method Codes (i.e., codes for hazardous waste treatment, disposal, and recycling systems)
1. 2. 3. 4.
20. Designated Facility Owner or Operator: Certification of receipt of hazardous materials covered by the manifest except as noted in Item 18a
Printed/Typed Name: [illegible] Signature: [illegible] Month: 11 Day: 26 Year:

EPA Form 8700-22 (Rev. 3-05) Previous editions are obsolete.

GENERATOR'S INITIAL COPY

Appendix B Laboratory Report

ANALYTICAL REPORT

PREPARED FOR

Attn: Ms. Abigal Jock
GEI Consultants Inc
950 Danby Road
Suite 201-F
Ithaca, New York 14850

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JOB DESCRIPTION

Unifirst – Liverpool, NY

JOB NUMBER

480-232963-1

Eurofins Buffalo
10 Hazelwood Drive
Amherst NY 14228-2298

See page two for job notes and contact information.

Eurofins Buffalo

Job Notes

This report may not be reproduced except in full, and with written approval from the laboratory. The results relate only to the samples tested. For questions please contact the Project Manager at the e-mail address or telephone number listed on this page.

The test results in this report relate only to the samples as received by the laboratory and will meet all requirements of the methodology, with any exceptions noted. This report shall not be reproduced except in full, without the express written approval of the laboratory. All questions should be directed to the Eurofins Environment Testing Northeast, LLC Project Manager.

Authorization

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Authorized for release by
Kavya Boyina, Project Manager
kavya.boyina@et.eurofinsus.com
Designee for
John Schove, Project Manager II
John.Schove@et.eurofinsus.com
(716)504-9838

Table of Contents

Definitions/Glossary

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Qualifiers

GC/MS VOA

Qualifier	Qualifier Description
F1	MS and/or MSD recovery exceeds control limits.
J	Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.
S1+	Surrogate recovery exceeds control limits, high biased.

GC VOA

Qualifier	Qualifier Description
J	Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.

LCMS

Qualifier	Qualifier Description
*-	LCS and/or LCSD is outside acceptance limits, low biased.
*5+	Isotope dilution analyte is outside acceptance limits, high biased.
I	Ion ratio outside limits
J	Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.

Glossary

Abbreviation	These commonly used abbreviations may or may not be present in this report.
☼	Listed under the "D" column to designate that the result is reported on a dry weight basis
%R	Percent Recovery
CFL	Contains Free Liquid
CFU	Colony Forming Unit
CNF	Contains No Free Liquid
DER	Duplicate Error Ratio (normalized absolute difference)
Dil Fac	Dilution Factor
DL	Detection Limit (DoD/DOE)
DL, RA, RE, IN	Indicates a Dilution, Re-analysis, Re-extraction, or additional Initial metals/anion analysis of the sample
DLC	Decision Level Concentration (Radiochemistry)
EDL	Estimated Detection Limit (Dioxin)
LOD	Limit of Detection (DoD/DOE)
LOQ	Limit of Quantitation (DoD/DOE)
MCL	EPA recommended "Maximum Contaminant Level"
MDA	Minimum Detectable Activity (Radiochemistry)
MDC	Minimum Detectable Concentration (Radiochemistry)
MDL	Method Detection Limit
ML	Minimum Level (Dioxin)
MPN	Most Probable Number
MQL	Method Quantitation Limit
NC	Not Calculated
ND	Not Detected at the reporting limit (or MDL or EDL if shown)
NEG	Negative / Absent
POS	Positive / Present
PQL	Practical Quantitation Limit
PRES	Presumptive
QC	Quality Control
RER	Relative Error Ratio (Radiochemistry)
RL	Reporting Limit or Requested Limit (Radiochemistry)
RPD	Relative Percent Difference, a measure of the relative difference between two points
TEF	Toxicity Equivalent Factor (Dioxin)
TEQ	Toxicity Equivalent Quotient (Dioxin)
TNTC	Too Numerous To Count

Job ID: 480-232963-1 Eurofins Buffalo

Job Narrative
480-232963-1

The analytical test results presented in this report meet all requirements of the associated regulatory program listed on the Accreditation/Certification Summary Page, unless otherwise noted. Data qualifiers and/or narrative comments are included to explain any exceptions, if applicable. Regulated compliance samples (e.g. SDWA, NPDES) must comply with associated agency requirements/permits.

- Matrix-specific batch QC (e.g., MS, MSD, SD) may not be reported when insufficient sample volume is available or when site-specific QC samples are not submitted. In such cases, a Laboratory Control Sample Duplicate (LCSD) may be analyzed to provide precision data for the batch.
- For samples analyzed using surrogate and/or isotope dilution analytes, any recoveries falling outside of established acceptance criteria are re-prepared and/or re-analyzed to confirm results, unless the deviation is due to sample dilution or otherwise explained in the case narrative.

Receipt
The samples were received on 9/26/2025 10:00 AM. Unless otherwise noted below, the samples arrived in good condition, and, where required, properly preserved and on ice. The temperatures of the 2 coolers at receipt time were 2.9°C and 3.2°C.

GC/MS VOA
Method 8260C: The following sample was diluted due to the nature of the sample matrix: MW1R (480-232963-1). Elevated reporting limits (RLs) are provided.

Method 8260C: The following samples were diluted to bring the concentration of target analytes within the calibration range: MW3S (480-232963-3), MW4S (480-232963-5), MW4I (480-232963-6), MW5S (480-232963-7), MW1D (480-232963-15), MW1E (480-232963-16), (480-232963-F-6 MS) and (480-232963-F-6 MSD). Elevated reporting limits (RLs) are provided.

Method 8260C: Surrogate recovery for the following sample was outside control limits: MW5S (480-232963-7). Evidence of matrix interference is present; therefore, re-analysis was not performed.

Method 8260C: The continuing calibration verification (CCV) associated with batch 758407 recovered above the upper control limit for Trichlorofluoromethane. The samples associated with this CCV were non-detects for the affected analytes; therefore, the data have been reported. The associated samples are:MW1R (480-232963-1), MW2R (480-232963-2), MW3S (480-232963-3), MW4S (480-232963-5), MW4I (480-232963-6), MW5S (480-232963-7), MW5I (480-232963-8), MW6 (480-232963-9), MW7 (480-232963-10), MW8 (480-232963-11), MW11 (480-232963-12), MW12S (480-232963-13), MW1C (480-232963-14), MW1D (480-232963-15), MW1E (480-232963-16) and TB (480-232963-17).

Method 8260C: The matrix spike / matrix spike duplicate (MS/MSD) recoveries for analytical batch 480-758407 were outside control limits for one or more analytes. See QC Sample Results for detail. Sample matrix interference and/or non-homogeneity are suspected because the associated laboratory control sample (LCS) recovery is within acceptance limits.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/ Glossary page.

GC VOA
Method RSK_175: The following samples were diluted to bring the concentration of target analytes within the calibration range: MW1R (480-232963-1), MW2R (480-232963-2), MW3S (480-232963-3), MW4S (480-232963-5) and MW4I (480-232963-6). Elevated reporting limits (RLs) are provided.

Method RSK_175: The %RPD between the primary and confirmation detector exceeded 40% for Ethene for the following sample: MW4S (480-232963-5). The lower value(s) has been reported and qualified in accordance with the laboratory's SOP.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/ Glossary page.

PFAS
Method 1633A: Analyte Perfluoroheptanesulfonic acid (PFHpS) and Perfluorononanoic acid (PFNA) was marked as non-detect for sample MW5S (480-232963-7) due to a detection below RL and failing ion ratio.

Method 1633A: The labeled isotopes 13C2 4:2 FTS are outside QC acceptance limits. Since the recovery in sample MW5S (480-232963-7) is less than 350%, the data is reported.

Eurofins Buffalo

Method 1633A: The transition mass ratio for Perfluorooctanesulfonic acid (PFOS) in sample MW3S (480-232963-3) in analytical batch 410-710552 was outside of the established ratio limits. However, analyst judgment was used to positively identify the analyte. The analyte is flagged appropriately to identify this exceedance.

Method 1633A: The labeled isotopes 13C2 4:2 FTS are outside QC acceptance limits. Since the recovery in sample MW3S (480-232963-3) is less than 350%, the data is reported.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/ Glossary page.

Metals
No additional analytical or quality issues were noted, other than those described above or in the Definitions/ Glossary page.

General Chemistry
Method 300.0_28D: The following samples were diluted due to the nature of the sample matrix: MW1R (480-232963-1), MW2R (480-232963-2), MW3S (480-232963-3) and MW3I (480-232963-4). Elevated reporting limits (RLs) are provided.

Method 300.0_28D: The following samples were diluted due to the nature of the sample matrix: MW4S (480-232963-5) and MW4I (480-232963-6). Elevated reporting limits (RLs) are provided.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/ Glossary page.

Detection Summary

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW1R — Lab Sample ID: 480-232963-1

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Xylenes, Total	5.1	J	10	3.3	ug/L	5		8260C	Total/NA
Ethane	5.2	J	7.5	1.5	ug/L	1		RSK-175	Total/NA
Methane - DL	1800		88	22	ug/L	22		RSK-175	Total/NA
Iron	0.61		0.050	0.019	mg/L	1		6010D	Total/NA
Iron, Dissolved	0.21		0.050	0.019	mg/L	1		6010D	Dissolved
Sulfate	30.6		10.0	1.7	mg/L	5		300.0	Total/NA

Client Sample ID: MW2R — Lab Sample ID: 480-232963-2

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
cis-1,2-Dichloroethene	39		1.0	0.81	ug/L	1		8260C	Total/NA
Tetrachloroethene	6.7		1.0	0.36	ug/L	1		8260C	Total/NA
trans-1,2-Dichloroethene	1.7		1.0	0.90	ug/L	1		8260C	Total/NA
Trichloroethene	4.4		1.0	0.46	ug/L	1		8260C	Total/NA
Vinyl chloride	3.9		1.0	0.90	ug/L	1		8260C	Total/NA
Methane - DL	890		88	22	ug/L	22		RSK-175	Total/NA
Iron	1.0		0.050	0.019	mg/L	1		6010D	Total/NA
Iron, Dissolved	1.2		0.050	0.019	mg/L	1		6010D	Dissolved
Sulfate	27.1		10.0	1.7	mg/L	5		300.0	Total/NA

Client Sample ID: MW3S — Lab Sample ID: 480-232963-3

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
cis-1,2-Dichloroethene	160000		2000	1600	ug/L	2000		8260C	Total/NA
Vinyl chloride	120000		2000	1800	ug/L	2000		8260C	Total/NA
Methane	6800		440	110	ug/L	110		RSK-175	Total/NA
Ethane	1100		830	170	ug/L	110		RSK-175	Total/NA
Ethene	29000		770	170	ug/L	110		RSK-175	Total/NA
Perfluorobutanesulfonic acid (PFBS)	4.5		1.7	0.41	ng/L	1		1633A	Total/NA
Perfluorobutanoic acid (PFBA)	5.1		3.3	0.91	ng/L	1		1633A	Total/NA
Perfluoroheptanoic acid (PFHpA)	5.7		1.7	0.66	ng/L	1		1633A	Total/NA
Perfluorohexanesulfonic acid (PFHxS)	3.4		1.7	0.66	ng/L	1		1633A	Total/NA
Perfluorohexanoic acid (PFHxA)	12		1.7	0.41	ng/L	1		1633A	Total/NA
Perfluorooctanesulfonic acid (PFOS)	6.6	I	1.7	1.1	ng/L	1		1633A	Total/NA
Perfluorooctanoic acid (PFOA)	15		1.7	0.74	ng/L	1		1633A	Total/NA
Perfluoropentanesulfonic acid (PFPeS)	1.7		1.7	0.41	ng/L	1		1633A	Total/NA
Perfluoropentanoic acid (PFPeA)	7.3		1.7	0.50	ng/L	1		1633A	Total/NA
Iron	91.6		0.050	0.019	mg/L	1		6010D	Total/NA
Iron, Dissolved	88.0		0.050	0.019	mg/L	1		6010D	Dissolved
Sulfate	25.0		20.0	3.5	mg/L	10		300.0	Total/NA

Client Sample ID: MW3I — Lab Sample ID: 480-232963-4

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
cis-1,2-Dichloroethene	2.9		1.0	0.81	ug/L	1		8260C	Total/NA
Trichloroethene	0.53	J	1.0	0.46	ug/L	1		8260C	Total/NA
Vinyl chloride	8.4		1.0	0.90	ug/L	1		8260C	Total/NA
Methane	140		4.0	1.0	ug/L	1		RSK-175	Total/NA
Perfluorobutanoic acid (PFBA)	0.97	J	3.1	0.85	ng/L	1		1633A	Total/NA
Iron	1.1		0.050	0.019	mg/L	1		6010D	Total/NA
Iron, Dissolved	1.0		0.050	0.019	mg/L	1		6010D	Dissolved
Sulfate	79.9		10.0	1.7	mg/L	5		300.0	Total/NA

This Detection Summary does not include radiochemical test results.

Eurofins Buffalo

Detection Summary

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW4S — Lab Sample ID: 480-232963-5

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
cis-1,2-Dichloroethene	8100		2000	1600	ug/L	2000		8260C	Total/NA
Tetrachloroethene	85000		2000	720	ug/L	2000		8260C	Total/NA
Trichloroethene	4100		2000	920	ug/L	2000		8260C	Total/NA
Ethane	390		7.5	1.5	ug/L	1		RSK-175	Total/NA
Ethene	13		7.0	1.5	ug/L	1		RSK-175	Total/NA
Methane - DL	4600		180	44	ug/L	44		RSK-175	Total/NA
Perfluorobutanesulfonic acid (PFBS)	4.4		1.5	0.39	ng/L	1		1633A	Total/NA
Perfluorobutanoic acid (PFBA)	12		3.1	0.85	ng/L	1		1633A	Total/NA
Perfluorodecanoic acid (PFDA)	1.5		1.5	0.39	ng/L	1		1633A	Total/NA
Perfluoroheptanoic acid (PFHpA)	11		1.5	0.62	ng/L	1		1633A	Total/NA
Perfluorohexanesulfonic acid (PFHxS)	3.7		1.5	0.62	ng/L	1		1633A	Total/NA
Perfluorohexanoic acid (PFHxA)	18		1.5	0.39	ng/L	1		1633A	Total/NA
Perfluorononanoic acid (PFNA)	4.1		1.5	0.39	ng/L	1		1633A	Total/NA
Perfluorooctanesulfonic acid (PFOS)	17		1.5	1.0	ng/L	1		1633A	Total/NA
Perfluorooctanoic acid (PFOA)	24		1.5	0.69	ng/L	1		1633A	Total/NA
Perfluoropentanesulfonic acid (PFPeS)	0.80	J	1.5	0.39	ng/L	1		1633A	Total/NA
Perfluoropentanoic acid (PFPeA)	25		1.5	0.46	ng/L	1		1633A	Total/NA
Iron	1.5		0.050	0.019	mg/L	1		6010D	Total/NA
Iron, Dissolved	0.081		0.050	0.019	mg/L	1		6010D	Dissolved
Sulfate	131		40.0	7.0	mg/L	20		300.0	Total/NA

Client Sample ID: MW4I — Lab Sample ID: 480-232963-6

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
1,1-Dichloroethene	36	J	40	12	ug/L	40		8260C	Total/NA
cis-1,2-Dichloroethene	3800	F1	40	32	ug/L	40		8260C	Total/NA
Tetrachloroethene	20	J	40	14	ug/L	40		8260C	Total/NA
Vinyl chloride	550		40	36	ug/L	40		8260C	Total/NA
Ethane	15		7.5	1.5	ug/L	1		RSK-175	Total/NA
Ethene	54		7.0	1.5	ug/L	1		RSK-175	Total/NA
Methane - DL	1400		88	22	ug/L	22		RSK-175	Total/NA
Perfluorobutanoic acid (PFBA)	1.1	J	3.0	0.83	ng/L	1		1633A	Total/NA
Iron	1.2		0.050	0.019	mg/L	1		6010D	Total/NA
Iron, Dissolved	1.3		0.050	0.019	mg/L	1		6010D	Dissolved
Sulfate	128		10.0	1.7	mg/L	5		300.0	Total/NA

Client Sample ID: MW5S — Lab Sample ID: 480-232963-7

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
2-Butanone (MEK)	34	J	100	13	ug/L	10		8260C	Total/NA
Acetone	86	J	100	30	ug/L	10		8260C	Total/NA
Benzene	18		10	4.1	ug/L	10		8260C	Total/NA
Cyclohexane	370		10	1.8	ug/L	10		8260C	Total/NA
Ethylbenzene	260		10	7.4	ug/L	10		8260C	Total/NA
Isopropylbenzene	110		10	7.9	ug/L	10		8260C	Total/NA
Methylcyclohexane	150		10	1.6	ug/L	10		8260C	Total/NA
Toluene	6.4	J	10	5.1	ug/L	10		8260C	Total/NA
Xylenes, Total	23		20	6.6	ug/L	10		8260C	Total/NA
Perfluorobutanesulfonic acid (PFBS)	0.85	J	1.8	0.44	ng/L	1		1633A	Total/NA
Perfluorobutanoic acid (PFBA)	9.1		3.5	0.97	ng/L	1		1633A	Total/NA
Perfluoroheptanoic acid (PFHpA)	2.6		1.8	0.70	ng/L	1		1633A	Total/NA

This Detection Summary does not include radiochemical test results.

Eurofins Buffalo

Client Sample ID: MW5S (Continued) — Lab Sample ID: 480-232963-7

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorohexanesulfonic acid (PFHxS)	1.4	J	1.8	0.70	ng/L	1		1633A	Total/NA
Perfluorohexanoic acid (PFHxA)	5.0		1.8	0.44	ng/L	1		1633A	Total/NA
Perfluorooctanesulfonamide (PFOSA)	0.48	J	1.8	0.44	ng/L	1		1633A	Total/NA
Perfluorooctanesulfonic acid (PFOS)	20		1.8	1.1	ng/L	1		1633A	Total/NA
Perfluorooctanoic acid (PFOA)	14		1.8	0.79	ng/L	1		1633A	Total/NA
Perfluoropentanoic acid (PFPeA)	8.1		1.8	0.53	ng/L	1		1633A	Total/NA

Client Sample ID: MW5I — Lab Sample ID: 480-232963-8

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	39		3.3	0.82	ng/L	1		1633A	Total/NA
Perfluorobutanoic acid (PFBA)	1.7	J	3.3	0.90	ng/L	1		1633A	Total/NA
Perfluoroheptanoic acid (PFHpA)	1.5	J	1.6	0.66	ng/L	1		1633A	Total/NA
Perfluorohexanesulfonic acid (PFHxS)	1.1	J	1.6	0.66	ng/L	1		1633A	Total/NA
Perfluorohexanoic acid (PFHxA)	4.4		1.6	0.41	ng/L	1		1633A	Total/NA
Perfluoropentanoic acid (PFPeA)	4.9		1.6	0.49	ng/L	1		1633A	Total/NA

Client Sample ID: MW6 — Lab Sample ID: 480-232963-9

No Detections.

Client Sample ID: MW7 — Lab Sample ID: 480-232963-10

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Vinyl chloride	13		1.0	0.90	ug/L	1		8260C	Total/NA

Client Sample ID: MW8 — Lab Sample ID: 480-232963-11

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Acetone	3.5	J	10	3.0	ug/L	1		8260C	Total/NA

Client Sample ID: MW11 — Lab Sample ID: 480-232963-12

No Detections.

Client Sample ID: MW12S — Lab Sample ID: 480-232963-13

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
cis-1,2-Dichloroethene	2.3		1.0	0.81	ug/L	1		8260C	Total/NA
Vinyl chloride	1.5		1.0	0.90	ug/L	1		8260C	Total/NA

Client Sample ID: MW1C — Lab Sample ID: 480-232963-14

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Acetone	5.6	J	10	3.0	ug/L	1		8260C	Total/NA

Client Sample ID: MW1D — Lab Sample ID: 480-232963-15

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
cis-1,2-Dichloroethene	990		20	16	ug/L	20		8260C	Total/NA
Tetrachloroethene	540		20	7.2	ug/L	20		8260C	Total/NA
Trichloroethene	210		20	9.2	ug/L	20		8260C	Total/NA
Vinyl chloride	28		20	18	ug/L	20		8260C	Total/NA

This Detection Summary does not include radiochemical test results.

Detection Summary

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW1E — Lab Sample ID: 480-232963-16

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
cis-1,2-Dichloroethene	2900		100	81	ug/L	100		8260C	Total/NA
Tetrachloroethene	3600		100	36	ug/L	100		8260C	Total/NA
Trichloroethene	1100		100	46	ug/L	100		8260C	Total/NA

Client Sample ID: TB — Lab Sample ID: 480-232963-17

No Detections.

This Detection Summary does not include radiochemical test results.

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW1R
Date Collected: 09/24/25 14:30
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-1
Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		5.0	4.1	ug/L			10/01/25 01:19	5
1,1,2,2-Tetrachloroethane	ND		5.0	1.1	ug/L			10/01/25 01:19	5
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		5.0	1.6	ug/L			10/01/25 01:19	5
1,1,2-Trichloroethane	ND		5.0	1.2	ug/L			10/01/25 01:19	5
1,1-Dichloroethane	ND		5.0	1.9	ug/L			10/01/25 01:19	5
1,1-Dichloroethene	ND		5.0	1.5	ug/L			10/01/25 01:19	5
1,2,4-Trichlorobenzene	ND		5.0	2.1	ug/L			10/01/25 01:19	5
1,2-Dibromo-3-Chloropropane	ND		5.0	2.0	ug/L			10/01/25 01:19	5
1,2-Dibromoethane	ND		5.0	3.7	ug/L			10/01/25 01:19	5
1,2-Dichlorobenzene	ND		5.0	4.0	ug/L			10/01/25 01:19	5
1,2-Dichloroethane	ND		5.0	1.1	ug/L			10/01/25 01:19	5
1,2-Dichloropropane	ND		5.0	3.6	ug/L			10/01/25 01:19	5
1,3-Dichlorobenzene	ND		5.0	3.9	ug/L			10/01/25 01:19	5
1,4-Dichlorobenzene	ND		5.0	4.2	ug/L			10/01/25 01:19	5
2-Butanone (MEK)	ND		50	6.6	ug/L			10/01/25 01:19	5
2-Hexanone	ND		25	6.2	ug/L			10/01/25 01:19	5
4-Methyl-2-pentanone (MIBK)	ND		25	11	ug/L			10/01/25 01:19	5
Acetone	ND		50	15	ug/L			10/01/25 01:19	5
Benzene	ND		5.0	2.1	ug/L			10/01/25 01:19	5
Bromodichloromethane	ND		5.0	2.0	ug/L			10/01/25 01:19	5
Bromoform	ND		5.0	1.3	ug/L			10/01/25 01:19	5
Bromomethane	ND		5.0	3.5	ug/L			10/01/25 01:19	5
Carbon disulfide	ND		5.0	0.95	ug/L			10/01/25 01:19	5
Carbon tetrachloride	ND		5.0	1.4	ug/L			10/01/25 01:19	5
Chlorobenzene	ND		5.0	3.8	ug/L			10/01/25 01:19	5
Chloroethane	ND		5.0	1.6	ug/L			10/01/25 01:19	5
Chloroform	ND		5.0	1.7	ug/L			10/01/25 01:19	5
Chloromethane	ND		5.0	1.8	ug/L			10/01/25 01:19	5
cis-1,2-Dichloroethene	ND		5.0	4.1	ug/L			10/01/25 01:19	5
cis-1,3-Dichloropropene	ND		5.0	1.8	ug/L			10/01/25 01:19	5
Cyclohexane	ND		5.0	0.90	ug/L			10/01/25 01:19	5
Dibromochloromethane	ND		5.0	1.6	ug/L			10/01/25 01:19	5
Dichlorodifluoromethane	ND		5.0	3.4	ug/L			10/01/25 01:19	5
Ethylbenzene	ND		5.0	3.7	ug/L			10/01/25 01:19	5
Isopropylbenzene	ND		5.0	4.0	ug/L			10/01/25 01:19	5
Methyl acetate	ND		13	6.5	ug/L			10/01/25 01:19	5
Methyl tert-butyl ether	ND		5.0	0.80	ug/L			10/01/25 01:19	5
Methylcyclohexane	ND		5.0	0.80	ug/L			10/01/25 01:19	5
Methylene Chloride	ND		5.0	2.2	ug/L			10/01/25 01:19	5
Styrene	ND		5.0	3.7	ug/L			10/01/25 01:19	5
Tetrachloroethene	ND		5.0	1.8	ug/L			10/01/25 01:19	5
Toluene	ND		5.0	2.6	ug/L			10/01/25 01:19	5
trans-1,2-Dichloroethene	ND		5.0	4.5	ug/L			10/01/25 01:19	5
trans-1,3-Dichloropropene	ND		5.0	1.9	ug/L			10/01/25 01:19	5
Trichloroethene	ND		5.0	2.3	ug/L			10/01/25 01:19	5
Trichlorofluoromethane	ND		5.0	4.4	ug/L			10/01/25 01:19	5
Vinyl chloride	ND		5.0	4.5	ug/L			10/01/25 01:19	5
Xylenes, Total	**5.1**	**J**	10	3.3	ug/L			10/01/25 01:19	5

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW1R

Date Collected: 09/24/25 14:30
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-1

Matrix: Water

Surrogate	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4 (Surr)	*107*		*77 - 120*		*10/01/25 01:19*	*5*
4-Bromofluorobenzene (Surr)	*107*		*73 - 120*		*10/01/25 01:19*	*5*
Dibromofluoromethane (Surr)	*107*		*75 - 123*		*10/01/25 01:19*	*5*
Toluene-d8 (Surr)	*101*		*80 - 120*		*10/01/25 01:19*	*5*

Method: RSK-175 - Dissolved Gases (GC)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Ethane	**5.2**	**J**	7.5	1.5	ug/L			10/01/25 10:47	1
Ethene	ND		7.0	1.5	ug/L			10/01/25 10:47	1

Method: RSK-175 - Dissolved Gases (GC) - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	**1800**		88	22	ug/L			10/01/25 12:40	22

Method: SW846 6010D - Metals (ICP)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron	**0.61**		0.050	0.019	mg/L		09/29/25 08:35	09/29/25 15:41	1

Method: SW846 6010D - Metals (ICP) - Dissolved

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron, Dissolved	**0.21**		0.050	0.019	mg/L		09/29/25 08:34	09/29/25 23:58	1

General Chemistry

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Sulfate (EPA 300.0)	**30.6**		10.0	1.7	mg/L			09/30/25 01:42	5

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW2R

Date Collected: 09/24/25 12:30
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-2

Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		1.0	0.82	ug/L			10/01/25 01:42	1
1,1,2,2-Tetrachloroethane	ND		1.0	0.21	ug/L			10/01/25 01:42	1
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1.0	0.31	ug/L			10/01/25 01:42	1
1,1,2-Trichloroethane	ND		1.0	0.23	ug/L			10/01/25 01:42	1
1,1-Dichloroethane	ND		1.0	0.38	ug/L			10/01/25 01:42	1
1,1-Dichloroethene	ND		1.0	0.29	ug/L			10/01/25 01:42	1
1,2,4-Trichlorobenzene	ND		1.0	0.41	ug/L			10/01/25 01:42	1
1,2-Dibromo-3-Chloropropane	ND		1.0	0.39	ug/L			10/01/25 01:42	1
1,2-Dibromoethane	ND		1.0	0.73	ug/L			10/01/25 01:42	1
1,2-Dichlorobenzene	ND		1.0	0.79	ug/L			10/01/25 01:42	1
1,2-Dichloroethane	ND		1.0	0.21	ug/L			10/01/25 01:42	1
1,2-Dichloropropane	ND		1.0	0.72	ug/L			10/01/25 01:42	1
1,3-Dichlorobenzene	ND		1.0	0.78	ug/L			10/01/25 01:42	1
1,4-Dichlorobenzene	ND		1.0	0.84	ug/L			10/01/25 01:42	1
2-Butanone (MEK)	ND		10	1.3	ug/L			10/01/25 01:42	1
2-Hexanone	ND		5.0	1.2	ug/L			10/01/25 01:42	1
4-Methyl-2-pentanone (MIBK)	ND		5.0	2.1	ug/L			10/01/25 01:42	1
Acetone	ND		10	3.0	ug/L			10/01/25 01:42	1
Benzene	ND		1.0	0.41	ug/L			10/01/25 01:42	1
Bromodichloromethane	ND		1.0	0.39	ug/L			10/01/25 01:42	1
Bromoform	ND		1.0	0.26	ug/L			10/01/25 01:42	1
Bromomethane	ND		1.0	0.69	ug/L			10/01/25 01:42	1
Carbon disulfide	ND		1.0	0.19	ug/L			10/01/25 01:42	1
Carbon tetrachloride	ND		1.0	0.27	ug/L			10/01/25 01:42	1
Chlorobenzene	ND		1.0	0.75	ug/L			10/01/25 01:42	1
Chloroethane	ND		1.0	0.32	ug/L			10/01/25 01:42	1
Chloroform	ND		1.0	0.34	ug/L			10/01/25 01:42	1
Chloromethane	ND		1.0	0.35	ug/L			10/01/25 01:42	1
cis-1,2-Dichloroethene	**39**		1.0	0.81	ug/L			10/01/25 01:42	1
cis-1,3-Dichloropropene	ND		1.0	0.36	ug/L			10/01/25 01:42	1
Cyclohexane	ND		1.0	0.18	ug/L			10/01/25 01:42	1
Dibromochloromethane	ND		1.0	0.32	ug/L			10/01/25 01:42	1
Dichlorodifluoromethane	ND		1.0	0.68	ug/L			10/01/25 01:42	1
Ethylbenzene	ND		1.0	0.74	ug/L			10/01/25 01:42	1
Isopropylbenzene	ND		1.0	0.79	ug/L			10/01/25 01:42	1
Methyl acetate	ND		2.5	1.3	ug/L			10/01/25 01:42	1
Methyl tert-butyl ether	ND		1.0	0.16	ug/L			10/01/25 01:42	1
Methylcyclohexane	ND		1.0	0.16	ug/L			10/01/25 01:42	1
Methylene Chloride	ND		1.0	0.44	ug/L			10/01/25 01:42	1
Styrene	ND		1.0	0.73	ug/L			10/01/25 01:42	1
Tetrachloroethene	**6.7**		1.0	0.36	ug/L			10/01/25 01:42	1
Toluene	ND		1.0	0.51	ug/L			10/01/25 01:42	1
trans-1,2-Dichloroethene	**1.7**		1.0	0.90	ug/L			10/01/25 01:42	1
trans-1,3-Dichloropropene	ND		1.0	0.37	ug/L			10/01/25 01:42	1
Trichloroethene	**4.4**		1.0	0.46	ug/L			10/01/25 01:42	1
Trichlorofluoromethane	ND		1.0	0.88	ug/L			10/01/25 01:42	1
Vinyl chloride	**3.9**		1.0	0.90	ug/L			10/01/25 01:42	1
Xylenes, Total	ND		2.0	0.66	ug/L			10/01/25 01:42	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW2R

Date Collected: 09/24/25 12:30
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-2

Matrix: Water

Surrogate	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4 (Surr)	*106*		*77 - 120*		*10/01/25 01:42*	*1*
4-Bromofluorobenzene (Surr)	*106*		*73 - 120*		*10/01/25 01:42*	*1*
Dibromofluoromethane (Surr)	*109*		*75 - 123*		*10/01/25 01:42*	*1*
Toluene-d8 (Surr)	*104*		*80 - 120*		*10/01/25 01:42*	*1*

Method: RSK-175 - Dissolved Gases (GC)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Ethane	ND		7.5	1.5	ug/L			10/01/25 11:06	1
Ethene	ND		7.0	1.5	ug/L			10/01/25 11:06	1

Method: RSK-175 - Dissolved Gases (GC) - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	**890**		88	22	ug/L			10/01/25 12:59	22

Method: SW846 6010D - Metals (ICP)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron	**1.0**		0.050	0.019	mg/L		09/29/25 08:35	09/29/25 15:43	1

Method: SW846 6010D - Metals (ICP) - Dissolved

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron, Dissolved	**1.2**		0.050	0.019	mg/L		09/29/25 08:34	09/30/25 00:15	1

General Chemistry

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Sulfate (EPA 300.0)	**27.1**		10.0	1.7	mg/L			09/30/25 01:56	5

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW3S
Date Collected: 09/24/25 14:15
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-3
Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		2000	1600	ug/L			10/01/25 02:05	2000
1,1,2,2-Tetrachloroethane	ND		2000	420	ug/L			10/01/25 02:05	2000
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		2000	620	ug/L			10/01/25 02:05	2000
1,1,2-Trichloroethane	ND		2000	460	ug/L			10/01/25 02:05	2000
1,1-Dichloroethane	ND		2000	760	ug/L			10/01/25 02:05	2000
1,1-Dichloroethene	ND		2000	580	ug/L			10/01/25 02:05	2000
1,2,4-Trichlorobenzene	ND		2000	820	ug/L			10/01/25 02:05	2000
1,2-Dibromo-3-Chloropropane	ND		2000	780	ug/L			10/01/25 02:05	2000
1,2-Dibromoethane	ND		2000	1500	ug/L			10/01/25 02:05	2000
1,2-Dichlorobenzene	ND		2000	1600	ug/L			10/01/25 02:05	2000
1,2-Dichloroethane	ND		2000	420	ug/L			10/01/25 02:05	2000
1,2-Dichloropropane	ND		2000	1400	ug/L			10/01/25 02:05	2000
1,3-Dichlorobenzene	ND		2000	1600	ug/L			10/01/25 02:05	2000
1,4-Dichlorobenzene	ND		2000	1700	ug/L			10/01/25 02:05	2000
2-Butanone (MEK)	ND		20000	2600	ug/L			10/01/25 02:05	2000
2-Hexanone	ND		10000	2500	ug/L			10/01/25 02:05	2000
4-Methyl-2-pentanone (MIBK)	ND		10000	4200	ug/L			10/01/25 02:05	2000
Acetone	ND		20000	6000	ug/L			10/01/25 02:05	2000
Benzene	ND		2000	820	ug/L			10/01/25 02:05	2000
Bromodichloromethane	ND		2000	780	ug/L			10/01/25 02:05	2000
Bromoform	ND		2000	520	ug/L			10/01/25 02:05	2000
Bromomethane	ND		2000	1400	ug/L			10/01/25 02:05	2000
Carbon disulfide	ND		2000	380	ug/L			10/01/25 02:05	2000
Carbon tetrachloride	ND		2000	540	ug/L			10/01/25 02:05	2000
Chlorobenzene	ND		2000	1500	ug/L			10/01/25 02:05	2000
Chloroethane	ND		2000	640	ug/L			10/01/25 02:05	2000
Chloroform	ND		2000	680	ug/L			10/01/25 02:05	2000
Chloromethane	ND		2000	700	ug/L			10/01/25 02:05	2000
cis-1,2-Dichloroethene	**160000**		2000	1600	ug/L			10/01/25 02:05	2000
cis-1,3-Dichloropropene	ND		2000	720	ug/L			10/01/25 02:05	2000
Cyclohexane	ND		2000	360	ug/L			10/01/25 02:05	2000
Dibromochloromethane	ND		2000	640	ug/L			10/01/25 02:05	2000
Dichlorodifluoromethane	ND		2000	1400	ug/L			10/01/25 02:05	2000
Ethylbenzene	ND		2000	1500	ug/L			10/01/25 02:05	2000
Isopropylbenzene	ND		2000	1600	ug/L			10/01/25 02:05	2000
Methyl acetate	ND		5000	2600	ug/L			10/01/25 02:05	2000
Methyl tert-butyl ether	ND		2000	320	ug/L			10/01/25 02:05	2000
Methylcyclohexane	ND		2000	320	ug/L			10/01/25 02:05	2000
Methylene Chloride	ND		2000	880	ug/L			10/01/25 02:05	2000
Styrene	ND		2000	1500	ug/L			10/01/25 02:05	2000
Tetrachloroethene	ND		2000	720	ug/L			10/01/25 02:05	2000
Toluene	ND		2000	1000	ug/L			10/01/25 02:05	2000
trans-1,2-Dichloroethene	ND		2000	1800	ug/L			10/01/25 02:05	2000
trans-1,3-Dichloropropene	ND		2000	740	ug/L			10/01/25 02:05	2000
Trichloroethene	ND		2000	920	ug/L			10/01/25 02:05	2000
Trichlorofluoromethane	ND		2000	1800	ug/L			10/01/25 02:05	2000
Vinyl chloride	**120000**		2000	1800	ug/L			10/01/25 02:05	2000
Xylenes, Total	ND		4000	1300	ug/L			10/01/25 02:05	2000

Client Sample ID: MW3S

Date Collected: 09/24/25 14:15
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-3

Matrix: Water

Surrogate	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4 (Surr)	*109*		*77 - 120*		*10/01/25 02:05*	*2000*
4-Bromofluorobenzene (Surr)	*105*		*73 - 120*		*10/01/25 02:05*	*2000*
Dibromofluoromethane (Surr)	*107*		*75 - 123*		*10/01/25 02:05*	*2000*
Toluene-d8 (Surr)	*101*		*80 - 120*		*10/01/25 02:05*	*2000*

Method: RSK-175 - Dissolved Gases (GC)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	**6800**		440	110	ug/L			10/01/25 11:25	110
Ethane	**1100**		830	170	ug/L			10/01/25 11:25	110
Ethene	**29000**		770	170	ug/L			10/01/25 11:25	110

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)	ND		3.3	0.83	ng/L		10/08/25 08:36	10/09/25 03:28	1
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)	ND		3.3	0.83	ng/L		10/08/25 08:36	10/09/25 03:28	1
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	ND		3.3	0.83	ng/L		10/08/25 08:36	10/09/25 03:28	1
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	ND		8.3	2.1	ng/L		10/08/25 08:36	10/09/25 03:28	1
3-Perfluoropentylpropanoic acid (5:3 FTCA)	ND		8.3	2.3	ng/L		10/08/25 08:36	10/09/25 03:28	1
3-Perfluoropropylpropanoic acid (3:3 FTCA)	ND		3.3	0.83	ng/L		10/08/25 08:36	10/09/25 03:28	1
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	ND		1.7	1.2	ng/L		10/08/25 08:36	10/09/25 03:28	1
N-ethylperfluorooctane sulfonamide (NEtFOSA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
N-methylperfluorooctane sulfonamide (NMeFOSA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluoro-3-methoxypropanoic acid (PFMPA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluoro-4-methoxybutanoic acid (PFMBA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluorobutanesulfonic acid (PFBS)	**4.5**		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluorobutanoic acid (PFBA)	**5.1**		3.3	0.91	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluorodecanesulfonic acid (PFDS)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluorodecanoic acid (PFDA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1

Client Sample ID: MW3S

Date Collected: 09/24/25 14:15
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-3

Matrix: Water

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorododecanesulfonic acid (PFDoS)	ND		1.7	0.50	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluorododecanoic acid (PFDoA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluoroheptanesulfonic acid (PFHpS)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluoroheptanoic acid (PFHpA)	**5.7**		1.7	0.66	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluorohexanesulfonic acid (PFHxS)	**3.4**		1.7	0.66	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluorohexanoic acid (PFHxA)	**12**		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluorononanesulfonic acid (PFNS)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluorononanoic acid (PFNA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluorooctanesulfonamide (PFOSA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluorooctanesulfonic acid (PFOS)	**6.6**	**I**	1.7	1.1	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluorooctanoic acid (PFOA)	**15**		1.7	0.74	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluoropentanesulfonic acid (PFPeS)	**1.7**		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluoropentanoic acid (PFPeA)	**7.3**		1.7	0.50	ng/L		10/08/25 08:36	10/09/25 03:28	1
Perfluoroundecanoic acid (PFUnA)	ND		1.7	0.41	ng/L		10/08/25 08:36	10/09/25 03:28	1

Isotope Dilution	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
13C2 4:2 FTS	*228*	**5+*	*40 - 200*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C2 6:2 FTS	*113*		*40 - 200*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C2 8:2 FTS	*92.4*		*40 - 300*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C2 PFDoA	*22.2*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C3 HFPO-DA	*84.7*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C3 PFBS	*108*		*40 - 135*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C3 PFHxS	*92.9*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C4 PFBA	*76.7*		*5 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C4 PFHpA	*94.9*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C5 PFHxA	*97.3*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C5 PFPeA	*65.6*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C6 PFDA	*84.8*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C7 PFUnA	*64.1*		*30 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C8 FOSA	*79.3*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C8 PFOA	*96.2*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C8 PFOS	*90.1*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
13C9 PFNA	*97.7*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
d3-NMeFOSAA	*58.0*		*40 - 170*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
d3-NMePFOSA	*59.0*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
d5-NEtFOSAA	*33.1*		*25 - 135*	*10/08/25 08:36*	*10/09/25 03:28*	*1*
d5-NEtPFOSA	*26.5*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:28*	*1*

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A - RE

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	ND		8.2	3.3	ng/L		10/01/25 08:35	10/02/25 03:58	1
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	ND		8.2	2.1	ng/L		10/01/25 08:35	10/02/25 03:58	1
Perfluorotetradecanoic acid (PFTeDA)	ND		1.6	0.41	ng/L		10/01/25 08:35	10/02/25 03:58	1
Perfluorotridecanoic acid (PFTrDA)	ND		1.6	0.41	ng/L		10/01/25 08:35	10/02/25 03:58	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW3S

Date Collected: 09/24/25 14:15
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-3

Matrix: Water

Isotope Dilution	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
13C2 PFDoA	*89.1*		*10 - 130*	*10/01/25 08:35*	*10/02/25 03:58*	*1*
13C2 PFTeDA	*80.9*		*10 - 130*	*10/01/25 08:35*	*10/02/25 03:58*	*1*
d7-N-MeFOSE-M	*85.0*		*10 - 130*	*10/01/25 08:35*	*10/02/25 03:58*	*1*
d9-N-EtFOSE-M	*56.5*		*10 - 130*	*10/01/25 08:35*	*10/02/25 03:58*	*1*

Method: SW846 6010D - Metals (ICP)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron	91.6		0.050	0.019	mg/L		09/29/25 08:35	09/29/25 15:45	1

Method: SW846 6010D - Metals (ICP) - Dissolved

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron, Dissolved	88.0		0.050	0.019	mg/L		09/29/25 08:34	09/30/25 00:17	1

General Chemistry

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Sulfate (EPA 300.0)	25.0		20.0	3.5	mg/L			09/30/25 02:11	10

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW3I
Date Collected: 09/24/25 11:50
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-4
Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		1.0	0.82	ug/L			10/01/25 16:02	1
1,1,2,2-Tetrachloroethane	ND		1.0	0.21	ug/L			10/01/25 16:02	1
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1.0	0.31	ug/L			10/01/25 16:02	1
1,1,2-Trichloroethane	ND		1.0	0.23	ug/L			10/01/25 16:02	1
1,1-Dichloroethane	ND		1.0	0.38	ug/L			10/01/25 16:02	1
1,1-Dichloroethene	ND		1.0	0.29	ug/L			10/01/25 16:02	1
1,2,4-Trichlorobenzene	ND		1.0	0.41	ug/L			10/01/25 16:02	1
1,2-Dibromo-3-Chloropropane	ND		1.0	0.39	ug/L			10/01/25 16:02	1
1,2-Dibromoethane	ND		1.0	0.73	ug/L			10/01/25 16:02	1
1,2-Dichlorobenzene	ND		1.0	0.79	ug/L			10/01/25 16:02	1
1,2-Dichloroethane	ND		1.0	0.21	ug/L			10/01/25 16:02	1
1,2-Dichloropropane	ND		1.0	0.72	ug/L			10/01/25 16:02	1
1,3-Dichlorobenzene	ND		1.0	0.78	ug/L			10/01/25 16:02	1
1,4-Dichlorobenzene	ND		1.0	0.84	ug/L			10/01/25 16:02	1
2-Butanone (MEK)	ND		10	1.3	ug/L			10/01/25 16:02	1
2-Hexanone	ND		5.0	1.2	ug/L			10/01/25 16:02	1
4-Methyl-2-pentanone (MIBK)	ND		5.0	2.1	ug/L			10/01/25 16:02	1
Acetone	ND		10	3.0	ug/L			10/01/25 16:02	1
Benzene	ND		1.0	0.41	ug/L			10/01/25 16:02	1
Bromodichloromethane	ND		1.0	0.39	ug/L			10/01/25 16:02	1
Bromoform	ND		1.0	0.26	ug/L			10/01/25 16:02	1
Bromomethane	ND		1.0	0.69	ug/L			10/01/25 16:02	1
Carbon disulfide	ND		1.0	0.19	ug/L			10/01/25 16:02	1
Carbon tetrachloride	ND		1.0	0.27	ug/L			10/01/25 16:02	1
Chlorobenzene	ND		1.0	0.75	ug/L			10/01/25 16:02	1
Chloroethane	ND		1.0	0.32	ug/L			10/01/25 16:02	1
Chloroform	ND		1.0	0.34	ug/L			10/01/25 16:02	1
Chloromethane	ND		1.0	0.35	ug/L			10/01/25 16:02	1
cis-1,2-Dichloroethene	**2.9**		1.0	0.81	ug/L			10/01/25 16:02	1
cis-1,3-Dichloropropene	ND		1.0	0.36	ug/L			10/01/25 16:02	1
Cyclohexane	ND		1.0	0.18	ug/L			10/01/25 16:02	1
Dibromochloromethane	ND		1.0	0.32	ug/L			10/01/25 16:02	1
Dichlorodifluoromethane	ND		1.0	0.68	ug/L			10/01/25 16:02	1
Ethylbenzene	ND		1.0	0.74	ug/L			10/01/25 16:02	1
Isopropylbenzene	ND		1.0	0.79	ug/L			10/01/25 16:02	1
Methyl acetate	ND		2.5	1.3	ug/L			10/01/25 16:02	1
Methyl tert-butyl ether	ND		1.0	0.16	ug/L			10/01/25 16:02	1
Methylcyclohexane	ND		1.0	0.16	ug/L			10/01/25 16:02	1
Methylene Chloride	ND		1.0	0.44	ug/L			10/01/25 16:02	1
Styrene	ND		1.0	0.73	ug/L			10/01/25 16:02	1
Tetrachloroethene	ND		1.0	0.36	ug/L			10/01/25 16:02	1
Toluene	ND		1.0	0.51	ug/L			10/01/25 16:02	1
trans-1,2-Dichloroethene	ND		1.0	0.90	ug/L			10/01/25 16:02	1
trans-1,3-Dichloropropene	ND		1.0	0.37	ug/L			10/01/25 16:02	1
Trichloroethene	**0.53**	**J**	1.0	0.46	ug/L			10/01/25 16:02	1
Trichlorofluoromethane	ND		1.0	0.88	ug/L			10/01/25 16:02	1
Vinyl chloride	**8.4**		1.0	0.90	ug/L			10/01/25 16:02	1
Xylenes, Total	ND		2.0	0.66	ug/L			10/01/25 16:02	1

Client Sample ID: MW3I

Date Collected: 09/24/25 11:50
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-4

Matrix: Water

Surrogate	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4 (Surr)	*105*		*77 - 120*		*10/01/25 16:02*	*1*
4-Bromofluorobenzene (Surr)	*104*		*73 - 120*		*10/01/25 16:02*	*1*
Dibromofluoromethane (Surr)	*104*		*75 - 123*		*10/01/25 16:02*	*1*
Toluene-d8 (Surr)	*100*		*80 - 120*		*10/01/25 16:02*	*1*

Method: RSK-175 - Dissolved Gases (GC)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	**140**		4.0	1.0	ug/L			10/01/25 11:44	1
Ethane	ND		7.5	1.5	ug/L			10/01/25 11:44	1
Ethene	ND		7.0	1.5	ug/L			10/01/25 11:44	1

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)	ND		3.1	0.78	ng/L		10/08/25 08:36	10/09/25 03:41	1
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)	ND		3.1	0.78	ng/L		10/08/25 08:36	10/09/25 03:41	1
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	ND		3.1	0.78	ng/L		10/08/25 08:36	10/09/25 03:41	1
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	ND		7.8	1.9	ng/L		10/08/25 08:36	10/09/25 03:41	1
3-Perfluoropentylpropanoic acid (5:3 FTCA)	ND		7.8	2.2	ng/L		10/08/25 08:36	10/09/25 03:41	1
3-Perfluoropropylpropanoic acid (3:3 FTCA)	ND		3.1	0.78	ng/L		10/08/25 08:36	10/09/25 03:41	1
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	ND		1.6	1.1	ng/L		10/08/25 08:36	10/09/25 03:41	1
N-ethylperfluorooctane sulfonamide (NEtFOSA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	ND		7.8	3.1	ng/L		10/08/25 08:36	10/09/25 03:41	1
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
N-methylperfluorooctane sulfonamide (NMeFOSA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	ND		7.8	1.9	ng/L		10/08/25 08:36	10/09/25 03:41	1
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluoro-3-methoxypropanoic acid (PFMPA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluoro-4-methoxybutanoic acid (PFMBA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorobutanesulfonic acid (PFBS)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorobutanoic acid (PFBA)	**0.97**	**J**	3.1	0.85	ng/L		10/08/25 08:36	10/09/25 03:41	1

Client Sample ID: MW3I
Date Collected: 09/24/25 11:50
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-4
Matrix: Water

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorodecanesulfonic acid (PFDS)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorodecanoic acid (PFDA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorododecanesulfonic acid (PFDoS)	ND		1.6	0.47	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorododecanoic acid (PFDoA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluoroheptanesulfonic acid (PFHpS)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluoroheptanoic acid (PFHpA)	ND		1.6	0.62	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorohexanesulfonic acid (PFHxS)	ND		1.6	0.62	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorohexanoic acid (PFHxA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorononanesulfonic acid (PFNS)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorononanoic acid (PFNA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorooctanesulfonamide (PFOSA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorooctanesulfonic acid (PFOS)	ND		1.6	1.0	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorooctanoic acid (PFOA)	ND		1.6	0.70	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluoropentanesulfonic acid (PFPeS)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluoropentanoic acid (PFPeA)	ND		1.6	0.47	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorotetradecanoic acid (PFTeDA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluorotridecanoic acid (PFTrDA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1
Perfluoroundecanoic acid (PFUnA)	ND		1.6	0.39	ng/L		10/08/25 08:36	10/09/25 03:41	1

Isotope Dilution	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
13C2 4:2 FTS	*108*		*40 - 200*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C2 6:2 FTS	*99.2*		*40 - 200*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C2 8:2 FTS	*94.6*		*40 - 300*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C2 PFDoA	*57.2*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C2 PFTeDA	*38.1*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C3 HFPO-DA	*91.0*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C3 PFBS	*116*		*40 - 135*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C3 PFHxS	*96.3*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C4 PFBA	*105*		*5 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C4 PFHpA	*94.7*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C5 PFHxA	*106*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C5 PFPeA	*108*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C6 PFDA	*88.2*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C7 PFUnA	*73.1*		*30 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C8 FOSA	*84.1*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C8 PFOA	*99.1*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C8 PFOS	*92.5*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
13C9 PFNA	*100*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
d3-NMeFOSAA	*68.6*		*40 - 170*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
d3-NMePFOSA	*62.2*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
d5-NEtFOSAA	*54.2*		*25 - 135*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
d5-NEtPFOSA	*52.7*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
d7-N-MeFOSE-M	*23.0*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*
d9-N-EtFOSE-M	*15.4*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:41*	*1*

Method: SW846 6010D - Metals (ICP)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron	1.1		0.050	0.019	mg/L		09/29/25 08:35	09/29/25 15:47	1

Client Sample Results

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW3I

Date Collected: 09/24/25 11:50
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-4

Matrix: Water

Method: SW846 6010D - Metals (ICP) - Dissolved

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron, Dissolved	1.0		0.050	0.019	mg/L		09/29/25 08:34	09/30/25 00:19	1

General Chemistry

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Sulfate (EPA 300.0)	79.9		10.0	1.7	mg/L			09/30/25 02:26	5

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW4S

Date Collected: 09/24/25 15:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-5

Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		2000	1600	ug/L			10/01/25 02:51	2000
1,1,2,2-Tetrachloroethane	ND		2000	420	ug/L			10/01/25 02:51	2000
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		2000	620	ug/L			10/01/25 02:51	2000
1,1,2-Trichloroethane	ND		2000	460	ug/L			10/01/25 02:51	2000
1,1-Dichloroethane	ND		2000	760	ug/L			10/01/25 02:51	2000
1,1-Dichloroethene	ND		2000	580	ug/L			10/01/25 02:51	2000
1,2,4-Trichlorobenzene	ND		2000	820	ug/L			10/01/25 02:51	2000
1,2-Dibromo-3-Chloropropane	ND		2000	780	ug/L			10/01/25 02:51	2000
1,2-Dibromoethane	ND		2000	1500	ug/L			10/01/25 02:51	2000
1,2-Dichlorobenzene	ND		2000	1600	ug/L			10/01/25 02:51	2000
1,2-Dichloroethane	ND		2000	420	ug/L			10/01/25 02:51	2000
1,2-Dichloropropane	ND		2000	1400	ug/L			10/01/25 02:51	2000
1,3-Dichlorobenzene	ND		2000	1600	ug/L			10/01/25 02:51	2000
1,4-Dichlorobenzene	ND		2000	1700	ug/L			10/01/25 02:51	2000
2-Butanone (MEK)	ND		20000	2600	ug/L			10/01/25 02:51	2000
2-Hexanone	ND		10000	2500	ug/L			10/01/25 02:51	2000
4-Methyl-2-pentanone (MIBK)	ND		10000	4200	ug/L			10/01/25 02:51	2000
Acetone	ND		20000	6000	ug/L			10/01/25 02:51	2000
Benzene	ND		2000	820	ug/L			10/01/25 02:51	2000
Bromodichloromethane	ND		2000	780	ug/L			10/01/25 02:51	2000
Bromoform	ND		2000	520	ug/L			10/01/25 02:51	2000
Bromomethane	ND		2000	1400	ug/L			10/01/25 02:51	2000
Carbon disulfide	ND		2000	380	ug/L			10/01/25 02:51	2000
Carbon tetrachloride	ND		2000	540	ug/L			10/01/25 02:51	2000
Chlorobenzene	ND		2000	1500	ug/L			10/01/25 02:51	2000
Chloroethane	ND		2000	640	ug/L			10/01/25 02:51	2000
Chloroform	ND		2000	680	ug/L			10/01/25 02:51	2000
Chloromethane	ND		2000	700	ug/L			10/01/25 02:51	2000
cis-1,2-Dichloroethene	**8100**		2000	1600	ug/L			10/01/25 02:51	2000
cis-1,3-Dichloropropene	ND		2000	720	ug/L			10/01/25 02:51	2000
Cyclohexane	ND		2000	360	ug/L			10/01/25 02:51	2000
Dibromochloromethane	ND		2000	640	ug/L			10/01/25 02:51	2000
Dichlorodifluoromethane	ND		2000	1400	ug/L			10/01/25 02:51	2000
Ethylbenzene	ND		2000	1500	ug/L			10/01/25 02:51	2000
Isopropylbenzene	ND		2000	1600	ug/L			10/01/25 02:51	2000
Methyl acetate	ND		5000	2600	ug/L			10/01/25 02:51	2000
Methyl tert-butyl ether	ND		2000	320	ug/L			10/01/25 02:51	2000
Methylcyclohexane	ND		2000	320	ug/L			10/01/25 02:51	2000
Methylene Chloride	ND		2000	880	ug/L			10/01/25 02:51	2000
Styrene	ND		2000	1500	ug/L			10/01/25 02:51	2000
Tetrachloroethene	**85000**		2000	720	ug/L			10/01/25 02:51	2000
Toluene	ND		2000	1000	ug/L			10/01/25 02:51	2000
trans-1,2-Dichloroethene	ND		2000	1800	ug/L			10/01/25 02:51	2000
trans-1,3-Dichloropropene	ND		2000	740	ug/L			10/01/25 02:51	2000
Trichloroethene	**4100**		2000	920	ug/L			10/01/25 02:51	2000
Trichlorofluoromethane	ND		2000	1800	ug/L			10/01/25 02:51	2000
Vinyl chloride	ND		2000	1800	ug/L			10/01/25 02:51	2000
Xylenes, Total	ND		4000	1300	ug/L			10/01/25 02:51	2000

Client Sample ID: MW4S

Date Collected: 09/24/25 15:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-5

Matrix: Water

Surrogate	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4 (Surr)	*109*		*77 - 120*		*10/01/25 02:51*	*2000*
4-Bromofluorobenzene (Surr)	*108*		*73 - 120*		*10/01/25 02:51*	*2000*
Dibromofluoromethane (Surr)	*108*		*75 - 123*		*10/01/25 02:51*	*2000*
Toluene-d8 (Surr)	*100*		*80 - 120*		*10/01/25 02:51*	*2000*

Method: RSK-175 - Dissolved Gases (GC)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Ethane	**390**		7.5	1.5	ug/L			10/01/25 12:03	1
Ethene	**13**		7.0	1.5	ug/L			10/01/25 12:03	1

Method: RSK-175 - Dissolved Gases (GC) - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	**4600**		180	44	ug/L			10/01/25 13:18	44

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)	ND		3.1	0.77	ng/L		10/08/25 08:36	10/09/25 03:55	1
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)	ND		3.1	0.77	ng/L		10/08/25 08:36	10/09/25 03:55	1
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	ND		3.1	0.77	ng/L		10/08/25 08:36	10/09/25 03:55	1
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	ND		7.7	1.9	ng/L		10/08/25 08:36	10/09/25 03:55	1
3-Perfluoropentylpropanoic acid (5:3 FTCA)	ND		7.7	2.2	ng/L		10/08/25 08:36	10/09/25 03:55	1
3-Perfluoropropylpropanoic acid (3:3 FTCA)	ND		3.1	0.77	ng/L		10/08/25 08:36	10/09/25 03:55	1
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	ND		1.5	1.1	ng/L		10/08/25 08:36	10/09/25 03:55	1
N-ethylperfluorooctane sulfonamide (NEtFOSA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	ND		7.7	3.1	ng/L		10/08/25 08:36	10/09/25 03:55	1
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
N-methylperfluorooctane sulfonamide (NMeFOSA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	ND		7.7	1.9	ng/L		10/08/25 08:36	10/09/25 03:55	1
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluoro-3-methoxypropanoic acid (PFMPA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluoro-4-methoxybutanoic acid (PFMBA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1

Client Sample ID: MW4S
Date Collected: 09/24/25 15:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-5
Matrix: Water

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorobutanesulfonic acid (PFBS)	**4.4**		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorobutanoic acid (PFBA)	**12**		3.1	0.85	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorodecanesulfonic acid (PFDS)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorodecanoic acid (PFDA)	**1.5**		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorododecanesulfonic acid (PFDoS)	ND		1.5	0.46	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorododecanoic acid (PFDoA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluoroheptanesulfonic acid (PFHpS)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluoroheptanoic acid (PFHpA)	**11**		1.5	0.62	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorohexanesulfonic acid (PFHxS)	**3.7**		1.5	0.62	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorohexanoic acid (PFHxA)	**18**		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorononanesulfonic acid (PFNS)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorononanoic acid (PFNA)	**4.1**		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorooctanesulfonamide (PFOSA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorooctanesulfonic acid (PFOS)	**17**		1.5	1.0	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorooctanoic acid (PFOA)	**24**		1.5	0.69	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluoropentanesulfonic acid (PFPeS)	**0.80**	**J**	1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluoropentanoic acid (PFPeA)	**25**		1.5	0.46	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorotetradecanoic acid (PFTeDA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluorotridecanoic acid (PFTrDA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1
Perfluoroundecanoic acid (PFUnA)	ND		1.5	0.39	ng/L		10/08/25 08:36	10/09/25 03:55	1

Isotope Dilution	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
13C2 4:2 FTS	*191*		*40 - 200*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C2 6:2 FTS	*139*		*40 - 200*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C2 8:2 FTS	*143*		*40 - 300*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C2 PFDoA	*81.8*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C2 PFTeDA	*61.9*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C3 HFPO-DA	*82.9*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C3 PFBS	*117*		*40 - 135*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C3 PFHxS	*99.9*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C4 PFBA	*109*		*5 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C4 PFHpA	*91.1*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C5 PFHxA	*97.5*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C5 PFPeA	*99.2*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C6 PFDA	*96.6*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C7 PFUnA	*85.7*		*30 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C8 FOSA	*94.4*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C8 PFOA	*102*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C8 PFOS	*104*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
13C9 PFNA	*100*		*40 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
d3-NMeFOSAA	*112*		*40 - 170*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
d3-NMePFOSA	*81.1*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
d5-NEtFOSAA	*87.5*		*25 - 135*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
d5-NEtPFOSA	*71.7*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
d7-N-MeFOSE-M	*57.3*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*
d9-N-EtFOSE-M	*44.0*		*10 - 130*	*10/08/25 08:36*	*10/09/25 03:55*	*1*

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW4S

Date Collected: 09/24/25 15:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-5

Matrix: Water

Method: SW846 6010D - Metals (ICP)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron	1.5		0.050	0.019	mg/L		09/29/25 08:35	09/29/25 15:49	1

Method: SW846 6010D - Metals (ICP) - Dissolved

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron, Dissolved	0.081		0.050	0.019	mg/L		09/29/25 08:34	09/30/25 00:24	1

General Chemistry

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Sulfate (EPA 300.0)	131		40.0	7.0	mg/L			09/30/25 16:37	20

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW4I
Date Collected: 09/24/25 14:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-6
Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		40	33	ug/L			10/01/25 03:14	40
1,1,2,2-Tetrachloroethane	ND		40	8.4	ug/L			10/01/25 03:14	40
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		40	12	ug/L			10/01/25 03:14	40
1,1,2-Trichloroethane	ND		40	9.2	ug/L			10/01/25 03:14	40
1,1-Dichloroethane	ND		40	15	ug/L			10/01/25 03:14	40
1,1-Dichloroethene	**36**	**J**	40	12	ug/L			10/01/25 03:14	40
1,2,4-Trichlorobenzene	ND		40	16	ug/L			10/01/25 03:14	40
1,2-Dibromo-3-Chloropropane	ND		40	16	ug/L			10/01/25 03:14	40
1,2-Dibromoethane	ND		40	29	ug/L			10/01/25 03:14	40
1,2-Dichlorobenzene	ND		40	32	ug/L			10/01/25 03:14	40
1,2-Dichloroethane	ND		40	8.4	ug/L			10/01/25 03:14	40
1,2-Dichloropropane	ND		40	29	ug/L			10/01/25 03:14	40
1,3-Dichlorobenzene	ND		40	31	ug/L			10/01/25 03:14	40
1,4-Dichlorobenzene	ND		40	34	ug/L			10/01/25 03:14	40
2-Butanone (MEK)	ND		400	53	ug/L			10/01/25 03:14	40
2-Hexanone	ND		200	50	ug/L			10/01/25 03:14	40
4-Methyl-2-pentanone (MIBK)	ND		200	84	ug/L			10/01/25 03:14	40
Acetone	ND		400	120	ug/L			10/01/25 03:14	40
Benzene	ND		40	16	ug/L			10/01/25 03:14	40
Bromodichloromethane	ND		40	16	ug/L			10/01/25 03:14	40
Bromoform	ND		40	10	ug/L			10/01/25 03:14	40
Bromomethane	ND		40	28	ug/L			10/01/25 03:14	40
Carbon disulfide	ND		40	7.6	ug/L			10/01/25 03:14	40
Carbon tetrachloride	ND		40	11	ug/L			10/01/25 03:14	40
Chlorobenzene	ND		40	30	ug/L			10/01/25 03:14	40
Chloroethane	ND		40	13	ug/L			10/01/25 03:14	40
Chloroform	ND		40	14	ug/L			10/01/25 03:14	40
Chloromethane	ND		40	14	ug/L			10/01/25 03:14	40
cis-1,2-Dichloroethene	**3800**	**F1**	40	32	ug/L			10/01/25 03:14	40
cis-1,3-Dichloropropene	ND		40	14	ug/L			10/01/25 03:14	40
Cyclohexane	ND		40	7.2	ug/L			10/01/25 03:14	40
Dibromochloromethane	ND		40	13	ug/L			10/01/25 03:14	40
Dichlorodifluoromethane	ND		40	27	ug/L			10/01/25 03:14	40
Ethylbenzene	ND		40	30	ug/L			10/01/25 03:14	40
Isopropylbenzene	ND		40	32	ug/L			10/01/25 03:14	40
Methyl acetate	ND		100	52	ug/L			10/01/25 03:14	40
Methyl tert-butyl ether	ND		40	6.4	ug/L			10/01/25 03:14	40
Methylcyclohexane	ND		40	6.4	ug/L			10/01/25 03:14	40
Methylene Chloride	ND		40	18	ug/L			10/01/25 03:14	40
Styrene	ND		40	29	ug/L			10/01/25 03:14	40
Tetrachloroethene	**20**	**J**	40	14	ug/L			10/01/25 03:14	40
Toluene	ND		40	20	ug/L			10/01/25 03:14	40
trans-1,2-Dichloroethene	ND		40	36	ug/L			10/01/25 03:14	40
trans-1,3-Dichloropropene	ND		40	15	ug/L			10/01/25 03:14	40
Trichloroethene	ND		40	18	ug/L			10/01/25 03:14	40
Trichlorofluoromethane	ND		40	35	ug/L			10/01/25 03:14	40
Vinyl chloride	**550**		40	36	ug/L			10/01/25 03:14	40
Xylenes, Total	ND		80	26	ug/L			10/01/25 03:14	40

Client Sample ID: MW4I

Date Collected: 09/24/25 14:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-6
Matrix: Water

Surrogate	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4 (Surr)	*119*		*77 - 120*		*10/01/25 03:14*	*40*
4-Bromofluorobenzene (Surr)	*114*		*73 - 120*		*10/01/25 03:14*	*40*
Dibromofluoromethane (Surr)	*117*		*75 - 123*		*10/01/25 03:14*	*40*
Toluene-d8 (Surr)	*108*		*80 - 120*		*10/01/25 03:14*	*40*

Method: RSK-175 - Dissolved Gases (GC)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Ethane	**15**		7.5	1.5	ug/L			10/01/25 12:21	1
Ethene	**54**		7.0	1.5	ug/L			10/01/25 12:21	1

Method: RSK-175 - Dissolved Gases (GC) - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	**1400**		88	22	ug/L			10/01/25 13:56	22

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)	ND		3.0	0.76	ng/L		10/08/25 08:36	10/09/25 04:08	1
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)	ND		3.0	0.76	ng/L		10/08/25 08:36	10/09/25 04:08	1
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	ND		3.0	0.76	ng/L		10/08/25 08:36	10/09/25 04:08	1
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	ND		7.6	1.9	ng/L		10/08/25 08:36	10/09/25 04:08	1
3-Perfluoropentylpropanoic acid (5:3 FTCA)	ND		7.6	2.1	ng/L		10/08/25 08:36	10/09/25 04:08	1
3-Perfluoropropylpropanoic acid (3:3 FTCA)	ND		3.0	0.76	ng/L		10/08/25 08:36	10/09/25 04:08	1
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	ND		1.5	1.1	ng/L		10/08/25 08:36	10/09/25 04:08	1
N-ethylperfluorooctane sulfonamide (NEtFOSA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	ND		7.6	3.0	ng/L		10/08/25 08:36	10/09/25 04:08	1
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
N-methylperfluorooctane sulfonamide (NMeFOSA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	ND		7.6	1.9	ng/L		10/08/25 08:36	10/09/25 04:08	1
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluoro-3-methoxypropanoic acid (PFMPA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluoro-4-methoxybutanoic acid (PFMBA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1

Client Sample ID: MW4I
Date Collected: 09/24/25 14:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-6
Matrix: Water

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorobutanesulfonic acid (PFBS)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorobutanoic acid (PFBA)	**1.1**	**J**	3.0	0.83	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorodecanesulfonic acid (PFDS)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorodecanoic acid (PFDA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorododecanesulfonic acid (PFDoS)	ND		1.5	0.45	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorododecanoic acid (PFDoA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluoroheptanesulfonic acid (PFHpS)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluoroheptanoic acid (PFHpA)	ND		1.5	0.61	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorohexanesulfonic acid (PFHxS)	ND		1.5	0.61	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorohexanoic acid (PFHxA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorononanesulfonic acid (PFNS)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorononanoic acid (PFNA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorooctanesulfonamide (PFOSA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorooctanesulfonic acid (PFOS)	ND		1.5	0.99	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorooctanoic acid (PFOA)	ND		1.5	0.68	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluoropentanesulfonic acid (PFPeS)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluoropentanoic acid (PFPeA)	ND		1.5	0.45	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorotetradecanoic acid (PFTeDA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluorotridecanoic acid (PFTrDA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1
Perfluoroundecanoic acid (PFUnA)	ND		1.5	0.38	ng/L		10/08/25 08:36	10/09/25 04:08	1

Isotope Dilution	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
13C2 4:2 FTS	*113*		*40 - 200*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C2 6:2 FTS	*92.9*		*40 - 200*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C2 8:2 FTS	*97.7*		*40 - 300*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C2 PFDoA	*80.5*		*10 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C2 PFTeDA	*66.6*		*10 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C3 HFPO-DA	*85.9*		*40 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C3 PFBS	*110*		*40 - 135*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C3 PFHxS	*92.8*		*40 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C4 PFBA	*99.7*		*5 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C4 PFHpA	*86.4*		*40 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C5 PFHxA	*99.5*		*40 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C5 PFPeA	*102*		*40 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C6 PFDA	*84.0*		*40 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C7 PFUnA	*78.0*		*30 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C8 FOSA	*80.0*		*40 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C8 PFOA	*88.8*		*40 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C8 PFOS	*88.8*		*40 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
13C9 PFNA	*93.4*		*40 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
d3-NMeFOSAA	*77.1*		*40 - 170*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
d3-NMePFOSA	*69.3*		*10 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
d5-NEtFOSAA	*66.6*		*25 - 135*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
d5-NEtPFOSA	*65.7*		*10 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
d7-N-MeFOSE-M	*49.2*		*10 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*
d9-N-EtFOSE-M	*38.6*		*10 - 130*	*10/08/25 08:36*	*10/09/25 04:08*	*1*

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW4I

Date Collected: 09/24/25 14:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-6

Matrix: Water

Method: SW846 6010D - Metals (ICP)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron	1.2		0.050	0.019	mg/L		09/29/25 08:35	09/29/25 15:52	1

Method: SW846 6010D - Metals (ICP) - Dissolved

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron, Dissolved	1.3		0.050	0.019	mg/L		09/29/25 08:34	09/30/25 00:22	1

General Chemistry

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Sulfate (EPA 300.0)	128		10.0	1.7	mg/L			09/30/25 16:52	5

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW5S

Date Collected: 09/24/25 09:50
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-7

Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		10	8.2	ug/L			10/01/25 03:38	10
1,1,2,2-Tetrachloroethane	ND		10	2.1	ug/L			10/01/25 03:38	10
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		10	3.1	ug/L			10/01/25 03:38	10
1,1,2-Trichloroethane	ND		10	2.3	ug/L			10/01/25 03:38	10
1,1-Dichloroethane	ND		10	3.8	ug/L			10/01/25 03:38	10
1,1-Dichloroethene	ND		10	2.9	ug/L			10/01/25 03:38	10
1,2,4-Trichlorobenzene	ND		10	4.1	ug/L			10/01/25 03:38	10
1,2-Dibromo-3-Chloropropane	ND		10	3.9	ug/L			10/01/25 03:38	10
1,2-Dibromoethane	ND		10	7.3	ug/L			10/01/25 03:38	10
1,2-Dichlorobenzene	ND		10	7.9	ug/L			10/01/25 03:38	10
1,2-Dichloroethane	ND		10	2.1	ug/L			10/01/25 03:38	10
1,2-Dichloropropane	ND		10	7.2	ug/L			10/01/25 03:38	10
1,3-Dichlorobenzene	ND		10	7.8	ug/L			10/01/25 03:38	10
1,4-Dichlorobenzene	ND		10	8.4	ug/L			10/01/25 03:38	10
2-Butanone (MEK)	**34**	**J**	100	13	ug/L			10/01/25 03:38	10
2-Hexanone	ND		50	12	ug/L			10/01/25 03:38	10
4-Methyl-2-pentanone (MIBK)	ND		50	21	ug/L			10/01/25 03:38	10
Acetone	**86**	**J**	100	30	ug/L			10/01/25 03:38	10
Benzene	**18**		10	4.1	ug/L			10/01/25 03:38	10
Bromodichloromethane	ND		10	3.9	ug/L			10/01/25 03:38	10
Bromoform	ND		10	2.6	ug/L			10/01/25 03:38	10
Bromomethane	ND		10	6.9	ug/L			10/01/25 03:38	10
Carbon disulfide	ND		10	1.9	ug/L			10/01/25 03:38	10
Carbon tetrachloride	ND		10	2.7	ug/L			10/01/25 03:38	10
Chlorobenzene	ND		10	7.5	ug/L			10/01/25 03:38	10
Chloroethane	ND		10	3.2	ug/L			10/01/25 03:38	10
Chloroform	ND		10	3.4	ug/L			10/01/25 03:38	10
Chloromethane	ND		10	3.5	ug/L			10/01/25 03:38	10
cis-1,2-Dichloroethene	ND		10	8.1	ug/L			10/01/25 03:38	10
cis-1,3-Dichloropropene	ND		10	3.6	ug/L			10/01/25 03:38	10
Cyclohexane	**370**		10	1.8	ug/L			10/01/25 03:38	10
Dibromochloromethane	ND		10	3.2	ug/L			10/01/25 03:38	10
Dichlorodifluoromethane	ND		10	6.8	ug/L			10/01/25 03:38	10
Ethylbenzene	**260**		10	7.4	ug/L			10/01/25 03:38	10
Isopropylbenzene	**110**		10	7.9	ug/L			10/01/25 03:38	10
Methyl acetate	ND		25	13	ug/L			10/01/25 03:38	10
Methyl tert-butyl ether	ND		10	1.6	ug/L			10/01/25 03:38	10
Methylcyclohexane	**150**		10	1.6	ug/L			10/01/25 03:38	10
Methylene Chloride	ND		10	4.4	ug/L			10/01/25 03:38	10
Styrene	ND		10	7.3	ug/L			10/01/25 03:38	10
Tetrachloroethene	ND		10	3.6	ug/L			10/01/25 03:38	10
Toluene	**6.4**	**J**	10	5.1	ug/L			10/01/25 03:38	10
trans-1,2-Dichloroethene	ND		10	9.0	ug/L			10/01/25 03:38	10
trans-1,3-Dichloropropene	ND		10	3.7	ug/L			10/01/25 03:38	10
Trichloroethene	ND		10	4.6	ug/L			10/01/25 03:38	10
Trichlorofluoromethane	ND		10	8.8	ug/L			10/01/25 03:38	10
Vinyl chloride	ND		10	9.0	ug/L			10/01/25 03:38	10
Xylenes, Total	**23**		20	6.6	ug/L			10/01/25 03:38	10

Client Sample ID: MW5S
Date Collected: 09/24/25 09:50
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-7
Matrix: Water

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	*124*	*S1+*	*77 - 120*		*10/01/25 03:38*	*10*
4-Bromofluorobenzene (Surr)	*108*		*73 - 120*		*10/01/25 03:38*	*10*
Dibromofluoromethane (Surr)	*114*		*75 - 123*		*10/01/25 03:38*	*10*
Toluene-d8 (Surr)	*102*		*80 - 120*		*10/01/25 03:38*	*10*

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)	ND		3.5	0.88	ng/L		10/03/25 13:46	10/06/25 06:17	1
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)	ND		3.5	0.88	ng/L		10/03/25 13:46	10/06/25 06:17	1
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	ND		3.5	0.88	ng/L		10/03/25 13:46	10/06/25 06:17	1
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	ND		8.8	2.2	ng/L		10/03/25 13:46	10/06/25 06:17	1
3-Perfluoropentylpropanoic acid (5:3 FTCA)	ND		8.8	2.5	ng/L		10/03/25 13:46	10/06/25 06:17	1
3-Perfluoropropylpropanoic acid (3:3 FTCA)	ND		3.5	0.88	ng/L		10/03/25 13:46	10/06/25 06:17	1
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	ND		1.8	1.2	ng/L		10/03/25 13:46	10/06/25 06:17	1
N-ethylperfluorooctane sulfonamide (NEtFOSA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	ND		8.8	3.5	ng/L		10/03/25 13:46	10/06/25 06:17	1
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
N-methylperfluorooctane sulfonamide (NMeFOSA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	ND		8.8	2.2	ng/L		10/03/25 13:46	10/06/25 06:17	1
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluoro-3-methoxypropanoic acid (PFMPA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluoro-4-methoxybutanoic acid (PFMBA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorobutanesulfonic acid (PFBS)	**0.85**	**J**	1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorobutanoic acid (PFBA)	**9.1**		3.5	0.97	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorodecanesulfonic acid (PFDS)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorodecanoic acid (PFDA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorododecanesulfonic acid (PFDoS)	ND		1.8	0.53	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorododecanoic acid (PFDoA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1

Client Sample ID: MW5S
Date Collected: 09/24/25 09:50
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-7
Matrix: Water

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanesulfonic acid (PFHpS)	ND		1.8	1.8	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluoroheptanoic acid (PFHpA)	**2.6**		1.8	0.70	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorohexanesulfonic acid (PFHxS)	**1.4**	**J**	1.8	0.70	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorohexanoic acid (PFHxA)	**5.0**		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorononanesulfonic acid (PFNS)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorononanoic acid (PFNA)	ND		1.8	1.8	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorooctanesulfonamide (PFOSA)	**0.48**	**J**	1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorooctanesulfonic acid (PFOS)	**20**		1.8	1.1	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorooctanoic acid (PFOA)	**14**		1.8	0.79	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluoropentanesulfonic acid (PFPeS)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluoropentanoic acid (PFPeA)	**8.1**		1.8	0.53	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorotetradecanoic acid (PFTeDA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluorotridecanoic acid (PFTrDA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1
Perfluoroundecanoic acid (PFUnA)	ND		1.8	0.44	ng/L		10/03/25 13:46	10/06/25 06:17	1

Isotope Dilution	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
13C2 4:2 FTS	*219*	**5+*	*40 - 200*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C2 6:2 FTS	*190*		*40 - 200*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C2 8:2 FTS	*125*		*40 - 300*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C2 PFDoA	*80.5*		*10 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C2 PFTeDA	*66.6*		*10 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C3 HFPO-DA	*95.1*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C3 PFBS	*103*		*40 - 135*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C3 PFHxS	*96.3*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C4 PFBA	*103*		*5 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C4 PFHpA	*100*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C5 PFHxA	*97.9*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C5 PFPeA	*86.5*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C6 PFDA	*95.1*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C7 PFUnA	*83.1*		*30 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C8 FOSA	*93.6*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C8 PFOA	*98.2*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C8 PFOS	*94.5*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
13C9 PFNA	*99.3*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
d3-NMeFOSAA	*89.9*		*40 - 170*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
d3-NMePFOSA	*73.0*		*10 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
d5-NEtFOSAA	*80.5*		*25 - 135*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
d5-NEtPFOSA	*63.1*		*10 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
d7-N-MeFOSE-M	*51.8*		*10 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*
d9-N-EtFOSE-M	*43.6*		*10 - 130*	*10/03/25 13:46*	*10/06/25 06:17*	*1*

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW5I
Date Collected: 09/24/25 10:40
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-8
Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		1.0	0.82	ug/L			10/01/25 04:01	1
1,1,2,2-Tetrachloroethane	ND		1.0	0.21	ug/L			10/01/25 04:01	1
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1.0	0.31	ug/L			10/01/25 04:01	1
1,1,2-Trichloroethane	ND		1.0	0.23	ug/L			10/01/25 04:01	1
1,1-Dichloroethane	ND		1.0	0.38	ug/L			10/01/25 04:01	1
1,1-Dichloroethene	ND		1.0	0.29	ug/L			10/01/25 04:01	1
1,2,4-Trichlorobenzene	ND		1.0	0.41	ug/L			10/01/25 04:01	1
1,2-Dibromo-3-Chloropropane	ND		1.0	0.39	ug/L			10/01/25 04:01	1
1,2-Dibromoethane	ND		1.0	0.73	ug/L			10/01/25 04:01	1
1,2-Dichlorobenzene	ND		1.0	0.79	ug/L			10/01/25 04:01	1
1,2-Dichloroethane	ND		1.0	0.21	ug/L			10/01/25 04:01	1
1,2-Dichloropropane	ND		1.0	0.72	ug/L			10/01/25 04:01	1
1,3-Dichlorobenzene	ND		1.0	0.78	ug/L			10/01/25 04:01	1
1,4-Dichlorobenzene	ND		1.0	0.84	ug/L			10/01/25 04:01	1
2-Butanone (MEK)	ND		10	1.3	ug/L			10/01/25 04:01	1
2-Hexanone	ND		5.0	1.2	ug/L			10/01/25 04:01	1
4-Methyl-2-pentanone (MIBK)	ND		5.0	2.1	ug/L			10/01/25 04:01	1
Acetone	ND		10	3.0	ug/L			10/01/25 04:01	1
Benzene	ND		1.0	0.41	ug/L			10/01/25 04:01	1
Bromodichloromethane	ND		1.0	0.39	ug/L			10/01/25 04:01	1
Bromoform	ND		1.0	0.26	ug/L			10/01/25 04:01	1
Bromomethane	ND		1.0	0.69	ug/L			10/01/25 04:01	1
Carbon disulfide	ND		1.0	0.19	ug/L			10/01/25 04:01	1
Carbon tetrachloride	ND		1.0	0.27	ug/L			10/01/25 04:01	1
Chlorobenzene	ND		1.0	0.75	ug/L			10/01/25 04:01	1
Chloroethane	ND		1.0	0.32	ug/L			10/01/25 04:01	1
Chloroform	ND		1.0	0.34	ug/L			10/01/25 04:01	1
Chloromethane	ND		1.0	0.35	ug/L			10/01/25 04:01	1
cis-1,2-Dichloroethene	ND		1.0	0.81	ug/L			10/01/25 04:01	1
cis-1,3-Dichloropropene	ND		1.0	0.36	ug/L			10/01/25 04:01	1
Cyclohexane	ND		1.0	0.18	ug/L			10/01/25 04:01	1
Dibromochloromethane	ND		1.0	0.32	ug/L			10/01/25 04:01	1
Dichlorodifluoromethane	ND		1.0	0.68	ug/L			10/01/25 04:01	1
Ethylbenzene	ND		1.0	0.74	ug/L			10/01/25 04:01	1
Isopropylbenzene	ND		1.0	0.79	ug/L			10/01/25 04:01	1
Methyl acetate	ND		2.5	1.3	ug/L			10/01/25 04:01	1
Methyl tert-butyl ether	ND		1.0	0.16	ug/L			10/01/25 04:01	1
Methylcyclohexane	ND		1.0	0.16	ug/L			10/01/25 04:01	1
Methylene Chloride	ND		1.0	0.44	ug/L			10/01/25 04:01	1
Styrene	ND		1.0	0.73	ug/L			10/01/25 04:01	1
Tetrachloroethene	ND		1.0	0.36	ug/L			10/01/25 04:01	1
Toluene	ND		1.0	0.51	ug/L			10/01/25 04:01	1
trans-1,2-Dichloroethene	ND		1.0	0.90	ug/L			10/01/25 04:01	1
trans-1,3-Dichloropropene	ND		1.0	0.37	ug/L			10/01/25 04:01	1
Trichloroethene	ND		1.0	0.46	ug/L			10/01/25 04:01	1
Trichlorofluoromethane	ND		1.0	0.88	ug/L			10/01/25 04:01	1
Vinyl chloride	ND		1.0	0.90	ug/L			10/01/25 04:01	1
Xylenes, Total	ND		2.0	0.66	ug/L			10/01/25 04:01	1

Client Sample ID: MW5I
Date Collected: 09/24/25 10:40
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-8
Matrix: Water

Surrogate	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4 (Surr)	*110*		*77 - 120*		*10/01/25 04:01*	*1*
4-Bromofluorobenzene (Surr)	*103*		*73 - 120*		*10/01/25 04:01*	*1*
Dibromofluoromethane (Surr)	*109*		*75 - 123*		*10/01/25 04:01*	*1*
Toluene-d8 (Surr)	*100*		*80 - 120*		*10/01/25 04:01*	*1*

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)	ND		3.3	0.82	ng/L		10/03/25 13:46	10/06/25 06:31	1
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)	ND		3.3	0.82	ng/L		10/03/25 13:46	10/06/25 06:31	1
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	**39**		3.3	0.82	ng/L		10/03/25 13:46	10/06/25 06:31	1
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	ND		8.2	2.1	ng/L		10/03/25 13:46	10/06/25 06:31	1
3-Perfluoropentylpropanoic acid (5:3 FTCA)	ND		8.2	2.3	ng/L		10/03/25 13:46	10/06/25 06:31	1
3-Perfluoropropylpropanoic acid (3:3 FTCA)	ND		3.3	0.82	ng/L		10/03/25 13:46	10/06/25 06:31	1
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	ND		1.6	1.1	ng/L		10/03/25 13:46	10/06/25 06:31	1
N-ethylperfluorooctane sulfonamide (NEtFOSA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	ND		8.2	3.3	ng/L		10/03/25 13:46	10/06/25 06:31	1
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
N-methylperfluorooctane sulfonamide (NMeFOSA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	ND		8.2	2.1	ng/L		10/03/25 13:46	10/06/25 06:31	1
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluoro-3-methoxypropanoic acid (PFMPA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluoro-4-methoxybutanoic acid (PFMBA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorobutanesulfonic acid (PFBS)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorobutanoic acid (PFBA)	**1.7**	**J**	3.3	0.90	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorodecanesulfonic acid (PFDS)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorodecanoic acid (PFDA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorododecanesulfonic acid (PFDoS)	ND		1.6	0.49	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorododecanoic acid (PFDoA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluoroheptanesulfonic acid (PFHpS)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1

Client Sample ID: MW5I
Date Collected: 09/24/25 10:40
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-8
Matrix: Water

Method: EPA 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	**1.5**	**J**	1.6	0.66	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorohexanesulfonic acid (PFHxS)	**1.1**	**J**	1.6	0.66	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorohexanoic acid (PFHxA)	**4.4**		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorononanesulfonic acid (PFNS)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorononanoic acid (PFNA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorooctanesulfonamide (PFOSA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorooctanesulfonic acid (PFOS)	ND		1.6	1.1	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorooctanoic acid (PFOA)	ND		1.6	0.74	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluoropentanesulfonic acid (PFPeS)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluoropentanoic acid (PFPeA)	**4.9**		1.6	0.49	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorotetradecanoic acid (PFTeDA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluorotridecanoic acid (PFTrDA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1
Perfluoroundecanoic acid (PFUnA)	ND		1.6	0.41	ng/L		10/03/25 13:46	10/06/25 06:31	1

Isotope Dilution	*%Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
13C2 4:2 FTS	*111*		*40 - 200*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C2 6:2 FTS	*92.6*		*40 - 200*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C2 8:2 FTS	*81.3*		*40 - 300*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C2 PFDoA	*87.4*		*10 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C2 PFTeDA	*66.0*		*10 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C3 HFPO-DA	*108*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C3 PFBS	*103*		*40 - 135*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C3 PFHxS	*96.6*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C4 PFBA	*107*		*5 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C4 PFHpA	*102*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C5 PFHxA	*96.2*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C5 PFPeA	*100*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C6 PFDA	*104*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C7 PFUnA	*90.2*		*30 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C8 FOSA	*87.5*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C8 PFOA	*101*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C8 PFOS	*94.0*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
13C9 PFNA	*95.8*		*40 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
d3-NMeFOSAA	*81.4*		*40 - 170*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
d3-NMePFOSA	*68.6*		*10 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
d5-NEtFOSAA	*73.0*		*25 - 135*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
d5-NEtPFOSA	*63.1*		*10 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
d7-N-MeFOSE-M	*21.5*		*10 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*
d9-N-EtFOSE-M	*14.5*		*10 - 130*	*10/03/25 13:46*	*10/06/25 06:31*	*1*

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW6

Date Collected: 09/23/25 14:20
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-9

Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		1.0	0.82	ug/L			10/01/25 04:24	1
1,1,2,2-Tetrachloroethane	ND		1.0	0.21	ug/L			10/01/25 04:24	1
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1.0	0.31	ug/L			10/01/25 04:24	1
1,1,2-Trichloroethane	ND		1.0	0.23	ug/L			10/01/25 04:24	1
1,1-Dichloroethane	ND		1.0	0.38	ug/L			10/01/25 04:24	1
1,1-Dichloroethene	ND		1.0	0.29	ug/L			10/01/25 04:24	1
1,2,4-Trichlorobenzene	ND		1.0	0.41	ug/L			10/01/25 04:24	1
1,2-Dibromo-3-Chloropropane	ND		1.0	0.39	ug/L			10/01/25 04:24	1
1,2-Dibromoethane	ND		1.0	0.73	ug/L			10/01/25 04:24	1
1,2-Dichlorobenzene	ND		1.0	0.79	ug/L			10/01/25 04:24	1
1,2-Dichloroethane	ND		1.0	0.21	ug/L			10/01/25 04:24	1
1,2-Dichloropropane	ND		1.0	0.72	ug/L			10/01/25 04:24	1
1,3-Dichlorobenzene	ND		1.0	0.78	ug/L			10/01/25 04:24	1
1,4-Dichlorobenzene	ND		1.0	0.84	ug/L			10/01/25 04:24	1
2-Butanone (MEK)	ND		10	1.3	ug/L			10/01/25 04:24	1
2-Hexanone	ND		5.0	1.2	ug/L			10/01/25 04:24	1
4-Methyl-2-pentanone (MIBK)	ND		5.0	2.1	ug/L			10/01/25 04:24	1
Acetone	ND		10	3.0	ug/L			10/01/25 04:24	1
Benzene	ND		1.0	0.41	ug/L			10/01/25 04:24	1
Bromodichloromethane	ND		1.0	0.39	ug/L			10/01/25 04:24	1
Bromoform	ND		1.0	0.26	ug/L			10/01/25 04:24	1
Bromomethane	ND		1.0	0.69	ug/L			10/01/25 04:24	1
Carbon disulfide	ND		1.0	0.19	ug/L			10/01/25 04:24	1
Carbon tetrachloride	ND		1.0	0.27	ug/L			10/01/25 04:24	1
Chlorobenzene	ND		1.0	0.75	ug/L			10/01/25 04:24	1
Chloroethane	ND		1.0	0.32	ug/L			10/01/25 04:24	1
Chloroform	ND		1.0	0.34	ug/L			10/01/25 04:24	1
Chloromethane	ND		1.0	0.35	ug/L			10/01/25 04:24	1
cis-1,2-Dichloroethene	ND		1.0	0.81	ug/L			10/01/25 04:24	1
cis-1,3-Dichloropropene	ND		1.0	0.36	ug/L			10/01/25 04:24	1
Cyclohexane	ND		1.0	0.18	ug/L			10/01/25 04:24	1
Dibromochloromethane	ND		1.0	0.32	ug/L			10/01/25 04:24	1
Dichlorodifluoromethane	ND		1.0	0.68	ug/L			10/01/25 04:24	1
Ethylbenzene	ND		1.0	0.74	ug/L			10/01/25 04:24	1
Isopropylbenzene	ND		1.0	0.79	ug/L			10/01/25 04:24	1
Methyl acetate	ND		2.5	1.3	ug/L			10/01/25 04:24	1
Methyl tert-butyl ether	ND		1.0	0.16	ug/L			10/01/25 04:24	1
Methylcyclohexane	ND		1.0	0.16	ug/L			10/01/25 04:24	1
Methylene Chloride	ND		1.0	0.44	ug/L			10/01/25 04:24	1
Styrene	ND		1.0	0.73	ug/L			10/01/25 04:24	1
Tetrachloroethene	ND		1.0	0.36	ug/L			10/01/25 04:24	1
Toluene	ND		1.0	0.51	ug/L			10/01/25 04:24	1
trans-1,2-Dichloroethene	ND		1.0	0.90	ug/L			10/01/25 04:24	1
trans-1,3-Dichloropropene	ND		1.0	0.37	ug/L			10/01/25 04:24	1
Trichloroethene	ND		1.0	0.46	ug/L			10/01/25 04:24	1
Trichlorofluoromethane	ND		1.0	0.88	ug/L			10/01/25 04:24	1
Vinyl chloride	ND		1.0	0.90	ug/L			10/01/25 04:24	1
Xylenes, Total	ND		2.0	0.66	ug/L			10/01/25 04:24	1

Client: GEI Consultants Inc
Job ID: 480-232963-1
Project/Site: Unifirst – Liverpool, NY

Client Sample ID: MW6
Lab Sample ID: 480-232963-9
Date Collected: 09/23/25 14:20
Matrix: Water
Date Received: 09/26/25 10:00

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	109		77 - 120		10/01/25 04:24	1
4-Bromofluorobenzene (Surr)	104		73 - 120		10/01/25 04:24	1
Dibromofluoromethane (Surr)	109		75 - 123		10/01/25 04:24	1
Toluene-d8 (Surr)	103		80 - 120		10/01/25 04:24	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW7

Date Collected: 09/23/25 15:35
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-10

Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		1.0	0.82	ug/L			10/01/25 04:47	1
1,1,2,2-Tetrachloroethane	ND		1.0	0.21	ug/L			10/01/25 04:47	1
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1.0	0.31	ug/L			10/01/25 04:47	1
1,1,2-Trichloroethane	ND		1.0	0.23	ug/L			10/01/25 04:47	1
1,1-Dichloroethane	ND		1.0	0.38	ug/L			10/01/25 04:47	1
1,1-Dichloroethene	ND		1.0	0.29	ug/L			10/01/25 04:47	1
1,2,4-Trichlorobenzene	ND		1.0	0.41	ug/L			10/01/25 04:47	1
1,2-Dibromo-3-Chloropropane	ND		1.0	0.39	ug/L			10/01/25 04:47	1
1,2-Dibromoethane	ND		1.0	0.73	ug/L			10/01/25 04:47	1
1,2-Dichlorobenzene	ND		1.0	0.79	ug/L			10/01/25 04:47	1
1,2-Dichloroethane	ND		1.0	0.21	ug/L			10/01/25 04:47	1
1,2-Dichloropropane	ND		1.0	0.72	ug/L			10/01/25 04:47	1
1,3-Dichlorobenzene	ND		1.0	0.78	ug/L			10/01/25 04:47	1
1,4-Dichlorobenzene	ND		1.0	0.84	ug/L			10/01/25 04:47	1
2-Butanone (MEK)	ND		10	1.3	ug/L			10/01/25 04:47	1
2-Hexanone	ND		5.0	1.2	ug/L			10/01/25 04:47	1
4-Methyl-2-pentanone (MIBK)	ND		5.0	2.1	ug/L			10/01/25 04:47	1
Acetone	ND		10	3.0	ug/L			10/01/25 04:47	1
Benzene	ND		1.0	0.41	ug/L			10/01/25 04:47	1
Bromodichloromethane	ND		1.0	0.39	ug/L			10/01/25 04:47	1
Bromoform	ND		1.0	0.26	ug/L			10/01/25 04:47	1
Bromomethane	ND		1.0	0.69	ug/L			10/01/25 04:47	1
Carbon disulfide	ND		1.0	0.19	ug/L			10/01/25 04:47	1
Carbon tetrachloride	ND		1.0	0.27	ug/L			10/01/25 04:47	1
Chlorobenzene	ND		1.0	0.75	ug/L			10/01/25 04:47	1
Chloroethane	ND		1.0	0.32	ug/L			10/01/25 04:47	1
Chloroform	ND		1.0	0.34	ug/L			10/01/25 04:47	1
Chloromethane	ND		1.0	0.35	ug/L			10/01/25 04:47	1
cis-1,2-Dichloroethene	ND		1.0	0.81	ug/L			10/01/25 04:47	1
cis-1,3-Dichloropropene	ND		1.0	0.36	ug/L			10/01/25 04:47	1
Cyclohexane	ND		1.0	0.18	ug/L			10/01/25 04:47	1
Dibromochloromethane	ND		1.0	0.32	ug/L			10/01/25 04:47	1
Dichlorodifluoromethane	ND		1.0	0.68	ug/L			10/01/25 04:47	1
Ethylbenzene	ND		1.0	0.74	ug/L			10/01/25 04:47	1
Isopropylbenzene	ND		1.0	0.79	ug/L			10/01/25 04:47	1
Methyl acetate	ND		2.5	1.3	ug/L			10/01/25 04:47	1
Methyl tert-butyl ether	ND		1.0	0.16	ug/L			10/01/25 04:47	1
Methylcyclohexane	ND		1.0	0.16	ug/L			10/01/25 04:47	1
Methylene Chloride	ND		1.0	0.44	ug/L			10/01/25 04:47	1
Styrene	ND		1.0	0.73	ug/L			10/01/25 04:47	1
Tetrachloroethene	ND		1.0	0.36	ug/L			10/01/25 04:47	1
Toluene	ND		1.0	0.51	ug/L			10/01/25 04:47	1
trans-1,2-Dichloroethene	ND		1.0	0.90	ug/L			10/01/25 04:47	1
trans-1,3-Dichloropropene	ND		1.0	0.37	ug/L			10/01/25 04:47	1
Trichloroethene	ND		1.0	0.46	ug/L			10/01/25 04:47	1
Trichlorofluoromethane	ND		1.0	0.88	ug/L			10/01/25 04:47	1
Vinyl chloride	**13**		1.0	0.90	ug/L			10/01/25 04:47	1
Xylenes, Total	ND		2.0	0.66	ug/L			10/01/25 04:47	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW7
Date Collected: 09/23/25 15:35
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-10
Matrix: Water

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	109		77 - 120		10/01/25 04:47	1
4-Bromofluorobenzene (Surr)	103		73 - 120		10/01/25 04:47	1
Dibromofluoromethane (Surr)	108		75 - 123		10/01/25 04:47	1
Toluene-d8 (Surr)	100		80 - 120		10/01/25 04:47	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW8

Date Collected: 09/24/25 10:30
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-11

Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		1.0	0.82	ug/L			10/01/25 05:10	1
1,1,2,2-Tetrachloroethane	ND		1.0	0.21	ug/L			10/01/25 05:10	1
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1.0	0.31	ug/L			10/01/25 05:10	1
1,1,2-Trichloroethane	ND		1.0	0.23	ug/L			10/01/25 05:10	1
1,1-Dichloroethane	ND		1.0	0.38	ug/L			10/01/25 05:10	1
1,1-Dichloroethene	ND		1.0	0.29	ug/L			10/01/25 05:10	1
1,2,4-Trichlorobenzene	ND		1.0	0.41	ug/L			10/01/25 05:10	1
1,2-Dibromo-3-Chloropropane	ND		1.0	0.39	ug/L			10/01/25 05:10	1
1,2-Dibromoethane	ND		1.0	0.73	ug/L			10/01/25 05:10	1
1,2-Dichlorobenzene	ND		1.0	0.79	ug/L			10/01/25 05:10	1
1,2-Dichloroethane	ND		1.0	0.21	ug/L			10/01/25 05:10	1
1,2-Dichloropropane	ND		1.0	0.72	ug/L			10/01/25 05:10	1
1,3-Dichlorobenzene	ND		1.0	0.78	ug/L			10/01/25 05:10	1
1,4-Dichlorobenzene	ND		1.0	0.84	ug/L			10/01/25 05:10	1
2-Butanone (MEK)	ND		10	1.3	ug/L			10/01/25 05:10	1
2-Hexanone	ND		5.0	1.2	ug/L			10/01/25 05:10	1
4-Methyl-2-pentanone (MIBK)	ND		5.0	2.1	ug/L			10/01/25 05:10	1
Acetone	**3.5**	**J**	10	3.0	ug/L			10/01/25 05:10	1
Benzene	ND		1.0	0.41	ug/L			10/01/25 05:10	1
Bromodichloromethane	ND		1.0	0.39	ug/L			10/01/25 05:10	1
Bromoform	ND		1.0	0.26	ug/L			10/01/25 05:10	1
Bromomethane	ND		1.0	0.69	ug/L			10/01/25 05:10	1
Carbon disulfide	ND		1.0	0.19	ug/L			10/01/25 05:10	1
Carbon tetrachloride	ND		1.0	0.27	ug/L			10/01/25 05:10	1
Chlorobenzene	ND		1.0	0.75	ug/L			10/01/25 05:10	1
Chloroethane	ND		1.0	0.32	ug/L			10/01/25 05:10	1
Chloroform	ND		1.0	0.34	ug/L			10/01/25 05:10	1
Chloromethane	ND		1.0	0.35	ug/L			10/01/25 05:10	1
cis-1,2-Dichloroethene	ND		1.0	0.81	ug/L			10/01/25 05:10	1
cis-1,3-Dichloropropene	ND		1.0	0.36	ug/L			10/01/25 05:10	1
Cyclohexane	ND		1.0	0.18	ug/L			10/01/25 05:10	1
Dibromochloromethane	ND		1.0	0.32	ug/L			10/01/25 05:10	1
Dichlorodifluoromethane	ND		1.0	0.68	ug/L			10/01/25 05:10	1
Ethylbenzene	ND		1.0	0.74	ug/L			10/01/25 05:10	1
Isopropylbenzene	ND		1.0	0.79	ug/L			10/01/25 05:10	1
Methyl acetate	ND		2.5	1.3	ug/L			10/01/25 05:10	1
Methyl tert-butyl ether	ND		1.0	0.16	ug/L			10/01/25 05:10	1
Methylcyclohexane	ND		1.0	0.16	ug/L			10/01/25 05:10	1
Methylene Chloride	ND		1.0	0.44	ug/L			10/01/25 05:10	1
Styrene	ND		1.0	0.73	ug/L			10/01/25 05:10	1
Tetrachloroethene	ND		1.0	0.36	ug/L			10/01/25 05:10	1
Toluene	ND		1.0	0.51	ug/L			10/01/25 05:10	1
trans-1,2-Dichloroethene	ND		1.0	0.90	ug/L			10/01/25 05:10	1
trans-1,3-Dichloropropene	ND		1.0	0.37	ug/L			10/01/25 05:10	1
Trichloroethene	ND		1.0	0.46	ug/L			10/01/25 05:10	1
Trichlorofluoromethane	ND		1.0	0.88	ug/L			10/01/25 05:10	1
Vinyl chloride	ND		1.0	0.90	ug/L			10/01/25 05:10	1
Xylenes, Total	ND		2.0	0.66	ug/L			10/01/25 05:10	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW8
Date Collected: 09/24/25 10:30
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-11
Matrix: Water

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	112		77 - 120		10/01/25 05:10	1
4-Bromofluorobenzene (Surr)	106		73 - 120		10/01/25 05:10	1
Dibromofluoromethane (Surr)	105		75 - 123		10/01/25 05:10	1
Toluene-d8 (Surr)	103		80 - 120		10/01/25 05:10	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW11

Date Collected: 09/24/25 16:05
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-12

Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		1.0	0.82	ug/L			10/01/25 05:33	1
1,1,2,2-Tetrachloroethane	ND		1.0	0.21	ug/L			10/01/25 05:33	1
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1.0	0.31	ug/L			10/01/25 05:33	1
1,1,2-Trichloroethane	ND		1.0	0.23	ug/L			10/01/25 05:33	1
1,1-Dichloroethane	ND		1.0	0.38	ug/L			10/01/25 05:33	1
1,1-Dichloroethene	ND		1.0	0.29	ug/L			10/01/25 05:33	1
1,2,4-Trichlorobenzene	ND		1.0	0.41	ug/L			10/01/25 05:33	1
1,2-Dibromo-3-Chloropropane	ND		1.0	0.39	ug/L			10/01/25 05:33	1
1,2-Dibromoethane	ND		1.0	0.73	ug/L			10/01/25 05:33	1
1,2-Dichlorobenzene	ND		1.0	0.79	ug/L			10/01/25 05:33	1
1,2-Dichloroethane	ND		1.0	0.21	ug/L			10/01/25 05:33	1
1,2-Dichloropropane	ND		1.0	0.72	ug/L			10/01/25 05:33	1
1,3-Dichlorobenzene	ND		1.0	0.78	ug/L			10/01/25 05:33	1
1,4-Dichlorobenzene	ND		1.0	0.84	ug/L			10/01/25 05:33	1
2-Butanone (MEK)	ND		10	1.3	ug/L			10/01/25 05:33	1
2-Hexanone	ND		5.0	1.2	ug/L			10/01/25 05:33	1
4-Methyl-2-pentanone (MIBK)	ND		5.0	2.1	ug/L			10/01/25 05:33	1
Acetone	ND		10	3.0	ug/L			10/01/25 05:33	1
Benzene	ND		1.0	0.41	ug/L			10/01/25 05:33	1
Bromodichloromethane	ND		1.0	0.39	ug/L			10/01/25 05:33	1
Bromoform	ND		1.0	0.26	ug/L			10/01/25 05:33	1
Bromomethane	ND		1.0	0.69	ug/L			10/01/25 05:33	1
Carbon disulfide	ND		1.0	0.19	ug/L			10/01/25 05:33	1
Carbon tetrachloride	ND		1.0	0.27	ug/L			10/01/25 05:33	1
Chlorobenzene	ND		1.0	0.75	ug/L			10/01/25 05:33	1
Chloroethane	ND		1.0	0.32	ug/L			10/01/25 05:33	1
Chloroform	ND		1.0	0.34	ug/L			10/01/25 05:33	1
Chloromethane	ND		1.0	0.35	ug/L			10/01/25 05:33	1
cis-1,2-Dichloroethene	ND		1.0	0.81	ug/L			10/01/25 05:33	1
cis-1,3-Dichloropropene	ND		1.0	0.36	ug/L			10/01/25 05:33	1
Cyclohexane	ND		1.0	0.18	ug/L			10/01/25 05:33	1
Dibromochloromethane	ND		1.0	0.32	ug/L			10/01/25 05:33	1
Dichlorodifluoromethane	ND		1.0	0.68	ug/L			10/01/25 05:33	1
Ethylbenzene	ND		1.0	0.74	ug/L			10/01/25 05:33	1
Isopropylbenzene	ND		1.0	0.79	ug/L			10/01/25 05:33	1
Methyl acetate	ND		2.5	1.3	ug/L			10/01/25 05:33	1
Methyl tert-butyl ether	ND		1.0	0.16	ug/L			10/01/25 05:33	1
Methylcyclohexane	ND		1.0	0.16	ug/L			10/01/25 05:33	1
Methylene Chloride	ND		1.0	0.44	ug/L			10/01/25 05:33	1
Styrene	ND		1.0	0.73	ug/L			10/01/25 05:33	1
Tetrachloroethene	ND		1.0	0.36	ug/L			10/01/25 05:33	1
Toluene	ND		1.0	0.51	ug/L			10/01/25 05:33	1
trans-1,2-Dichloroethene	ND		1.0	0.90	ug/L			10/01/25 05:33	1
trans-1,3-Dichloropropene	ND		1.0	0.37	ug/L			10/01/25 05:33	1
Trichloroethene	ND		1.0	0.46	ug/L			10/01/25 05:33	1
Trichlorofluoromethane	ND		1.0	0.88	ug/L			10/01/25 05:33	1
Vinyl chloride	ND		1.0	0.90	ug/L			10/01/25 05:33	1
Xylenes, Total	ND		2.0	0.66	ug/L			10/01/25 05:33	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW11
Date Collected: 09/24/25 16:05
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-12
Matrix: Water

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	110		77 - 120		10/01/25 05:33	1
4-Bromofluorobenzene (Surr)	104		73 - 120		10/01/25 05:33	1
Dibromofluoromethane (Surr)	107		75 - 123		10/01/25 05:33	1
Toluene-d8 (Surr)	98		80 - 120		10/01/25 05:33	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW12S

Date Collected: 09/24/25 16:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-13

Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		1.0	0.82	ug/L			10/01/25 05:56	1
1,1,2,2-Tetrachloroethane	ND		1.0	0.21	ug/L			10/01/25 05:56	1
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1.0	0.31	ug/L			10/01/25 05:56	1
1,1,2-Trichloroethane	ND		1.0	0.23	ug/L			10/01/25 05:56	1
1,1-Dichloroethane	ND		1.0	0.38	ug/L			10/01/25 05:56	1
1,1-Dichloroethene	ND		1.0	0.29	ug/L			10/01/25 05:56	1
1,2,4-Trichlorobenzene	ND		1.0	0.41	ug/L			10/01/25 05:56	1
1,2-Dibromo-3-Chloropropane	ND		1.0	0.39	ug/L			10/01/25 05:56	1
1,2-Dibromoethane	ND		1.0	0.73	ug/L			10/01/25 05:56	1
1,2-Dichlorobenzene	ND		1.0	0.79	ug/L			10/01/25 05:56	1
1,2-Dichloroethane	ND		1.0	0.21	ug/L			10/01/25 05:56	1
1,2-Dichloropropane	ND		1.0	0.72	ug/L			10/01/25 05:56	1
1,3-Dichlorobenzene	ND		1.0	0.78	ug/L			10/01/25 05:56	1
1,4-Dichlorobenzene	ND		1.0	0.84	ug/L			10/01/25 05:56	1
2-Butanone (MEK)	ND		10	1.3	ug/L			10/01/25 05:56	1
2-Hexanone	ND		5.0	1.2	ug/L			10/01/25 05:56	1
4-Methyl-2-pentanone (MIBK)	ND		5.0	2.1	ug/L			10/01/25 05:56	1
Acetone	ND		10	3.0	ug/L			10/01/25 05:56	1
Benzene	ND		1.0	0.41	ug/L			10/01/25 05:56	1
Bromodichloromethane	ND		1.0	0.39	ug/L			10/01/25 05:56	1
Bromoform	ND		1.0	0.26	ug/L			10/01/25 05:56	1
Bromomethane	ND		1.0	0.69	ug/L			10/01/25 05:56	1
Carbon disulfide	ND		1.0	0.19	ug/L			10/01/25 05:56	1
Carbon tetrachloride	ND		1.0	0.27	ug/L			10/01/25 05:56	1
Chlorobenzene	ND		1.0	0.75	ug/L			10/01/25 05:56	1
Chloroethane	ND		1.0	0.32	ug/L			10/01/25 05:56	1
Chloroform	ND		1.0	0.34	ug/L			10/01/25 05:56	1
Chloromethane	ND		1.0	0.35	ug/L			10/01/25 05:56	1
cis-1,2-Dichloroethene	**2.3**		1.0	0.81	ug/L			10/01/25 05:56	1
cis-1,3-Dichloropropene	ND		1.0	0.36	ug/L			10/01/25 05:56	1
Cyclohexane	ND		1.0	0.18	ug/L			10/01/25 05:56	1
Dibromochloromethane	ND		1.0	0.32	ug/L			10/01/25 05:56	1
Dichlorodifluoromethane	ND		1.0	0.68	ug/L			10/01/25 05:56	1
Ethylbenzene	ND		1.0	0.74	ug/L			10/01/25 05:56	1
Isopropylbenzene	ND		1.0	0.79	ug/L			10/01/25 05:56	1
Methyl acetate	ND		2.5	1.3	ug/L			10/01/25 05:56	1
Methyl tert-butyl ether	ND		1.0	0.16	ug/L			10/01/25 05:56	1
Methylcyclohexane	ND		1.0	0.16	ug/L			10/01/25 05:56	1
Methylene Chloride	ND		1.0	0.44	ug/L			10/01/25 05:56	1
Styrene	ND		1.0	0.73	ug/L			10/01/25 05:56	1
Tetrachloroethene	ND		1.0	0.36	ug/L			10/01/25 05:56	1
Toluene	ND		1.0	0.51	ug/L			10/01/25 05:56	1
trans-1,2-Dichloroethene	ND		1.0	0.90	ug/L			10/01/25 05:56	1
trans-1,3-Dichloropropene	ND		1.0	0.37	ug/L			10/01/25 05:56	1
Trichloroethene	ND		1.0	0.46	ug/L			10/01/25 05:56	1
Trichlorofluoromethane	ND		1.0	0.88	ug/L			10/01/25 05:56	1
Vinyl chloride	**1.5**		1.0	0.90	ug/L			10/01/25 05:56	1
Xylenes, Total	ND		2.0	0.66	ug/L			10/01/25 05:56	1

Client: GEI Consultants Inc
Job ID: 480-232963-1
Project/Site: Unifirst – Liverpool, NY

Client Sample ID: MW12S
Lab Sample ID: 480-232963-13
Date Collected: 09/24/25 16:00
Matrix: Water
Date Received: 09/26/25 10:00

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	113		77 - 120		10/01/25 05:56	1
4-Bromofluorobenzene (Surr)	103		73 - 120		10/01/25 05:56	1
Dibromofluoromethane (Surr)	111		75 - 123		10/01/25 05:56	1
Toluene-d8 (Surr)	103		80 - 120		10/01/25 05:56	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW1C

Date Collected: 09/25/25 12:15
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-14

Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		1.0	0.82	ug/L			10/01/25 06:19	1
1,1,2,2-Tetrachloroethane	ND		1.0	0.21	ug/L			10/01/25 06:19	1
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1.0	0.31	ug/L			10/01/25 06:19	1
1,1,2-Trichloroethane	ND		1.0	0.23	ug/L			10/01/25 06:19	1
1,1-Dichloroethane	ND		1.0	0.38	ug/L			10/01/25 06:19	1
1,1-Dichloroethene	ND		1.0	0.29	ug/L			10/01/25 06:19	1
1,2,4-Trichlorobenzene	ND		1.0	0.41	ug/L			10/01/25 06:19	1
1,2-Dibromo-3-Chloropropane	ND		1.0	0.39	ug/L			10/01/25 06:19	1
1,2-Dibromoethane	ND		1.0	0.73	ug/L			10/01/25 06:19	1
1,2-Dichlorobenzene	ND		1.0	0.79	ug/L			10/01/25 06:19	1
1,2-Dichloroethane	ND		1.0	0.21	ug/L			10/01/25 06:19	1
1,2-Dichloropropane	ND		1.0	0.72	ug/L			10/01/25 06:19	1
1,3-Dichlorobenzene	ND		1.0	0.78	ug/L			10/01/25 06:19	1
1,4-Dichlorobenzene	ND		1.0	0.84	ug/L			10/01/25 06:19	1
2-Butanone (MEK)	ND		10	1.3	ug/L			10/01/25 06:19	1
2-Hexanone	ND		5.0	1.2	ug/L			10/01/25 06:19	1
4-Methyl-2-pentanone (MIBK)	ND		5.0	2.1	ug/L			10/01/25 06:19	1
Acetone	**5.6**	**J**	10	3.0	ug/L			10/01/25 06:19	1
Benzene	ND		1.0	0.41	ug/L			10/01/25 06:19	1
Bromodichloromethane	ND		1.0	0.39	ug/L			10/01/25 06:19	1
Bromoform	ND		1.0	0.26	ug/L			10/01/25 06:19	1
Bromomethane	ND		1.0	0.69	ug/L			10/01/25 06:19	1
Carbon disulfide	ND		1.0	0.19	ug/L			10/01/25 06:19	1
Carbon tetrachloride	ND		1.0	0.27	ug/L			10/01/25 06:19	1
Chlorobenzene	ND		1.0	0.75	ug/L			10/01/25 06:19	1
Chloroethane	ND		1.0	0.32	ug/L			10/01/25 06:19	1
Chloroform	ND		1.0	0.34	ug/L			10/01/25 06:19	1
Chloromethane	ND		1.0	0.35	ug/L			10/01/25 06:19	1
cis-1,2-Dichloroethene	ND		1.0	0.81	ug/L			10/01/25 06:19	1
cis-1,3-Dichloropropene	ND		1.0	0.36	ug/L			10/01/25 06:19	1
Cyclohexane	ND		1.0	0.18	ug/L			10/01/25 06:19	1
Dibromochloromethane	ND		1.0	0.32	ug/L			10/01/25 06:19	1
Dichlorodifluoromethane	ND		1.0	0.68	ug/L			10/01/25 06:19	1
Ethylbenzene	ND		1.0	0.74	ug/L			10/01/25 06:19	1
Isopropylbenzene	ND		1.0	0.79	ug/L			10/01/25 06:19	1
Methyl acetate	ND		2.5	1.3	ug/L			10/01/25 06:19	1
Methyl tert-butyl ether	ND		1.0	0.16	ug/L			10/01/25 06:19	1
Methylcyclohexane	ND		1.0	0.16	ug/L			10/01/25 06:19	1
Methylene Chloride	ND		1.0	0.44	ug/L			10/01/25 06:19	1
Styrene	ND		1.0	0.73	ug/L			10/01/25 06:19	1
Tetrachloroethene	ND		1.0	0.36	ug/L			10/01/25 06:19	1
Toluene	ND		1.0	0.51	ug/L			10/01/25 06:19	1
trans-1,2-Dichloroethene	ND		1.0	0.90	ug/L			10/01/25 06:19	1
trans-1,3-Dichloropropene	ND		1.0	0.37	ug/L			10/01/25 06:19	1
Trichloroethene	ND		1.0	0.46	ug/L			10/01/25 06:19	1
Trichlorofluoromethane	ND		1.0	0.88	ug/L			10/01/25 06:19	1
Vinyl chloride	ND		1.0	0.90	ug/L			10/01/25 06:19	1
Xylenes, Total	ND		2.0	0.66	ug/L			10/01/25 06:19	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW1C
Date Collected: 09/25/25 12:15
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-14
Matrix: Water

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	110		77 - 120		10/01/25 06:19	1
4-Bromofluorobenzene (Surr)	102		73 - 120		10/01/25 06:19	1
Dibromofluoromethane (Surr)	107		75 - 123		10/01/25 06:19	1
Toluene-d8 (Surr)	102		80 - 120		10/01/25 06:19	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW1D

Date Collected: 09/25/25 09:15
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-15

Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		20	16	ug/L			10/01/25 06:43	20
1,1,2,2-Tetrachloroethane	ND		20	4.2	ug/L			10/01/25 06:43	20
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		20	6.2	ug/L			10/01/25 06:43	20
1,1,2-Trichloroethane	ND		20	4.6	ug/L			10/01/25 06:43	20
1,1-Dichloroethane	ND		20	7.6	ug/L			10/01/25 06:43	20
1,1-Dichloroethene	ND		20	5.8	ug/L			10/01/25 06:43	20
1,2,4-Trichlorobenzene	ND		20	8.2	ug/L			10/01/25 06:43	20
1,2-Dibromo-3-Chloropropane	ND		20	7.8	ug/L			10/01/25 06:43	20
1,2-Dibromoethane	ND		20	15	ug/L			10/01/25 06:43	20
1,2-Dichlorobenzene	ND		20	16	ug/L			10/01/25 06:43	20
1,2-Dichloroethane	ND		20	4.2	ug/L			10/01/25 06:43	20
1,2-Dichloropropane	ND		20	14	ug/L			10/01/25 06:43	20
1,3-Dichlorobenzene	ND		20	16	ug/L			10/01/25 06:43	20
1,4-Dichlorobenzene	ND		20	17	ug/L			10/01/25 06:43	20
2-Butanone (MEK)	ND		200	26	ug/L			10/01/25 06:43	20
2-Hexanone	ND		100	25	ug/L			10/01/25 06:43	20
4-Methyl-2-pentanone (MIBK)	ND		100	42	ug/L			10/01/25 06:43	20
Acetone	ND		200	60	ug/L			10/01/25 06:43	20
Benzene	ND		20	8.2	ug/L			10/01/25 06:43	20
Bromodichloromethane	ND		20	7.8	ug/L			10/01/25 06:43	20
Bromoform	ND		20	5.2	ug/L			10/01/25 06:43	20
Bromomethane	ND		20	14	ug/L			10/01/25 06:43	20
Carbon disulfide	ND		20	3.8	ug/L			10/01/25 06:43	20
Carbon tetrachloride	ND		20	5.4	ug/L			10/01/25 06:43	20
Chlorobenzene	ND		20	15	ug/L			10/01/25 06:43	20
Chloroethane	ND		20	6.4	ug/L			10/01/25 06:43	20
Chloroform	ND		20	6.8	ug/L			10/01/25 06:43	20
Chloromethane	ND		20	7.0	ug/L			10/01/25 06:43	20
cis-1,2-Dichloroethene	**990**		20	16	ug/L			10/01/25 06:43	20
cis-1,3-Dichloropropene	ND		20	7.2	ug/L			10/01/25 06:43	20
Cyclohexane	ND		20	3.6	ug/L			10/01/25 06:43	20
Dibromochloromethane	ND		20	6.4	ug/L			10/01/25 06:43	20
Dichlorodifluoromethane	ND		20	14	ug/L			10/01/25 06:43	20
Ethylbenzene	ND		20	15	ug/L			10/01/25 06:43	20
Isopropylbenzene	ND		20	16	ug/L			10/01/25 06:43	20
Methyl acetate	ND		50	26	ug/L			10/01/25 06:43	20
Methyl tert-butyl ether	ND		20	3.2	ug/L			10/01/25 06:43	20
Methylcyclohexane	ND		20	3.2	ug/L			10/01/25 06:43	20
Methylene Chloride	ND		20	8.8	ug/L			10/01/25 06:43	20
Styrene	ND		20	15	ug/L			10/01/25 06:43	20
Tetrachloroethene	**540**		20	7.2	ug/L			10/01/25 06:43	20
Toluene	ND		20	10	ug/L			10/01/25 06:43	20
trans-1,2-Dichloroethene	ND		20	18	ug/L			10/01/25 06:43	20
trans-1,3-Dichloropropene	ND		20	7.4	ug/L			10/01/25 06:43	20
Trichloroethene	**210**		20	9.2	ug/L			10/01/25 06:43	20
Trichlorofluoromethane	ND		20	18	ug/L			10/01/25 06:43	20
Vinyl chloride	**28**		20	18	ug/L			10/01/25 06:43	20
Xylenes, Total	ND		40	13	ug/L			10/01/25 06:43	20

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW1D
Date Collected: 09/25/25 09:15
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-15
Matrix: Water

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	110		77 - 120		10/01/25 06:43	20
4-Bromofluorobenzene (Surr)	107		73 - 120		10/01/25 06:43	20
Dibromofluoromethane (Surr)	108		75 - 123		10/01/25 06:43	20
Toluene-d8 (Surr)	103		80 - 120		10/01/25 06:43	20

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW1E

Date Collected: 09/25/25 11:05
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-16

Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		100	82	ug/L			10/01/25 07:06	100
1,1,2,2-Tetrachloroethane	ND		100	21	ug/L			10/01/25 07:06	100
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		100	31	ug/L			10/01/25 07:06	100
1,1,2-Trichloroethane	ND		100	23	ug/L			10/01/25 07:06	100
1,1-Dichloroethane	ND		100	38	ug/L			10/01/25 07:06	100
1,1-Dichloroethene	ND		100	29	ug/L			10/01/25 07:06	100
1,2,4-Trichlorobenzene	ND		100	41	ug/L			10/01/25 07:06	100
1,2-Dibromo-3-Chloropropane	ND		100	39	ug/L			10/01/25 07:06	100
1,2-Dibromoethane	ND		100	73	ug/L			10/01/25 07:06	100
1,2-Dichlorobenzene	ND		100	79	ug/L			10/01/25 07:06	100
1,2-Dichloroethane	ND		100	21	ug/L			10/01/25 07:06	100
1,2-Dichloropropane	ND		100	72	ug/L			10/01/25 07:06	100
1,3-Dichlorobenzene	ND		100	78	ug/L			10/01/25 07:06	100
1,4-Dichlorobenzene	ND		100	84	ug/L			10/01/25 07:06	100
2-Butanone (MEK)	ND		1000	130	ug/L			10/01/25 07:06	100
2-Hexanone	ND		500	120	ug/L			10/01/25 07:06	100
4-Methyl-2-pentanone (MIBK)	ND		500	210	ug/L			10/01/25 07:06	100
Acetone	ND		1000	300	ug/L			10/01/25 07:06	100
Benzene	ND		100	41	ug/L			10/01/25 07:06	100
Bromodichloromethane	ND		100	39	ug/L			10/01/25 07:06	100
Bromoform	ND		100	26	ug/L			10/01/25 07:06	100
Bromomethane	ND		100	69	ug/L			10/01/25 07:06	100
Carbon disulfide	ND		100	19	ug/L			10/01/25 07:06	100
Carbon tetrachloride	ND		100	27	ug/L			10/01/25 07:06	100
Chlorobenzene	ND		100	75	ug/L			10/01/25 07:06	100
Chloroethane	ND		100	32	ug/L			10/01/25 07:06	100
Chloroform	ND		100	34	ug/L			10/01/25 07:06	100
Chloromethane	ND		100	35	ug/L			10/01/25 07:06	100
cis-1,2-Dichloroethene	**2900**		100	81	ug/L			10/01/25 07:06	100
cis-1,3-Dichloropropene	ND		100	36	ug/L			10/01/25 07:06	100
Cyclohexane	ND		100	18	ug/L			10/01/25 07:06	100
Dibromochloromethane	ND		100	32	ug/L			10/01/25 07:06	100
Dichlorodifluoromethane	ND		100	68	ug/L			10/01/25 07:06	100
Ethylbenzene	ND		100	74	ug/L			10/01/25 07:06	100
Isopropylbenzene	ND		100	79	ug/L			10/01/25 07:06	100
Methyl acetate	ND		250	130	ug/L			10/01/25 07:06	100
Methyl tert-butyl ether	ND		100	16	ug/L			10/01/25 07:06	100
Methylcyclohexane	ND		100	16	ug/L			10/01/25 07:06	100
Methylene Chloride	ND		100	44	ug/L			10/01/25 07:06	100
Styrene	ND		100	73	ug/L			10/01/25 07:06	100
Tetrachloroethene	**3600**		100	36	ug/L			10/01/25 07:06	100
Toluene	ND		100	51	ug/L			10/01/25 07:06	100
trans-1,2-Dichloroethene	ND		100	90	ug/L			10/01/25 07:06	100
trans-1,3-Dichloropropene	ND		100	37	ug/L			10/01/25 07:06	100
Trichloroethene	**1100**		100	46	ug/L			10/01/25 07:06	100
Trichlorofluoromethane	ND		100	88	ug/L			10/01/25 07:06	100
Vinyl chloride	ND		100	90	ug/L			10/01/25 07:06	100
Xylenes, Total	ND		200	66	ug/L			10/01/25 07:06	100

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW1E
Date Collected: 09/25/25 11:05
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-16
Matrix: Water

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	109		77 - 120		10/01/25 07:06	100
4-Bromofluorobenzene (Surr)	103		73 - 120		10/01/25 07:06	100
Dibromofluoromethane (Surr)	111		75 - 123		10/01/25 07:06	100
Toluene-d8 (Surr)	97		80 - 120		10/01/25 07:06	100

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: TB

Date Collected: 09/25/25 00:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-17

Matrix: Water

Method: SW846 8260C - Volatile Organic Compounds by GC/MS

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		1.0	0.82	ug/L			10/01/25 07:29	1
1,1,2,2-Tetrachloroethane	ND		1.0	0.21	ug/L			10/01/25 07:29	1
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1.0	0.31	ug/L			10/01/25 07:29	1
1,1,2-Trichloroethane	ND		1.0	0.23	ug/L			10/01/25 07:29	1
1,1-Dichloroethane	ND		1.0	0.38	ug/L			10/01/25 07:29	1
1,1-Dichloroethene	ND		1.0	0.29	ug/L			10/01/25 07:29	1
1,2,4-Trichlorobenzene	ND		1.0	0.41	ug/L			10/01/25 07:29	1
1,2-Dibromo-3-Chloropropane	ND		1.0	0.39	ug/L			10/01/25 07:29	1
1,2-Dibromoethane	ND		1.0	0.73	ug/L			10/01/25 07:29	1
1,2-Dichlorobenzene	ND		1.0	0.79	ug/L			10/01/25 07:29	1
1,2-Dichloroethane	ND		1.0	0.21	ug/L			10/01/25 07:29	1
1,2-Dichloropropane	ND		1.0	0.72	ug/L			10/01/25 07:29	1
1,3-Dichlorobenzene	ND		1.0	0.78	ug/L			10/01/25 07:29	1
1,4-Dichlorobenzene	ND		1.0	0.84	ug/L			10/01/25 07:29	1
2-Butanone (MEK)	ND		10	1.3	ug/L			10/01/25 07:29	1
2-Hexanone	ND		5.0	1.2	ug/L			10/01/25 07:29	1
4-Methyl-2-pentanone (MIBK)	ND		5.0	2.1	ug/L			10/01/25 07:29	1
Acetone	ND		10	3.0	ug/L			10/01/25 07:29	1
Benzene	ND		1.0	0.41	ug/L			10/01/25 07:29	1
Bromodichloromethane	ND		1.0	0.39	ug/L			10/01/25 07:29	1
Bromoform	ND		1.0	0.26	ug/L			10/01/25 07:29	1
Bromomethane	ND		1.0	0.69	ug/L			10/01/25 07:29	1
Carbon disulfide	ND		1.0	0.19	ug/L			10/01/25 07:29	1
Carbon tetrachloride	ND		1.0	0.27	ug/L			10/01/25 07:29	1
Chlorobenzene	ND		1.0	0.75	ug/L			10/01/25 07:29	1
Chloroethane	ND		1.0	0.32	ug/L			10/01/25 07:29	1
Chloroform	ND		1.0	0.34	ug/L			10/01/25 07:29	1
Chloromethane	ND		1.0	0.35	ug/L			10/01/25 07:29	1
cis-1,2-Dichloroethene	ND		1.0	0.81	ug/L			10/01/25 07:29	1
cis-1,3-Dichloropropene	ND		1.0	0.36	ug/L			10/01/25 07:29	1
Cyclohexane	ND		1.0	0.18	ug/L			10/01/25 07:29	1
Dibromochloromethane	ND		1.0	0.32	ug/L			10/01/25 07:29	1
Dichlorodifluoromethane	ND		1.0	0.68	ug/L			10/01/25 07:29	1
Ethylbenzene	ND		1.0	0.74	ug/L			10/01/25 07:29	1
Isopropylbenzene	ND		1.0	0.79	ug/L			10/01/25 07:29	1
Methyl acetate	ND		2.5	1.3	ug/L			10/01/25 07:29	1
Methyl tert-butyl ether	ND		1.0	0.16	ug/L			10/01/25 07:29	1
Methylcyclohexane	ND		1.0	0.16	ug/L			10/01/25 07:29	1
Methylene Chloride	ND		1.0	0.44	ug/L			10/01/25 07:29	1
Styrene	ND		1.0	0.73	ug/L			10/01/25 07:29	1
Tetrachloroethene	ND		1.0	0.36	ug/L			10/01/25 07:29	1
Toluene	ND		1.0	0.51	ug/L			10/01/25 07:29	1
trans-1,2-Dichloroethene	ND		1.0	0.90	ug/L			10/01/25 07:29	1
trans-1,3-Dichloropropene	ND		1.0	0.37	ug/L			10/01/25 07:29	1
Trichloroethene	ND		1.0	0.46	ug/L			10/01/25 07:29	1
Trichlorofluoromethane	ND		1.0	0.88	ug/L			10/01/25 07:29	1
Vinyl chloride	ND		1.0	0.90	ug/L			10/01/25 07:29	1
Xylenes, Total	ND		2.0	0.66	ug/L			10/01/25 07:29	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: TB
Date Collected: 09/25/25 00:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-17
Matrix: Water

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	107		77 - 120		10/01/25 07:29	1
4-Bromofluorobenzene (Surr)	105		73 - 120		10/01/25 07:29	1
Dibromofluoromethane (Surr)	104		75 - 123		10/01/25 07:29	1
Toluene-d8 (Surr)	102		80 - 120		10/01/25 07:29	1

Surrogate Summary

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Method: 8260C - Volatile Organic Compounds by GC/MS

Matrix: Water

Prep Type: Total/NA

		Percent Surrogate Recovery (Acceptance Limits)			
		DCA	BFB	DBFM	TOL
Lab Sample ID	Client Sample ID	(77-120)	(73-120)	(75-123)	(80-120)
480-232963-1	MW1R	107	107	107	101
480-232963-2	MW2R	106	106	109	104
480-232963-3	MW3S	109	105	107	101
480-232963-4	MW3I	105	104	104	100
480-232963-5	MW4S	109	108	108	100
480-232963-6	MW4I	119	114	117	108
480-232963-6 MS	MW4I	108	107	105	104
480-232963-6 MSD	MW4I	106	105	108	106
480-232963-7	MW5S	124 S1+	108	114	102
480-232963-8	MW5I	110	103	109	100
480-232963-9	MW6	109	104	109	103
480-232963-10	MW7	109	103	108	100
480-232963-11	MW8	112	106	105	103
480-232963-12	MW11	110	104	107	98
480-232963-13	MW12S	113	103	111	103
480-232963-14	MW1C	110	102	107	102
480-232963-15	MW1D	110	107	108	103
480-232963-16	MW1E	109	103	111	97
480-232963-17	TB	107	105	104	102
LCS 480-758407/6	Lab Control Sample	109	108	108	103
LCS 480-758501/6	Lab Control Sample	114	105	111	101
MB 480-758407/8	Method Blank	113	108	111	102
MB 480-758501/8	Method Blank	105	101	102	97

Surrogate Legend

DCA = 1,2-Dichloroethane-d4 (Surr)
BFB = 4-Bromofluorobenzene (Surr)
DBFM = Dibromofluoromethane (Surr)
TOL = Toluene-d8 (Surr)

Isotope Dilution Summary

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A

Matrix: Water **Prep Type: Total/NA**

		Percent Isotope Dilution Recovery (Acceptance Limits)							
Lab Sample ID	Client Sample ID	PFDoA (10-130)	PFTDA (10-130)	NMFM (10-130)	NEFM (10-130)				
480-232963-3 - RE	MW3S	89.1	80.9	85.0	56.5				

Surrogate Legend

PFDoA = 13C2 PFDoA
PFTDA = 13C2 PFTeDA
NMFM = d7-N-MeFOSE-M
NEFM = d9-N-EtFOSE-M

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A

Matrix: Water **Prep Type: Total/NA**

		Percent Isotope Dilution Recovery (Acceptance Limits)							
Lab Sample ID	Client Sample ID	M242FTS (40-200)	M262FTS (40-200)	M282FTS (40-300)	PFDoA (10-130)	PFTDA (10-130)	HFPODA (40-130)	C3PFBS (40-135)	C3PFHS (40-130)
480-232963-3	MW3S	228 *5+	113	92.4	22.2		84.7	108	92.9
480-232963-4	MW3I	108	99.2	94.6	57.2	38.1	91.0	116	96.3
480-232963-5	MW4S	191	139	143	81.8	61.9	82.9	117	99.9
480-232963-6	MW4I	113	92.9	97.7	80.5	66.6	85.9	110	92.8
480-232963-7	MW5S	219 *5+	190	125	80.5	66.6	95.1	103	96.3
480-232963-8	MW5I	111	92.6	81.3	87.4	66.0	108	103	96.6
LCS 410-707351/17-A	Lab Control Sample	112	121	126	92.9	85.7	115	108	105
LCS 410-708885/16-A	Lab Control Sample	89.7	82.4	74.5	89.9	77.4	108	94.6	91.5
LCS 410-710704/6-A	Lab Control Sample	111	93.4	110	93.9	79.6	91.7	110	95.4
LCSD 410-707351/18-A	Lab Control Sample Dup	114	120	124	97.5	90.1	113	107	107
LCSD 410-708885/17-A	Lab Control Sample Dup	89.8	84.8	81.1	96.3	82.1	107	99.8	96.3
LCSD 410-710704/7-A	Lab Control Sample Dup	109	94.3	103	87.9	63.6	88.6	100	89.7
LLCS 410-708885/18-A	Lab Control Sample	89.8	85.9	78.6	98.4	81.4	97.0	97.5	95.9
LLCS 410-710704/8-A	Lab Control Sample	112	101	113	89.1	73.8	84.6	108	94.4
MB 410-707351/16-A	Method Blank	123	134	141	97.7	86.4	111	106	105
MB 410-708885/15-A	Method Blank	86.4	86.3	77.6	93.1	74.6	101	99.4	96.6
MB 410-710704/5-A	Method Blank	107	98.6	105	89.7	67.0	89.7	103	87.4

		Percent Isotope Dilution Recovery (Acceptance Limits)							
Lab Sample ID	Client Sample ID	PFBA (5-130)	C4PFHA (40-130)	13C5PHA (40-130)	PFPeA (40-130)	C6PFDA (40-130)	13C7PUA (30-130)	PFOSA (40-130)	C8PFOA (40-130)
480-232963-3	MW3S	76.7	94.9	97.3	65.6	84.8	64.1	79.3	96.2
480-232963-4	MW3I	105	94.7	106	108	88.2	73.1	84.1	99.1
480-232963-5	MW4S	109	91.1	97.5	99.2	96.6	85.7	94.4	102
480-232963-6	MW4I	99.7	86.4	99.5	102	84.0	78.0	80.0	88.8
480-232963-7	MW5S	103	100	97.9	86.5	95.1	83.1	93.6	98.2
480-232963-8	MW5I	107	102	96.2	100	104	90.2	87.5	101
LCS 410-707351/17-A	Lab Control Sample	110	106	104	103	99.0	86.9	106	102
LCS 410-708885/16-A	Lab Control Sample	101	95.8	96.5	99.4	92.2	92.5	81.7	93.9
LCS 410-710704/6-A	Lab Control Sample	103	93.7	96.2	102	96.0	89.6	88.0	93.0
LCSD 410-707351/18-A	Lab Control Sample Dup	108	112	106	108	101	90.8	104	110
LCSD 410-708885/17-A	Lab Control Sample Dup	104	102	98.5	103	98.3	92.3	83.7	97.5
LCSD 410-710704/7-A	Lab Control Sample Dup	97.7	92.3	100	97.4	90.7	86.1	86.6	88.4
LLCS 410-708885/18-A	Lab Control Sample	98.0	94.9	89.3	94.2	97.1	94.8	77.2	90.4
LLCS 410-710704/8-A	Lab Control Sample	101	93.5	100	101	91.0	90.9	87.3	90.3
MB 410-707351/16-A	Method Blank	114	108	108	104	101	90.6	111	112
MB 410-708885/15-A	Method Blank	104	102	98.5	100	98.2	92.2	83.8	94.1

Isotope Dilution Summary

Client: GEI Consultants Inc
Job ID: 480-232963-1
Project/Site: Unifirst – Liverpool, NY

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Matrix: Water **Prep Type: Total/NA**

		Percent Isotope Dilution Recovery (Acceptance Limits)							
Lab Sample ID	Client Sample ID	PFBA (5-130)	C4PFHA (40-130)	13C5PHA (40-130)	PFPeA (40-130)	C6PFDA (40-130)	13C7PUA (30-130)	PFOSA (40-130)	C8PFOA (40-130)
MB 410-710704/5-A	Method Blank	102	96.9	99.0	101	95.0	89.0	90.8	92.9

		Percent Isotope Dilution Recovery (Acceptance Limits)							
Lab Sample ID	Client Sample ID	C8PFOS (40-130)	C9PFNA (40-130)	d3NMFOS (40-170)	d3NMFSA (10-130)	d5NEFOS (25-135)	d5NPFSA (10-130)	NMFM (10-130)	NEFM (10-130)
480-232963-3	MW3S	90.1	97.7	58.0	59.0	33.1	26.5		
480-232963-4	MW3I	92.5	100	68.6	62.2	54.2	52.7	23.0	15.4
480-232963-5	MW4S	104	100	112	81.1	87.5	71.7	57.3	44.0
480-232963-6	MW4I	88.8	93.4	77.1	69.3	66.6	65.7	49.2	38.6
480-232963-7	MW5S	94.5	99.3	89.9	73.0	80.5	63.1	51.8	43.6
480-232963-8	MW5I	94.0	95.8	81.4	68.6	73.0	63.1	21.5	14.5
LCS 410-707351/17-A	Lab Control Sample	96.9	103	102	89.7	96.3	83.0	99.7	71.9
LCS 410-708885/16-A	Lab Control Sample	91.7	88.0	76.0	70.7	70.2	64.8	61.6	54.0
LCS 410-710704/6-A	Lab Control Sample	98.1	103	88.0	78.5	78.6	78.4	72.7	58.2
LCSD 410-707351/18-A	Lab Control Sample Dup	104	103	102	93.0	97.0	86.9	113	105
LCSD 410-708885/17-A	Lab Control Sample Dup	97.5	100	77.4	70.0	73.9	64.9	59.4	52.1
LCSD 410-710704/7-A	Lab Control Sample Dup	95.5	93.5	87.5	72.3	76.6	71.0	61.1	55.0
LLCS 410-708885/18-A	Lab Control Sample	93.8	91.9	80.3	66.5	78.0	63.7	59.3	54.7
LLCS 410-710704/8-A	Lab Control Sample	99.2	95.0	98.0	78.3	83.9	77.3	67.4	55.8
MB 410-707351/16-A	Method Blank	102	108	105	87.5	99.9	79.2	108	89.6
MB 410-708885/15-A	Method Blank	95.6	93.6	81.4	74.1	77.1	68.1	50.9	43.2
MB 410-710704/5-A	Method Blank	96.1	92.7	88.9	76.5	79.1	73.8	56.6	43.1

Surrogate Legend

M242FTS = 13C2 4:2 FTS
M262FTS = 13C2 6:2 FTS
M282FTS = 13C2 8:2 FTS
PFDoA = 13C2 PFDoA
PFTDA = 13C2 PFTeDA
HFPODA = 13C3 HFPO-DA
C3PFBS = 13C3 PFBS
C3PFHS = 13C3 PFHxS
PFBA = 13C4 PFBA
C4PFHA = 13C4 PFHpA
13C5PHA = 13C5 PFHxA
PFPeA = 13C5 PFPeA
C6PFDA = 13C6 PFDA
13C7PUA = 13C7 PFUnA
PFOSA = 13C8 FOSA
C8PFOA = 13C8 PFOA
C8PFOS = 13C8 PFOS
C9PFNA = 13C9 PFNA
d3NMFOS = d3-NMeFOSAA
d3NMFSA = d3-NMePFOSA
d5NEFOS = d5-NEtFOSAA
d5NPFSA = d5-NEtPFOSA
NMFM = d7-N-MeFOSE-M
NEFM = d9-N-EtFOSE-M

Method: 8260C - Volatile Organic Compounds by GC/MS

Lab Sample ID: MB 480-758407/8
Matrix: Water
Analysis Batch: 758407

Client Sample ID: Method Blank
Prep Type: Total/NA

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		1.0	0.82	ug/L			10/01/25 00:55	1
1,1,2,2-Tetrachloroethane	ND		1.0	0.21	ug/L			10/01/25 00:55	1
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1.0	0.31	ug/L			10/01/25 00:55	1
1,1,2-Trichloroethane	ND		1.0	0.23	ug/L			10/01/25 00:55	1
1,1-Dichloroethane	ND		1.0	0.38	ug/L			10/01/25 00:55	1
1,1-Dichloroethene	ND		1.0	0.29	ug/L			10/01/25 00:55	1
1,2,4-Trichlorobenzene	ND		1.0	0.41	ug/L			10/01/25 00:55	1
1,2-Dibromo-3-Chloropropane	ND		1.0	0.39	ug/L			10/01/25 00:55	1
1,2-Dibromoethane	ND		1.0	0.73	ug/L			10/01/25 00:55	1
1,2-Dichlorobenzene	ND		1.0	0.79	ug/L			10/01/25 00:55	1
1,2-Dichloroethane	ND		1.0	0.21	ug/L			10/01/25 00:55	1
1,2-Dichloropropane	ND		1.0	0.72	ug/L			10/01/25 00:55	1
1,3-Dichlorobenzene	ND		1.0	0.78	ug/L			10/01/25 00:55	1
1,4-Dichlorobenzene	ND		1.0	0.84	ug/L			10/01/25 00:55	1
2-Butanone (MEK)	ND		10	1.3	ug/L			10/01/25 00:55	1
2-Hexanone	ND		5.0	1.2	ug/L			10/01/25 00:55	1
4-Methyl-2-pentanone (MIBK)	ND		5.0	2.1	ug/L			10/01/25 00:55	1
Acetone	ND		10	3.0	ug/L			10/01/25 00:55	1
Benzene	ND		1.0	0.41	ug/L			10/01/25 00:55	1
Bromodichloromethane	ND		1.0	0.39	ug/L			10/01/25 00:55	1
Bromoform	ND		1.0	0.26	ug/L			10/01/25 00:55	1
Bromomethane	ND		1.0	0.69	ug/L			10/01/25 00:55	1
Carbon disulfide	ND		1.0	0.19	ug/L			10/01/25 00:55	1
Carbon tetrachloride	ND		1.0	0.27	ug/L			10/01/25 00:55	1
Chlorobenzene	ND		1.0	0.75	ug/L			10/01/25 00:55	1
Chloroethane	ND		1.0	0.32	ug/L			10/01/25 00:55	1
Chloroform	ND		1.0	0.34	ug/L			10/01/25 00:55	1
Chloromethane	ND		1.0	0.35	ug/L			10/01/25 00:55	1
cis-1,2-Dichloroethene	ND		1.0	0.81	ug/L			10/01/25 00:55	1
cis-1,3-Dichloropropene	ND		1.0	0.36	ug/L			10/01/25 00:55	1
Cyclohexane	ND		1.0	0.18	ug/L			10/01/25 00:55	1
Dibromochloromethane	ND		1.0	0.32	ug/L			10/01/25 00:55	1
Dichlorodifluoromethane	ND		1.0	0.68	ug/L			10/01/25 00:55	1
Ethylbenzene	ND		1.0	0.74	ug/L			10/01/25 00:55	1
Isopropylbenzene	ND		1.0	0.79	ug/L			10/01/25 00:55	1
Methyl acetate	ND		2.5	1.3	ug/L			10/01/25 00:55	1
Methyl tert-butyl ether	ND		1.0	0.16	ug/L			10/01/25 00:55	1
Methylcyclohexane	ND		1.0	0.16	ug/L			10/01/25 00:55	1
Methylene Chloride	ND		1.0	0.44	ug/L			10/01/25 00:55	1
Styrene	ND		1.0	0.73	ug/L			10/01/25 00:55	1
Tetrachloroethene	ND		1.0	0.36	ug/L			10/01/25 00:55	1
Toluene	ND		1.0	0.51	ug/L			10/01/25 00:55	1
trans-1,2-Dichloroethene	ND		1.0	0.90	ug/L			10/01/25 00:55	1
trans-1,3-Dichloropropene	ND		1.0	0.37	ug/L			10/01/25 00:55	1
Trichloroethene	ND		1.0	0.46	ug/L			10/01/25 00:55	1
Trichlorofluoromethane	ND		1.0	0.88	ug/L			10/01/25 00:55	1
Vinyl chloride	ND		1.0	0.90	ug/L			10/01/25 00:55	1
Xylenes, Total	ND		2.0	0.66	ug/L			10/01/25 00:55	1

Method: 8260C - Volatile Organic Compounds by GC/MS (Continued)

Lab Sample ID: MB 480-758407/8
Matrix: Water
Analysis Batch: 758407

Client Sample ID: Method Blank
Prep Type: Total/NA

Surrogate	MB %Recovery	MB Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	113		77 - 120		10/01/25 00:55	1
4-Bromofluorobenzene (Surr)	108		73 - 120		10/01/25 00:55	1
Dibromofluoromethane (Surr)	111		75 - 123		10/01/25 00:55	1
Toluene-d8 (Surr)	102		80 - 120		10/01/25 00:55	1

Lab Sample ID: LCS 480-758407/6
Matrix: Water
Analysis Batch: 758407

Client Sample ID: Lab Control Sample
Prep Type: Total/NA

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
1,1,1-Trichloroethane	25.0	25.1		ug/L		101	73 - 126
1,1,2,2-Tetrachloroethane	25.0	22.0		ug/L		88	76 - 120
1,1,2-Trichloro-1,2,2-trifluoroethane	25.0	25.9		ug/L		104	61 - 148
1,1,2-Trichloroethane	25.0	23.2		ug/L		93	76 - 122
1,1-Dichloroethane	25.0	23.3		ug/L		93	77 - 120
1,1-Dichloroethene	25.0	23.5		ug/L		94	66 - 127
1,2,4-Trichlorobenzene	25.0	21.9		ug/L		88	79 - 122
1,2-Dibromo-3-Chloropropane	25.0	20.2		ug/L		81	56 - 134
1,2-Dibromoethane	25.0	23.2		ug/L		93	77 - 120
1,2-Dichlorobenzene	25.0	22.5		ug/L		90	80 - 124
1,2-Dichloroethane	25.0	24.9		ug/L		100	75 - 120
1,2-Dichloropropane	25.0	23.4		ug/L		93	76 - 120
1,3-Dichlorobenzene	25.0	22.2		ug/L		89	77 - 120
1,4-Dichlorobenzene	25.0	21.6		ug/L		86	80 - 120
2-Butanone (MEK)	125	122		ug/L		98	57 - 140
2-Hexanone	125	127		ug/L		101	65 - 127
4-Methyl-2-pentanone (MIBK)	125	115		ug/L		92	71 - 125
Acetone	125	132		ug/L		106	56 - 142
Benzene	25.0	23.5		ug/L		94	71 - 124
Bromodichloromethane	25.0	25.5		ug/L		102	80 - 122
Bromoform	25.0	30.9		ug/L		123	61 - 132
Bromomethane	25.0	27.6		ug/L		110	55 - 144
Carbon disulfide	25.0	23.6		ug/L		94	59 - 134
Carbon tetrachloride	25.0	26.1		ug/L		104	72 - 134
Chlorobenzene	25.0	23.1		ug/L		92	80 - 120
Chloroethane	25.0	27.3		ug/L		109	69 - 136
Chloroform	25.0	25.0		ug/L		100	73 - 127
Chloromethane	25.0	26.4		ug/L		105	68 - 124
cis-1,2-Dichloroethene	25.0	23.5		ug/L		94	74 - 124
cis-1,3-Dichloropropene	25.0	21.6		ug/L		86	74 - 124
Cyclohexane	25.0	23.1		ug/L		93	59 - 135
Dibromochloromethane	25.0	27.0		ug/L		108	75 - 125
Dichlorodifluoromethane	25.0	25.8		ug/L		103	59 - 135
Ethylbenzene	25.0	23.5		ug/L		94	77 - 123
Isopropylbenzene	25.0	21.8		ug/L		87	77 - 122
Methyl acetate	50.0	46.0		ug/L		92	74 - 133
Methyl tert-butyl ether	25.0	22.3		ug/L		89	77 - 120
Methylcyclohexane	25.0	24.0		ug/L		96	68 - 134

Method: 8260C - Volatile Organic Compounds by GC/MS (Continued)

Lab Sample ID: LCS 480-758407/6
Matrix: Water
Analysis Batch: 758407

Client Sample ID: Lab Control Sample
Prep Type: Total/NA

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
Methylene Chloride	25.0	25.0		ug/L		100	75 - 124
Styrene	25.0	22.9		ug/L		92	80 - 120
Tetrachloroethene	25.0	24.1		ug/L		96	74 - 122
Toluene	25.0	23.1		ug/L		92	80 - 122
trans-1,2-Dichloroethene	25.0	24.4		ug/L		98	73 - 127
trans-1,3-Dichloropropene	25.0	21.6		ug/L		86	80 - 120
Trichloroethene	25.0	24.4		ug/L		98	74 - 123
Trichlorofluoromethane	25.0	31.0		ug/L		124	62 - 150
Vinyl chloride	25.0	26.7		ug/L		107	65 - 133
Xylenes, Total	50.0	46.9		ug/L		94	76 - 122

Surrogate	*LCS %Recovery*	*LCS Qualifier*	*Limits*
1,2-Dichloroethane-d4 (Surr)	*109*		*77 - 120*
4-Bromofluorobenzene (Surr)	*108*		*73 - 120*
Dibromofluoromethane (Surr)	*108*		*75 - 123*
Toluene-d8 (Surr)	*103*		*80 - 120*

Lab Sample ID: 480-232963-6 MS
Matrix: Water
Analysis Batch: 758407

Client Sample ID: MW4I
Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MS Result	MS Qualifier	Unit	D	%Rec	%Rec Limits
1,1,1-Trichloroethane	ND		1000	1040		ug/L		104	73 - 126
1,1,2,2-Tetrachloroethane	ND		1000	901		ug/L		90	76 - 120
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1000	1080		ug/L		108	61 - 148
1,1,2-Trichloroethane	ND		1000	951		ug/L		95	76 - 122
1,1-Dichloroethane	ND		1000	929		ug/L		93	77 - 120
1,1-Dichloroethene	36	J	1000	983		ug/L		95	66 - 127
1,2,4-Trichlorobenzene	ND		1000	848		ug/L		85	79 - 122
1,2-Dibromo-3-Chloropropane	ND		1000	884		ug/L		88	56 - 134
1,2-Dibromoethane	ND		1000	924		ug/L		92	77 - 120
1,2-Dichlorobenzene	ND		1000	916		ug/L		92	80 - 124
1,2-Dichloroethane	ND		1000	992		ug/L		99	75 - 120
1,2-Dichloropropane	ND		1000	927		ug/L		93	76 - 120
1,3-Dichlorobenzene	ND		1000	894		ug/L		89	77 - 120
1,4-Dichlorobenzene	ND		1000	879		ug/L		88	78 - 124
2-Butanone (MEK)	ND		5000	4890		ug/L		98	57 - 140
2-Hexanone	ND		5000	5110		ug/L		102	65 - 127
4-Methyl-2-pentanone (MIBK)	ND		5000	4750		ug/L		95	71 - 125
Acetone	ND		5000	4860		ug/L		97	56 - 142
Benzene	ND		1000	952		ug/L		95	71 - 124
Bromodichloromethane	ND		1000	1040		ug/L		104	80 - 122
Bromoform	ND		1000	1200		ug/L		120	61 - 132
Bromomethane	ND		1000	1120		ug/L		112	55 - 144
Carbon disulfide	ND		1000	987		ug/L		99	59 - 134
Carbon tetrachloride	ND		1000	1070		ug/L		107	72 - 134
Chlorobenzene	ND		1000	924		ug/L		92	80 - 120
Chloroethane	ND		1000	1140		ug/L		114	69 - 136

Method: 8260C - Volatile Organic Compounds by GC/MS (Continued)

Lab Sample ID: 480-232963-6 MS
Matrix: Water
Analysis Batch: 758407

Client Sample ID: MW4I
Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MS Result	MS Qualifier	Unit	D	%Rec	%Rec Limits
Chloroform	ND		1000	995		ug/L		100	73 - 127
Chloromethane	ND		1000	1090		ug/L		109	68 - 124
cis-1,2-Dichloroethene	3800	F1	1000	3500	F1	ug/L		-30	74 - 124
cis-1,3-Dichloropropene	ND		1000	846		ug/L		85	74 - 124
Cyclohexane	ND		1000	967		ug/L		97	59 - 135
Dibromochloromethane	ND		1000	1080		ug/L		108	75 - 125
Dichlorodifluoromethane	ND		1000	1080		ug/L		108	59 - 135
Ethylbenzene	ND		1000	946		ug/L		95	77 - 123
Isopropylbenzene	ND		1000	903		ug/L		90	77 - 122
Methyl acetate	ND		2000	1830		ug/L		92	74 - 133
Methyl tert-butyl ether	ND		1000	896		ug/L		90	77 - 120
Methylcyclohexane	ND		1000	972		ug/L		97	68 - 134
Methylene Chloride	ND		1000	1010		ug/L		101	75 - 124
Styrene	ND		1000	929		ug/L		93	80 - 120
Tetrachloroethene	20	J	1000	1000		ug/L		98	74 - 122
Toluene	ND		1000	927		ug/L		93	80 - 122
trans-1,2-Dichloroethene	ND		1000	998		ug/L		100	73 - 127
trans-1,3-Dichloropropene	ND		1000	803		ug/L		80	80 - 120
Trichloroethene	ND		1000	987		ug/L		99	74 - 123
Trichlorofluoromethane	ND		1000	1300		ug/L		130	62 - 150
Vinyl chloride	550		1000	1440		ug/L		90	65 - 133
Xylenes, Total	ND		2000	1880		ug/L		94	76 - 122

Surrogate	*MS %Recovery*	*MS Qualifier*	*Limits*
1,2-Dichloroethane-d4 (Surr)	*108*		*77 - 120*
4-Bromofluorobenzene (Surr)	*107*		*73 - 120*
Dibromofluoromethane (Surr)	*105*		*75 - 123*
Toluene-d8 (Surr)	*104*		*80 - 120*

Lab Sample ID: 480-232963-6 MSD
Matrix: Water
Analysis Batch: 758407

Client Sample ID: MW4I
Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MSD Result	MSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
1,1,1-Trichloroethane	ND		1000	991		ug/L		99	73 - 126	4	15
1,1,2,2-Tetrachloroethane	ND		1000	940		ug/L		94	76 - 120	4	15
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1000	907		ug/L		91	61 - 148	18	20
1,1,2-Trichloroethane	ND		1000	950		ug/L		95	76 - 122	0	15
1,1-Dichloroethane	ND		1000	911		ug/L		91	77 - 120	2	20
1,1-Dichloroethene	36	J	1000	958		ug/L		92	66 - 127	3	16
1,2,4-Trichlorobenzene	ND		1000	881		ug/L		88	79 - 122	4	20
1,2-Dibromo-3-Chloropropane	ND		1000	895		ug/L		90	56 - 134	1	15
1,2-Dibromoethane	ND		1000	901		ug/L		90	77 - 120	3	15
1,2-Dichlorobenzene	ND		1000	923		ug/L		92	80 - 124	1	20
1,2-Dichloroethane	ND		1000	961		ug/L		96	75 - 120	3	20
1,2-Dichloropropane	ND		1000	911		ug/L		91	76 - 120	2	20
1,3-Dichlorobenzene	ND		1000	920		ug/L		92	77 - 120	3	20
1,4-Dichlorobenzene	ND		1000	891		ug/L		89	78 - 124	1	20

Method: 8260C - Volatile Organic Compounds by GC/MS (Continued)

Lab Sample ID: 480-232963-6 MSD
Matrix: Water
Analysis Batch: 758407

Client Sample ID: MW4I
Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MSD Result	MSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
2-Butanone (MEK)	ND		5000	4790		ug/L		96	57 - 140	2	20
2-Hexanone	ND		5000	5120		ug/L		102	65 - 127	0	15
4-Methyl-2-pentanone (MIBK)	ND		5000	4630		ug/L		93	71 - 125	3	35
Acetone	ND		5000	4740		ug/L		95	56 - 142	2	15
Benzene	ND		1000	912		ug/L		91	71 - 124	4	13
Bromodichloromethane	ND		1000	1020		ug/L		102	80 - 122	1	15
Bromoform	ND		1000	1190		ug/L		119	61 - 132	0	15
Bromomethane	ND		1000	1100		ug/L		110	55 - 144	2	15
Carbon disulfide	ND		1000	932		ug/L		93	59 - 134	6	15
Carbon tetrachloride	ND		1000	981		ug/L		98	72 - 134	8	15
Chlorobenzene	ND		1000	904		ug/L		90	80 - 120	2	25
Chloroethane	ND		1000	1110		ug/L		111	69 - 136	3	15
Chloroform	ND		1000	993		ug/L		99	73 - 127	0	20
Chloromethane	ND		1000	1030		ug/L		103	68 - 124	6	15
cis-1,2-Dichloroethene	3800	F1	1000	3330	F1	ug/L		-46	74 - 124	5	15
cis-1,3-Dichloropropene	ND		1000	803		ug/L		80	74 - 124	5	15
Cyclohexane	ND		1000	931		ug/L		93	59 - 135	4	20
Dibromochloromethane	ND		1000	1090		ug/L		109	75 - 125	1	15
Dichlorodifluoromethane	ND		1000	1050		ug/L		105	59 - 135	2	20
Ethylbenzene	ND		1000	933		ug/L		93	77 - 123	1	15
Isopropylbenzene	ND		1000	926		ug/L		93	77 - 122	2	20
Methyl acetate	ND		2000	1790		ug/L		90	74 - 133	2	20
Methyl tert-butyl ether	ND		1000	879		ug/L		88	77 - 120	2	37
Methylcyclohexane	ND		1000	937		ug/L		94	68 - 134	4	20
Methylene Chloride	ND		1000	975		ug/L		98	75 - 124	4	15
Styrene	ND		1000	909		ug/L		91	80 - 120	2	20
Tetrachloroethene	20	J	1000	947		ug/L		93	74 - 122	6	20
Toluene	ND		1000	921		ug/L		92	80 - 122	1	15
trans-1,2-Dichloroethene	ND		1000	926		ug/L		93	73 - 127	7	20
trans-1,3-Dichloropropene	ND		1000	795		ug/L		80	80 - 120	1	15
Trichloroethene	ND		1000	947		ug/L		95	74 - 123	4	16
Trichlorofluoromethane	ND		1000	1280		ug/L		128	62 - 150	1	20
Vinyl chloride	550		1000	1400		ug/L		85	65 - 133	3	15
Xylenes, Total	ND		2000	1840		ug/L		92	76 - 122	2	16

Surrogate	*MSD %Recovery*	*MSD Qualifier*	*Limits*
1,2-Dichloroethane-d4 (Surr)	*106*		*77 - 120*
4-Bromofluorobenzene (Surr)	*105*		*73 - 120*
Dibromofluoromethane (Surr)	*108*		*75 - 123*
Toluene-d8 (Surr)	*106*		*80 - 120*

Lab Sample ID: MB 480-758501/8
Matrix: Water
Analysis Batch: 758501

Client Sample ID: Method Blank
Prep Type: Total/NA

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1-Trichloroethane	ND		1.0	0.82	ug/L			10/01/25 12:44	1
1,1,2,2-Tetrachloroethane	ND		1.0	0.21	ug/L			10/01/25 12:44	1

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Method: 8260C - Volatile Organic Compounds by GC/MS (Continued)

Lab Sample ID: MB 480-758501/8
Matrix: Water
Analysis Batch: 758501

Client Sample ID: Method Blank
Prep Type: Total/NA

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		1.0	0.31	ug/L			10/01/25 12:44	1
1,1,2-Trichloroethane	ND		1.0	0.23	ug/L			10/01/25 12:44	1
1,1-Dichloroethane	ND		1.0	0.38	ug/L			10/01/25 12:44	1
1,1-Dichloroethene	ND		1.0	0.29	ug/L			10/01/25 12:44	1
1,2,4-Trichlorobenzene	ND		1.0	0.41	ug/L			10/01/25 12:44	1
1,2-Dibromo-3-Chloropropane	ND		1.0	0.39	ug/L			10/01/25 12:44	1
1,2-Dibromoethane	ND		1.0	0.73	ug/L			10/01/25 12:44	1
1,2-Dichlorobenzene	ND		1.0	0.79	ug/L			10/01/25 12:44	1
1,2-Dichloroethane	ND		1.0	0.21	ug/L			10/01/25 12:44	1
1,2-Dichloropropane	ND		1.0	0.72	ug/L			10/01/25 12:44	1
1,3-Dichlorobenzene	ND		1.0	0.78	ug/L			10/01/25 12:44	1
1,4-Dichlorobenzene	ND		1.0	0.84	ug/L			10/01/25 12:44	1
2-Butanone (MEK)	ND		10	1.3	ug/L			10/01/25 12:44	1
2-Hexanone	ND		5.0	1.2	ug/L			10/01/25 12:44	1
4-Methyl-2-pentanone (MIBK)	ND		5.0	2.1	ug/L			10/01/25 12:44	1
Acetone	ND		10	3.0	ug/L			10/01/25 12:44	1
Benzene	ND		1.0	0.41	ug/L			10/01/25 12:44	1
Bromodichloromethane	ND		1.0	0.39	ug/L			10/01/25 12:44	1
Bromoform	ND		1.0	0.26	ug/L			10/01/25 12:44	1
Bromomethane	ND		1.0	0.69	ug/L			10/01/25 12:44	1
Carbon disulfide	ND		1.0	0.19	ug/L			10/01/25 12:44	1
Carbon tetrachloride	ND		1.0	0.27	ug/L			10/01/25 12:44	1
Chlorobenzene	ND		1.0	0.75	ug/L			10/01/25 12:44	1
Chloroethane	ND		1.0	0.32	ug/L			10/01/25 12:44	1
Chloroform	ND		1.0	0.34	ug/L			10/01/25 12:44	1
Chloromethane	ND		1.0	0.35	ug/L			10/01/25 12:44	1
cis-1,2-Dichloroethene	ND		1.0	0.81	ug/L			10/01/25 12:44	1
cis-1,3-Dichloropropene	ND		1.0	0.36	ug/L			10/01/25 12:44	1
Cyclohexane	ND		1.0	0.18	ug/L			10/01/25 12:44	1
Dibromochloromethane	ND		1.0	0.32	ug/L			10/01/25 12:44	1
Dichlorodifluoromethane	ND		1.0	0.68	ug/L			10/01/25 12:44	1
Ethylbenzene	ND		1.0	0.74	ug/L			10/01/25 12:44	1
Isopropylbenzene	ND		1.0	0.79	ug/L			10/01/25 12:44	1
Methyl acetate	ND		2.5	1.3	ug/L			10/01/25 12:44	1
Methyl tert-butyl ether	ND		1.0	0.16	ug/L			10/01/25 12:44	1
Methylcyclohexane	ND		1.0	0.16	ug/L			10/01/25 12:44	1
Methylene Chloride	ND		1.0	0.44	ug/L			10/01/25 12:44	1
Styrene	ND		1.0	0.73	ug/L			10/01/25 12:44	1
Tetrachloroethene	ND		1.0	0.36	ug/L			10/01/25 12:44	1
Toluene	ND		1.0	0.51	ug/L			10/01/25 12:44	1
trans-1,2-Dichloroethene	ND		1.0	0.90	ug/L			10/01/25 12:44	1
trans-1,3-Dichloropropene	ND		1.0	0.37	ug/L			10/01/25 12:44	1
Trichloroethene	ND		1.0	0.46	ug/L			10/01/25 12:44	1
Trichlorofluoromethane	ND		1.0	0.88	ug/L			10/01/25 12:44	1
Vinyl chloride	ND		1.0	0.90	ug/L			10/01/25 12:44	1
Xylenes, Total	ND		2.0	0.66	ug/L			10/01/25 12:44	1

Method: 8260C - Volatile Organic Compounds by GC/MS (Continued)

Lab Sample ID: MB 480-758501/8
Matrix: Water
Analysis Batch: 758501

Client Sample ID: Method Blank
Prep Type: Total/NA

Surrogate	MB %Recovery	MB Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	105		77 - 120		10/01/25 12:44	1
4-Bromofluorobenzene (Surr)	101		73 - 120		10/01/25 12:44	1
Dibromofluoromethane (Surr)	102		75 - 123		10/01/25 12:44	1
Toluene-d8 (Surr)	97		80 - 120		10/01/25 12:44	1

Lab Sample ID: LCS 480-758501/6
Matrix: Water
Analysis Batch: 758501

Client Sample ID: Lab Control Sample
Prep Type: Total/NA

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
1,1,1-Trichloroethane	25.0	25.9		ug/L		104	73 - 126
1,1,2,2-Tetrachloroethane	25.0	21.9		ug/L		88	76 - 120
1,1,2-Trichloro-1,2,2-trifluoroethane	25.0	26.3		ug/L		105	61 - 148
1,1,2-Trichloroethane	25.0	21.4		ug/L		86	76 - 122
1,1-Dichloroethane	25.0	23.5		ug/L		94	77 - 120
1,1-Dichloroethene	25.0	23.7		ug/L		95	66 - 127
1,2,4-Trichlorobenzene	25.0	21.3		ug/L		85	79 - 122
1,2-Dibromo-3-Chloropropane	25.0	20.8		ug/L		83	56 - 134
1,2-Dibromoethane	25.0	22.3		ug/L		89	77 - 120
1,2-Dichlorobenzene	25.0	22.6		ug/L		90	80 - 124
1,2-Dichloroethane	25.0	24.6		ug/L		98	75 - 120
1,2-Dichloropropane	25.0	23.4		ug/L		94	76 - 120
1,3-Dichlorobenzene	25.0	22.1		ug/L		88	77 - 120
1,4-Dichlorobenzene	25.0	21.8		ug/L		87	80 - 120
2-Butanone (MEK)	125	124		ug/L		99	57 - 140
2-Hexanone	125	124		ug/L		99	65 - 127
4-Methyl-2-pentanone (MIBK)	125	111		ug/L		89	71 - 125
Acetone	125	143		ug/L		115	56 - 142
Benzene	25.0	23.5		ug/L		94	71 - 124
Bromodichloromethane	25.0	26.8		ug/L		107	80 - 122
Bromoform	25.0	29.2		ug/L		117	61 - 132
Bromomethane	25.0	28.4		ug/L		113	55 - 144
Carbon disulfide	25.0	23.8		ug/L		95	59 - 134
Carbon tetrachloride	25.0	26.1		ug/L		104	72 - 134
Chlorobenzene	25.0	22.6		ug/L		90	80 - 120
Chloroethane	25.0	27.1		ug/L		108	69 - 136
Chloroform	25.0	24.7		ug/L		99	73 - 127
Chloromethane	25.0	26.7		ug/L		107	68 - 124
cis-1,2-Dichloroethene	25.0	23.6		ug/L		95	74 - 124
cis-1,3-Dichloropropene	25.0	22.5		ug/L		90	74 - 124
Cyclohexane	25.0	23.8		ug/L		95	59 - 135
Dibromochloromethane	25.0	26.9		ug/L		108	75 - 125
Dichlorodifluoromethane	25.0	26.3		ug/L		105	59 - 135
Ethylbenzene	25.0	22.8		ug/L		91	77 - 123
Isopropylbenzene	25.0	21.9		ug/L		88	77 - 122
Methyl acetate	50.0	45.2		ug/L		90	74 - 133
Methyl tert-butyl ether	25.0	22.3		ug/L		89	77 - 120
Methylcyclohexane	25.0	24.6		ug/L		98	68 - 134

Method: 8260C - Volatile Organic Compounds by GC/MS (Continued)

Lab Sample ID: LCS 480-758501/6
Matrix: Water
Analysis Batch: 758501

Client Sample ID: Lab Control Sample
Prep Type: Total/NA

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
Methylene Chloride	25.0	24.9		ug/L		100	75 - 124
Styrene	25.0	22.1		ug/L		88	80 - 120
Tetrachloroethene	25.0	23.9		ug/L		95	74 - 122
Toluene	25.0	21.8		ug/L		87	80 - 122
trans-1,2-Dichloroethene	25.0	24.1		ug/L		96	73 - 127
trans-1,3-Dichloropropene	25.0	20.5		ug/L		82	80 - 120
Trichloroethene	25.0	24.3		ug/L		97	74 - 123
Trichlorofluoromethane	25.0	32.3		ug/L		129	62 - 150
Vinyl chloride	25.0	27.2		ug/L		109	65 - 133
Xylenes, Total	50.0	44.9		ug/L		90	76 - 122

Surrogate	*LCS %Recovery*	*LCS Qualifier*	*Limits*
1,2-Dichloroethane-d4 (Surr)	*114*		*77 - 120*
4-Bromofluorobenzene (Surr)	*105*		*73 - 120*
Dibromofluoromethane (Surr)	*111*		*75 - 123*
Toluene-d8 (Surr)	*101*		*80 - 120*

Method: RSK-175 - Dissolved Gases (GC)

Lab Sample ID: MB 480-758460/4
Matrix: Water
Analysis Batch: 758460

Client Sample ID: Method Blank
Prep Type: Total/NA

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	ND		4.0	1.0	ug/L			10/01/25 07:12	1
Ethane	ND		7.5	1.5	ug/L			10/01/25 07:12	1
Ethene	ND		7.0	1.5	ug/L			10/01/25 07:12	1

Lab Sample ID: LCS 480-758460/5
Matrix: Water
Analysis Batch: 758460

Client Sample ID: Lab Control Sample
Prep Type: Total/NA

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
Methane	19.5	18.6		ug/L		96	85 - 120
Ethane	36.5	35.1		ug/L		96	79 - 120
Ethene	34.0	32.2		ug/L		95	85 - 120

Lab Sample ID: LCSD 480-758460/6
Matrix: Water
Analysis Batch: 758460

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA

Analyte	Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
Methane	19.5	19.7		ug/L		101	85 - 120	6	50
Ethane	36.5	36.9		ug/L		101	79 - 120	5	50
Ethene	34.0	33.3		ug/L		98	85 - 120	3	50

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A

Lab Sample ID: MB 410-707351/16-A
Matrix: Water
Analysis Batch: 707164

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 707351

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)	ND		4.0	1.0	ng/L		10/01/25 08:35	10/02/25 02:22	1
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)	ND		4.0	1.0	ng/L		10/01/25 08:35	10/02/25 02:22	1
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	ND		4.0	1.0	ng/L		10/01/25 08:35	10/02/25 02:22	1
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	ND		10	2.5	ng/L		10/01/25 08:35	10/02/25 02:22	1
3-Perfluoropentylpropanoic acid (5:3 FTCA)	ND		10	2.8	ng/L		10/01/25 08:35	10/02/25 02:22	1
3-Perfluoropropylpropanoic acid (3:3 FTCA)	ND		4.0	1.0	ng/L		10/01/25 08:35	10/02/25 02:22	1
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	ND		2.0	1.4	ng/L		10/01/25 08:35	10/02/25 02:22	1
N-ethylperfluorooctane sulfonamide (NEtFOSA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	ND		10	4.0	ng/L		10/01/25 08:35	10/02/25 02:22	1
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
N-methylperfluorooctane sulfonamide (NMeFOSA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	ND		10	2.5	ng/L		10/01/25 08:35	10/02/25 02:22	1
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluoro-3-methoxypropanoic acid (PFMPA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluoro-4-methoxybutanoic acid (PFMBA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorobutanesulfonic acid (PFBS)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorobutanoic acid (PFBA)	ND		4.0	1.1	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorodecanesulfonic acid (PFDS)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorodecanoic acid (PFDA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorododecanesulfonic acid (PFDoS)	ND		2.0	0.60	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorododecanoic acid (PFDoA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluoroheptanesulfonic acid (PFHpS)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluoroheptanoic acid (PFHpA)	ND		2.0	0.80	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorohexanesulfonic acid (PFHxS)	ND		2.0	0.80	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorohexanoic acid (PFHxA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorononanesulfonic acid (PFNS)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorononanoic acid (PFNA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: MB 410-707351/16-A
Matrix: Water
Analysis Batch: 707164

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 707351

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanesulfonamide (PFOSA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorooctanesulfonic acid (PFOS)	ND		2.0	1.3	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorooctanoic acid (PFOA)	ND		2.0	0.90	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluoropentanesulfonic acid (PFPeS)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluoropentanoic acid (PFPeA)	ND		2.0	0.60	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorotetradecanoic acid (PFTeDA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluorotridecanoic acid (PFTrDA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1
Perfluoroundecanoic acid (PFUnA)	ND		2.0	0.50	ng/L		10/01/25 08:35	10/02/25 02:22	1

Isotope Dilution	*MB %Recovery*	*MB Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
13C2 4:2 FTS	*123*		*40 - 200*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C2 6:2 FTS	*134*		*40 - 200*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C2 8:2 FTS	*141*		*40 - 300*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C2 PFDoA	*97.7*		*10 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C2 PFTeDA	*86.4*		*10 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C3 HFPO-DA	*111*		*40 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C3 PFBS	*106*		*40 - 135*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C3 PFHxS	*105*		*40 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C4 PFBA	*114*		*5 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C4 PFHpA	*108*		*40 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C5 PFHxA	*108*		*40 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C5 PFPeA	*104*		*40 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C6 PFDA	*101*		*40 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C7 PFUnA	*90.6*		*30 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C8 FOSA	*111*		*40 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C8 PFOA	*112*		*40 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C8 PFOS	*102*		*40 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
13C9 PFNA	*108*		*40 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
d3-NMeFOSAA	*105*		*40 - 170*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
d3-NMePFOSA	*87.5*		*10 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
d5-NEtFOSAA	*99.9*		*25 - 135*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
d5-NEtPFOSA	*79.2*		*10 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
d7-N-MeFOSE-M	*108*		*10 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*
d9-N-EtFOSE-M	*89.6*		*10 - 130*	*10/01/25 08:35*	*10/02/25 02:22*	*1*

Lab Sample ID: LCS 410-707351/17-A
Matrix: Water
Analysis Batch: 707164

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 707351

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)	37.8	31.3		ng/L		83	55 - 160
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)	76.8	66.8		ng/L		87	60 - 150
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)	75.0	67.0		ng/L		89	70 - 145
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	76.2	66.8		ng/L		88	65 - 155

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LCS 410-707351/17-A
Matrix: Water
Analysis Batch: 707164

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 707351

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	200	116		ng/L		58	50 - 145
3-Perfluoropentylpropanoic acid (5:3 FTCA)	200	122	*-	ng/L		61	70 - 135
3-Perfluoropropylpropanoic acid (3:3 FTCA)	80.0	63.0		ng/L		79	65 - 130
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	37.8	30.5		ng/L		81	65 - 145
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	37.4	32.3		ng/L		87	70 - 155
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	30.0	24.4		ng/L		81	70 - 140
N-ethylperfluorooctane sulfonamide (NEtFOSA)	40.0	35.6		ng/L		89	65 - 145
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	200	186		ng/L		93	70 - 135
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	40.0	36.0		ng/L		90	70 - 145
N-methylperfluorooctane sulfonamide (NMeFOSA)	40.0	35.1		ng/L		88	60 - 150
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	200	169		ng/L		85	70 - 145
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	40.0	34.0		ng/L		85	50 - 140
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	40.0	32.7		ng/L		82	50 - 150
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	35.7	32.9		ng/L		92	70 - 140
Perfluoro-3-methoxypropanoic acid (PFMPA)	40.0	37.2		ng/L		93	55 - 140
Perfluoro-4-methoxybutanoic acid (PFMBA)	40.0	33.8		ng/L		84	60 - 150
Perfluorobutanesulfonic acid (PFBS)	35.5	31.1		ng/L		88	60 - 145
Perfluorobutanoic acid (PFBA)	80.0	64.8		ng/L		81	70 - 140
Perfluorodecanesulfonic acid (PFDS)	38.6	31.1		ng/L		81	60 - 145
Perfluorodecanoic acid (PFDA)	40.0	32.1		ng/L		80	70 - 140
Perfluorododecanesulfonic acid (PFDoS)	38.8	28.7		ng/L		74	50 - 145
Perfluorododecanoic acid (PFDoA)	40.0	35.6		ng/L		89	70 - 140
Perfluoroheptanesulfonic acid (PFHpS)	38.2	36.3		ng/L		95	70 - 150
Perfluoroheptanoic acid (PFHpA)	40.0	34.7		ng/L		87	70 - 150
Perfluorohexanesulfonic acid (PFHxS)	36.5	30.8		ng/L		84	65 - 145
Perfluorohexanoic acid (PFHxA)	40.0	33.8		ng/L		84	70 - 145
Perfluorononanesulfonic acid (PFNS)	38.5	33.6		ng/L		87	65 - 145
Perfluorononanoic acid (PFNA)	40.0	32.9		ng/L		82	70 - 150
Perfluorooctanesulfonamide (PFOSA)	40.0	35.4		ng/L		89	70 - 145

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LCS 410-707351/17-A
Matrix: Water
Analysis Batch: 707164

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 707351

Analyte			Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits		
Perfluorooctanesulfonic acid (PFOS)			37.2	31.9		ng/L		86	55 - 150		
Perfluorooctanoic acid (PFOA)			40.0	34.1		ng/L		85	70 - 150		
Perfluoropentanesulfonic acid (PFPeS)			37.6	32.2		ng/L		86	65 - 140		
Perfluoropentanoic acid (PFPeA)			40.0	34.4		ng/L		86	65 - 135		
Perfluorotetradecanoic acid (PFTeDA)			40.0	36.6		ng/L		92	60 - 140		
Perfluorotridecanoic acid (PFTrDA)			40.0	33.6		ng/L		84	65 - 140		
Perfluoroundecanoic acid (PFUnA)			40.0	35.4		ng/L		89	70 - 145		

Isotope Dilution	*LCS %Recovery*	*LCS Qualifier*	*Limits*
13C2 4:2 FTS	*112*		*40 - 200*
13C2 6:2 FTS	*121*		*40 - 200*
13C2 8:2 FTS	*126*		*40 - 300*
13C2 PFDoA	*92.9*		*10 - 130*
13C2 PFTeDA	*85.7*		*10 - 130*
13C3 HFPO-DA	*115*		*40 - 130*
13C3 PFBS	*108*		*40 - 135*
13C3 PFHxS	*105*		*40 - 130*
13C4 PFBA	*110*		*5 - 130*
13C4 PFHpA	*106*		*40 - 130*
13C5 PFHxA	*104*		*40 - 130*
13C5 PFPeA	*103*		*40 - 130*
13C6 PFDA	*99.0*		*40 - 130*
13C7 PFUnA	*86.9*		*30 - 130*
13C8 FOSA	*106*		*40 - 130*
13C8 PFOA	*102*		*40 - 130*
13C8 PFOS	*96.9*		*40 - 130*
13C9 PFNA	*103*		*40 - 130*
d3-NMeFOSAA	*102*		*40 - 170*
d3-NMePFOSA	*89.7*		*10 - 130*
d5-NEtFOSAA	*96.3*		*25 - 135*
d5-NEtPFOSA	*83.0*		*10 - 130*
d7-N-MeFOSE-M	*99.7*		*10 - 130*
d9-N-EtFOSE-M	*71.9*		*10 - 130*

Lab Sample ID: LCSD 410-707351/18-A
Matrix: Water
Analysis Batch: 707164

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA
Prep Batch: 707351

Analyte			Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)			37.8	29.9		ng/L		79	55 - 160	5	30
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)			76.8	67.5		ng/L		88	60 - 150	1	30
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)			75.0	64.5		ng/L		86	70 - 145	4	30

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LCSD 410-707351/18-A
Matrix: Water
Analysis Batch: 707164

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA
Prep Batch: 707351

Analyte			Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)			76.2	66.8		ng/L		88	65 - 155	0	30
3-Perfluoroheptylpropanoic acid (7:3 FTCA)			200	124		ng/L		62	50 - 145	6	30
3-Perfluoropentylpropanoic acid (5:3 FTCA)			200	133	*-	ng/L		66	70 - 135	8	30
3-Perfluoropropylpropanoic acid (3:3 FTCA)			80.0	80.6		ng/L		101	65 - 130	25	30
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)			37.8	33.1		ng/L		87	65 - 145	8	30
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)			37.4	31.0		ng/L		83	70 - 155	4	30
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)			30.0	25.8		ng/L		86	70 - 140	6	30
N-ethylperfluorooctane sulfonamide (NEtFOSA)			40.0	36.3		ng/L		91	65 - 145	2	30
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)			200	162		ng/L		81	70 - 135	14	30
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)			40.0	34.9		ng/L		87	70 - 145	3	30
N-methylperfluorooctane sulfonamide (NMeFOSA)			40.0	34.5		ng/L		86	60 - 150	2	30
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)			200	166		ng/L		83	70 - 145	2	30
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)			40.0	34.1		ng/L		85	50 - 140	0	30
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)			40.0	32.1		ng/L		80	50 - 150	2	30
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)			35.7	33.7		ng/L		94	70 - 140	2	30
Perfluoro-3-methoxypropanoic acid (PFMPA)			40.0	39.4		ng/L		98	55 - 140	6	30
Perfluoro-4-methoxybutanoic acid (PFMBA)			40.0	33.2		ng/L		83	60 - 150	2	30
Perfluorobutanesulfonic acid (PFBS)			35.5	30.8		ng/L		87	60 - 145	1	30
Perfluorobutanoic acid (PFBA)			80.0	65.8		ng/L		82	70 - 140	2	30
Perfluorodecanesulfonic acid (PFDS)			38.6	31.1		ng/L		81	60 - 145	0	30
Perfluorodecanoic acid (PFDA)			40.0	32.1		ng/L		80	70 - 140	0	30
Perfluorododecanesulfonic acid (PFDoS)			38.8	28.8		ng/L		74	50 - 145	0	30
Perfluorododecanoic acid (PFDoA)			40.0	37.7		ng/L		94	70 - 140	6	30
Perfluoroheptanesulfonic acid (PFHpS)			38.2	34.4		ng/L		90	70 - 150	5	30
Perfluoroheptanoic acid (PFHpA)			40.0	34.3		ng/L		86	70 - 150	1	30
Perfluorohexanesulfonic acid (PFHxS)			36.5	30.9		ng/L		85	65 - 145	0	30
Perfluorohexanoic acid (PFHxA)			40.0	34.0		ng/L		85	70 - 145	1	30
Perfluorononanesulfonic acid (PFNS)			38.5	32.3		ng/L		84	65 - 145	4	30
Perfluorononanoic acid (PFNA)			40.0	34.7		ng/L		87	70 - 150	5	30

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LCSD 410-707351/18-A
Matrix: Water
Analysis Batch: 707164

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA
Prep Batch: 707351

Analyte			Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
Perfluorooctanesulfonamide (PFOSA)			40.0	35.7		ng/L		89	70 - 145	1	30
Perfluorooctanesulfonic acid (PFOS)			37.2	30.5		ng/L		82	55 - 150	4	30
Perfluorooctanoic acid (PFOA)			40.0	33.2		ng/L		83	70 - 150	3	30
Perfluoropentanesulfonic acid (PFPeS)			37.6	31.7		ng/L		84	65 - 140	2	30
Perfluoropentanoic acid (PFPeA)			40.0	35.2		ng/L		88	65 - 135	2	30
Perfluorotetradecanoic acid (PFTeDA)			40.0	34.2		ng/L		86	60 - 140	7	30
Perfluorotridecanoic acid (PFTrDA)			40.0	33.3		ng/L		83	65 - 140	1	30
Perfluoroundecanoic acid (PFUnA)			40.0	36.3		ng/L		91	70 - 145	3	30

Isotope Dilution	*LCSD %Recovery*	*LCSD Qualifier*	*Limits*
13C2 4:2 FTS	*114*		*40 - 200*
13C2 6:2 FTS	*120*		*40 - 200*
13C2 8:2 FTS	*124*		*40 - 300*
13C2 PFDoA	*97.5*		*10 - 130*
13C2 PFTeDA	*90.1*		*10 - 130*
13C3 HFPO-DA	*113*		*40 - 130*
13C3 PFBS	*107*		*40 - 135*
13C3 PFHxS	*107*		*40 - 130*
13C4 PFBA	*108*		*5 - 130*
13C4 PFHpA	*112*		*40 - 130*
13C5 PFHxA	*106*		*40 - 130*
13C5 PFPeA	*108*		*40 - 130*
13C6 PFDA	*101*		*40 - 130*
13C7 PFUnA	*90.8*		*30 - 130*
13C8 FOSA	*104*		*40 - 130*
13C8 PFOA	*110*		*40 - 130*
13C8 PFOS	*104*		*40 - 130*
13C9 PFNA	*103*		*40 - 130*
d3-NMeFOSAA	*102*		*40 - 170*
d3-NMePFOSA	*93.0*		*10 - 130*
d5-NEtFOSAA	*97.0*		*25 - 135*
d5-NEtPFOSA	*86.9*		*10 - 130*
d7-N-MeFOSE-M	*113*		*10 - 130*
d9-N-EtFOSE-M	*105*		*10 - 130*

Lab Sample ID: MB 410-708885/15-A
Matrix: Water
Analysis Batch: 709341

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 708885

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)	ND		4.0	1.0	ng/L		10/03/25 13:46	10/06/25 04:01	1

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: MB 410-708885/15-A
Matrix: Water
Analysis Batch: 709341

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 708885

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)	ND		4.0	1.0	ng/L		10/03/25 13:46	10/06/25 04:01	1
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	ND		4.0	1.0	ng/L		10/03/25 13:46	10/06/25 04:01	1
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	ND		10	2.5	ng/L		10/03/25 13:46	10/06/25 04:01	1
3-Perfluoropentylpropanoic acid (5:3 FTCA)	ND		10	2.8	ng/L		10/03/25 13:46	10/06/25 04:01	1
3-Perfluoropropylpropanoic acid (3:3 FTCA)	ND		4.0	1.0	ng/L		10/03/25 13:46	10/06/25 04:01	1
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	ND		2.0	1.4	ng/L		10/03/25 13:46	10/06/25 04:01	1
N-ethylperfluorooctane sulfonamide (NEtFOSA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	ND		10	4.0	ng/L		10/03/25 13:46	10/06/25 04:01	1
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
N-methylperfluorooctane sulfonamide (NMeFOSA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	ND		10	2.5	ng/L		10/03/25 13:46	10/06/25 04:01	1
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluoro-3-methoxypropanoic acid (PFMPA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluoro-4-methoxybutanoic acid (PFMBA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorobutanesulfonic acid (PFBS)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorobutanoic acid (PFBA)	ND		4.0	1.1	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorodecanesulfonic acid (PFDS)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorodecanoic acid (PFDA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorododecanesulfonic acid (PFDoS)	ND		2.0	0.60	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorododecanoic acid (PFDoA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluoroheptanesulfonic acid (PFHpS)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluoroheptanoic acid (PFHpA)	ND		2.0	0.80	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorohexanesulfonic acid (PFHxS)	ND		2.0	0.80	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorohexanoic acid (PFHxA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorononanesulfonic acid (PFNS)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorononanoic acid (PFNA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorooctanesulfonamide (PFOSA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorooctanesulfonic acid (PFOS)	ND		2.0	1.3	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorooctanoic acid (PFOA)	ND		2.0	0.90	ng/L		10/03/25 13:46	10/06/25 04:01	1

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: MB 410-708885/15-A
Matrix: Water
Analysis Batch: 709341

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 708885

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoropentanesulfonic acid (PFPeS)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluoropentanoic acid (PFPeA)	ND		2.0	0.60	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorotetradecanoic acid (PFTeDA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluorotridecanoic acid (PFTrDA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1
Perfluoroundecanoic acid (PFUnA)	ND		2.0	0.50	ng/L		10/03/25 13:46	10/06/25 04:01	1

Isotope Dilution	*MB %Recovery*	*MB Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
13C2 4:2 FTS	*86.4*		*40 - 200*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C2 6:2 FTS	*86.3*		*40 - 200*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C2 8:2 FTS	*77.6*		*40 - 300*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C2 PFDoA	*93.1*		*10 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C2 PFTeDA	*74.6*		*10 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C3 HFPO-DA	*101*		*40 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C3 PFBS	*99.4*		*40 - 135*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C3 PFHxS	*96.6*		*40 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C4 PFBA	*104*		*5 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C4 PFHpA	*102*		*40 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C5 PFHxA	*98.5*		*40 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C5 PFPeA	*100*		*40 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C6 PFDA	*98.2*		*40 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C7 PFUnA	*92.2*		*30 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C8 FOSA	*83.8*		*40 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C8 PFOA	*94.1*		*40 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C8 PFOS	*95.6*		*40 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
13C9 PFNA	*93.6*		*40 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
d3-NMeFOSAA	*81.4*		*40 - 170*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
d3-NMePFOSA	*74.1*		*10 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
d5-NEtFOSAA	*77.1*		*25 - 135*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
d5-NEtPFOSA	*68.1*		*10 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
d7-N-MeFOSE-M	*50.9*		*10 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*
d9-N-EtFOSE-M	*43.2*		*10 - 130*	*10/03/25 13:46*	*10/06/25 04:01*	*1*

Lab Sample ID: LCS 410-708885/16-A
Matrix: Water
Analysis Batch: 709341

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 708885

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)	37.8	32.3		ng/L		86	55 - 160
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)	76.8	66.4		ng/L		86	60 - 150
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)	75.0	65.5		ng/L		87	70 - 145
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	76.2	70.0		ng/L		92	65 - 155
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	200	133		ng/L		66	50 - 145
3-Perfluoropentylpropanoic acid (5:3 FTCA)	200	147		ng/L		73	70 - 135

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LCS 410-708885/16-A
Matrix: Water
Analysis Batch: 709341

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 708885

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
3-Perfluoropropylpropanoic acid (3:3 FTCA)	80.0	66.9		ng/L		84	65 - 130
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	37.8	32.0		ng/L		85	65 - 145
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	37.4	33.8		ng/L		90	70 - 155
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	30.0	26.7		ng/L		89	70 - 140
N-ethylperfluorooctane sulfonamide (NEtFOSA)	40.0	38.6		ng/L		97	65 - 145
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	200	186		ng/L		93	70 - 135
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	40.0	37.6		ng/L		94	70 - 145
N-methylperfluorooctane sulfonamide (NMeFOSA)	40.0	38.0		ng/L		95	60 - 150
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	200	178		ng/L		89	70 - 145
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	40.0	36.0		ng/L		90	50 - 140
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	40.0	38.1		ng/L		95	50 - 150
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	35.7	34.1		ng/L		96	70 - 140
Perfluoro-3-methoxypropanoic acid (PFMPA)	40.0	38.6		ng/L		97	55 - 140
Perfluoro-4-methoxybutanoic acid (PFMBA)	40.0	35.8		ng/L		90	60 - 150
Perfluorobutanesulfonic acid (PFBS)	35.5	32.4		ng/L		91	60 - 145
Perfluorobutanoic acid (PFBA)	80.0	70.0		ng/L		88	70 - 140
Perfluorodecanesulfonic acid (PFDS)	38.6	32.0		ng/L		83	60 - 145
Perfluorodecanoic acid (PFDA)	40.0	34.2		ng/L		86	70 - 140
Perfluorododecanesulfonic acid (PFDoS)	38.8	30.1		ng/L		78	50 - 145
Perfluorododecanoic acid (PFDoA)	40.0	39.8		ng/L		99	70 - 140
Perfluoroheptanesulfonic acid (PFHpS)	38.2	36.2		ng/L		95	70 - 150
Perfluoroheptanoic acid (PFHpA)	40.0	36.7		ng/L		92	70 - 150
Perfluorohexanesulfonic acid (PFHxS)	36.5	32.5		ng/L		89	65 - 145
Perfluorohexanoic acid (PFHxA)	40.0	35.7		ng/L		89	70 - 145
Perfluorononanesulfonic acid (PFNS)	38.5	33.5		ng/L		87	65 - 145
Perfluorononanoic acid (PFNA)	40.0	38.4		ng/L		96	70 - 150
Perfluorooctanesulfonamide (PFOSA)	40.0	38.5		ng/L		96	70 - 145
Perfluorooctanesulfonic acid (PFOS)	37.2	31.8		ng/L		86	55 - 150
Perfluorooctanoic acid (PFOA)	40.0	35.0		ng/L		88	70 - 150
Perfluoropentanesulfonic acid (PFPeS)	37.6	34.6		ng/L		92	65 - 140

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LCS 410-708885/16-A
Matrix: Water
Analysis Batch: 709341

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 708885

Analyte			Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits		
Perfluoropentanoic acid (PFPeA)			40.0	38.5		ng/L		96	65 - 135		
Perfluorotetradecanoic acid (PFTeDA)			40.0	38.0		ng/L		95	60 - 140		
Perfluorotridecanoic acid (PFTrDA)			40.0	35.4		ng/L		89	65 - 140		
Perfluoroundecanoic acid (PFUnA)			40.0	37.3		ng/L		93	70 - 145		

Isotope Dilution	*LCS %Recovery*	*LCS Qualifier*	*Limits*
13C2 4:2 FTS	*89.7*		*40 - 200*
13C2 6:2 FTS	*82.4*		*40 - 200*
13C2 8:2 FTS	*74.5*		*40 - 300*
13C2 PFDoA	*89.9*		*10 - 130*
13C2 PFTeDA	*77.4*		*10 - 130*
13C3 HFPO-DA	*108*		*40 - 130*
13C3 PFBS	*94.6*		*40 - 135*
13C3 PFHxS	*91.5*		*40 - 130*
13C4 PFBA	*101*		*5 - 130*
13C4 PFHpA	*95.8*		*40 - 130*
13C5 PFHxA	*96.5*		*40 - 130*
13C5 PFPeA	*99.4*		*40 - 130*
13C6 PFDA	*92.2*		*40 - 130*
13C7 PFUnA	*92.5*		*30 - 130*
13C8 FOSA	*81.7*		*40 - 130*
13C8 PFOA	*93.9*		*40 - 130*
13C8 PFOS	*91.7*		*40 - 130*
13C9 PFNA	*88.0*		*40 - 130*
d3-NMeFOSAA	*76.0*		*40 - 170*
d3-NMePFOSA	*70.7*		*10 - 130*
d5-NEtFOSAA	*70.2*		*25 - 135*
d5-NEtPFOSA	*64.8*		*10 - 130*
d7-N-MeFOSE-M	*61.6*		*10 - 130*
d9-N-EtFOSE-M	*54.0*		*10 - 130*

Lab Sample ID: LCSD 410-708885/17-A
Matrix: Water
Analysis Batch: 709341

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA
Prep Batch: 708885

Analyte			Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)			37.8	34.2		ng/L		90	55 - 160	6	30
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)			76.8	74.6		ng/L		97	60 - 150	12	30
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)			75.0	72.2		ng/L		96	70 - 145	10	30
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)			76.2	71.9		ng/L		94	65 - 155	3	30
3-Perfluoroheptylpropanoic acid (7:3 FTCA)			200	149		ng/L		74	50 - 145	12	30

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LCSD 410-708885/17-A
Matrix: Water
Analysis Batch: 709341

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA
Prep Batch: 708885

Analyte	Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
3-Perfluoropentylpropanoic acid (5:3 FTCA)	200	159		ng/L		80	70 - 135	8	30
3-Perfluoropropylpropanoic acid (3:3 FTCA)	80.0	71.4		ng/L		89	65 - 130	7	30
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	37.8	35.5		ng/L		94	65 - 145	10	30
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	37.4	36.0		ng/L		96	70 - 155	7	30
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	30.0	28.4		ng/L		95	70 - 140	6	30
N-ethylperfluorooctane sulfonamide (NEtFOSA)	40.0	40.1		ng/L		100	65 - 145	4	30
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	200	196		ng/L		98	70 - 135	5	30
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	40.0	40.3		ng/L		101	70 - 145	7	30
N-methylperfluorooctane sulfonamide (NMeFOSA)	40.0	39.9		ng/L		100	60 - 150	5	30
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	200	186		ng/L		93	70 - 145	5	30
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	40.0	40.2		ng/L		100	50 - 140	11	30
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	40.0	36.8		ng/L		92	50 - 150	3	30
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	35.7	38.5		ng/L		108	70 - 140	12	30
Perfluoro-3-methoxypropanoic acid (PFMPA)	40.0	40.0		ng/L		100	55 - 140	3	30
Perfluoro-4-methoxybutanoic acid (PFMBA)	40.0	37.9		ng/L		95	60 - 150	6	30
Perfluorobutanesulfonic acid (PFBS)	35.5	34.9		ng/L		98	60 - 145	8	30
Perfluorobutanoic acid (PFBA)	80.0	74.7		ng/L		93	70 - 140	7	30
Perfluorodecanesulfonic acid (PFDS)	38.6	34.2		ng/L		89	60 - 145	7	30
Perfluorodecanoic acid (PFDA)	40.0	36.9		ng/L		92	70 - 140	7	30
Perfluorododecanesulfonic acid (PFDoS)	38.8	32.5		ng/L		84	50 - 145	8	30
Perfluorododecanoic acid (PFDoA)	40.0	40.3		ng/L		101	70 - 140	1	30
Perfluoroheptanesulfonic acid (PFHpS)	38.2	38.8		ng/L		102	70 - 150	7	30
Perfluoroheptanoic acid (PFHpA)	40.0	39.2		ng/L		98	70 - 150	7	30
Perfluorohexanesulfonic acid (PFHxS)	36.5	35.6		ng/L		98	65 - 145	9	30
Perfluorohexanoic acid (PFHxA)	40.0	37.8		ng/L		95	70 - 145	6	30
Perfluorononanesulfonic acid (PFNS)	38.5	36.1		ng/L		94	65 - 145	8	30
Perfluorononanoic acid (PFNA)	40.0	39.9		ng/L		100	70 - 150	4	30
Perfluorooctanesulfonamide (PFOSA)	40.0	41.0		ng/L		102	70 - 145	6	30
Perfluorooctanesulfonic acid (PFOS)	37.2	34.3		ng/L		92	55 - 150	7	30
Perfluorooctanoic acid (PFOA)	40.0	37.6		ng/L		94	70 - 150	7	30

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LCSD 410-708885/17-A
Matrix: Water
Analysis Batch: 709341

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA
Prep Batch: 708885

Analyte			Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
Perfluoropentanesulfonic acid (PFPeS)			37.6	37.4		ng/L		99	65 - 140	8	30
Perfluoropentanoic acid (PFPeA)			40.0	40.8		ng/L		102	65 - 135	6	30
Perfluorotetradecanoic acid (PFTeDA)			40.0	40.5		ng/L		101	60 - 140	6	30
Perfluorotridecanoic acid (PFTrDA)			40.0	37.0		ng/L		93	65 - 140	4	30
Perfluoroundecanoic acid (PFUnA)			40.0	39.3		ng/L		98	70 - 145	5	30

Isotope Dilution	*LCSD %Recovery*	*LCSD Qualifier*	*Limits*
13C2 4:2 FTS	*89.8*		*40 - 200*
13C2 6:2 FTS	*84.8*		*40 - 200*
13C2 8:2 FTS	*81.1*		*40 - 300*
13C2 PFDoA	*96.3*		*10 - 130*
13C2 PFTeDA	*82.1*		*10 - 130*
13C3 HFPO-DA	*107*		*40 - 130*
13C3 PFBS	*99.8*		*40 - 135*
13C3 PFHxS	*96.3*		*40 - 130*
13C4 PFBA	*104*		*5 - 130*
13C4 PFHpA	*102*		*40 - 130*
13C5 PFHxA	*98.5*		*40 - 130*
13C5 PFPeA	*103*		*40 - 130*
13C6 PFDA	*98.3*		*40 - 130*
13C7 PFUnA	*92.3*		*30 - 130*
13C8 FOSA	*83.7*		*40 - 130*
13C8 PFOA	*97.5*		*40 - 130*
13C8 PFOS	*97.5*		*40 - 130*
13C9 PFNA	*100*		*40 - 130*
d3-NMeFOSAA	*77.4*		*40 - 170*
d3-NMePFOSA	*70.0*		*10 - 130*
d5-NEtFOSAA	*73.9*		*25 - 135*
d5-NEtPFOSA	*64.9*		*10 - 130*
d7-N-MeFOSE-M	*59.4*		*10 - 130*
d9-N-EtFOSE-M	*52.1*		*10 - 130*

Lab Sample ID: LLCS 410-708885/18-A
Matrix: Water
Analysis Batch: 709341

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 708885

Analyte			Spike Added	LLCS Result	LLCS Qualifier	Unit	D	%Rec	%Rec Limits
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)			3.78	3.23		ng/L		85	55 - 160
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)			7.68	6.15		ng/L		80	60 - 150
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)			7.50	6.57		ng/L		88	70 - 145
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)			7.62	6.90		ng/L		91	65 - 155

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LLCS 410-708885/18-A
Matrix: Water
Analysis Batch: 709341

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 708885

Analyte	Spike Added	LLCS Result	LLCS Qualifier	Unit	D	%Rec	%Rec Limits
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	20.0	15.2		ng/L		76	50 - 145
3-Perfluoropentylpropanoic acid (5:3 FTCA)	20.0	16.7		ng/L		84	70 - 135
3-Perfluoropropylpropanoic acid (3:3 FTCA)	8.00	6.52		ng/L		82	65 - 130
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	3.78	3.22		ng/L		85	65 - 145
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	3.74	3.28		ng/L		88	70 - 155
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	3.00	2.59		ng/L		86	70 - 140
N-ethylperfluorooctane sulfonamide (NEtFOSA)	4.00	3.21		ng/L		80	65 - 145
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	20.0	18.4		ng/L		92	70 - 135
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	4.00	3.66		ng/L		92	70 - 145
N-methylperfluorooctane sulfonamide (NMeFOSA)	4.00	3.50		ng/L		87	60 - 150
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	20.0	16.9		ng/L		84	70 - 145
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	4.00	2.70		ng/L		68	50 - 140
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	4.00	3.32		ng/L		83	50 - 150
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	3.57	3.19		ng/L		90	70 - 140
Perfluoro-3-methoxypropanoic acid (PFMPA)	4.00	3.54		ng/L		89	55 - 140
Perfluoro-4-methoxybutanoic acid (PFMBA)	4.00	3.22		ng/L		81	60 - 150
Perfluorobutanesulfonic acid (PFBS)	3.55	2.98		ng/L		84	60 - 145
Perfluorobutanoic acid (PFBA)	8.00	6.43		ng/L		80	70 - 140
Perfluorodecanesulfonic acid (PFDS)	3.86	2.99		ng/L		78	60 - 145
Perfluorodecanoic acid (PFDA)	4.00	3.46		ng/L		87	70 - 140
Perfluorododecanesulfonic acid (PFDoS)	3.88	2.89		ng/L		75	50 - 145
Perfluorododecanoic acid (PFDoA)	4.00	3.50		ng/L		87	70 - 140
Perfluoroheptanesulfonic acid (PFHpS)	3.82	3.40		ng/L		89	70 - 150
Perfluoroheptanoic acid (PFHpA)	4.00	3.65		ng/L		91	70 - 150
Perfluorohexanesulfonic acid (PFHxS)	3.65	3.23		ng/L		89	65 - 145
Perfluorohexanoic acid (PFHxA)	4.00	3.72		ng/L		93	70 - 145
Perfluorononanesulfonic acid (PFNS)	3.85	2.87		ng/L		75	65 - 145
Perfluorononanoic acid (PFNA)	4.00	3.47		ng/L		87	70 - 150
Perfluorooctanesulfonamide (PFOSA)	4.00	3.81		ng/L		95	70 - 145

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LLCS 410-708885/18-A
Matrix: Water
Analysis Batch: 709341

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 708885

Analyte	Spike Added	LLCS Result	LLCS Qualifier	Unit	D	%Rec	%Rec Limits
Perfluorooctanesulfonic acid (PFOS)	3.72	3.32		ng/L		89	55 - 150
Perfluorooctanoic acid (PFOA)	4.00	3.34		ng/L		84	70 - 150
Perfluoropentanesulfonic acid (PFPeS)	3.76	3.17		ng/L		84	65 - 140
Perfluoropentanoic acid (PFPeA)	4.00	3.49		ng/L		87	65 - 135
Perfluorotetradecanoic acid (PFTeDA)	4.00	3.79		ng/L		95	60 - 140
Perfluorotridecanoic acid (PFTrDA)	4.00	3.20		ng/L		80	65 - 140
Perfluoroundecanoic acid (PFUnA)	4.00	3.51		ng/L		88	70 - 145

Isotope Dilution	*LLCS %Recovery*	*LLCS Qualifier*	*Limits*
13C2 4:2 FTS	*89.8*		*40 - 200*
13C2 6:2 FTS	*85.9*		*40 - 200*
13C2 8:2 FTS	*78.6*		*40 - 300*
13C2 PFDoA	*98.4*		*10 - 130*
13C2 PFTeDA	*81.4*		*10 - 130*
13C3 HFPO-DA	*97.0*		*40 - 130*
13C3 PFBS	*97.5*		*40 - 135*
13C3 PFHxS	*95.9*		*40 - 130*
13C4 PFBA	*98.0*		*5 - 130*
13C4 PFHpA	*94.9*		*40 - 130*
13C5 PFHxA	*89.3*		*40 - 130*
13C5 PFPeA	*94.2*		*40 - 130*
13C6 PFDA	*97.1*		*40 - 130*
13C7 PFUnA	*94.8*		*30 - 130*
13C8 FOSA	*77.2*		*40 - 130*
13C8 PFOA	*90.4*		*40 - 130*
13C8 PFOS	*93.8*		*40 - 130*
13C9 PFNA	*91.9*		*40 - 130*
d3-NMeFOSAA	*80.3*		*40 - 170*
d3-NMePFOSA	*66.5*		*10 - 130*
d5-NEtFOSAA	*78.0*		*25 - 135*
d5-NEtPFOSA	*63.7*		*10 - 130*
d7-N-MeFOSE-M	*59.3*		*10 - 130*
d9-N-EtFOSE-M	*54.7*		*10 - 130*

Lab Sample ID: MB 410-710704/5-A
Matrix: Water
Analysis Batch: 710552

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 710704

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)	ND		4.0	1.0	ng/L		10/08/25 08:36	10/08/25 22:56	1
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)	ND		4.0	1.0	ng/L		10/08/25 08:36	10/08/25 22:56	1

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: MB 410-710704/5-A
Matrix: Water
Analysis Batch: 710552

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 710704

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	ND		4.0	1.0	ng/L		10/08/25 08:36	10/08/25 22:56	1
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	ND		10	2.5	ng/L		10/08/25 08:36	10/08/25 22:56	1
3-Perfluoropentylpropanoic acid (5:3 FTCA)	ND		10	2.8	ng/L		10/08/25 08:36	10/08/25 22:56	1
3-Perfluoropropylpropanoic acid (3:3 FTCA)	ND		4.0	1.0	ng/L		10/08/25 08:36	10/08/25 22:56	1
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	ND		2.0	1.4	ng/L		10/08/25 08:36	10/08/25 22:56	1
N-ethylperfluorooctane sulfonamide (NEtFOSA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	ND		10	4.0	ng/L		10/08/25 08:36	10/08/25 22:56	1
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
N-methylperfluorooctane sulfonamide (NMeFOSA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	ND		10	2.5	ng/L		10/08/25 08:36	10/08/25 22:56	1
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluoro-3-methoxypropanoic acid (PFMPA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluoro-4-methoxybutanoic acid (PFMBA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorobutanesulfonic acid (PFBS)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorobutanoic acid (PFBA)	ND		4.0	1.1	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorodecanesulfonic acid (PFDS)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorodecanoic acid (PFDA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorododecanesulfonic acid (PFDoS)	ND		2.0	0.60	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorododecanoic acid (PFDoA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluoroheptanesulfonic acid (PFHpS)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluoroheptanoic acid (PFHpA)	ND		2.0	0.80	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorohexanesulfonic acid (PFHxS)	ND		2.0	0.80	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorohexanoic acid (PFHxA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorononanesulfonic acid (PFNS)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorononanoic acid (PFNA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorooctanesulfonamide (PFOSA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorooctanesulfonic acid (PFOS)	ND		2.0	1.3	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorooctanoic acid (PFOA)	ND		2.0	0.90	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluoropentanesulfonic acid (PFPeS)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluoropentanoic acid (PFPeA)	ND		2.0	0.60	ng/L		10/08/25 08:36	10/08/25 22:56	1

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: MB 410-710704/5-A
Matrix: Water
Analysis Batch: 710552

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 710704

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorotetradecanoic acid (PFTeDA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluorotridecanoic acid (PFTrDA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1
Perfluoroundecanoic acid (PFUnA)	ND		2.0	0.50	ng/L		10/08/25 08:36	10/08/25 22:56	1

Isotope Dilution	*MB %Recovery*	*MB Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
13C2 4:2 FTS	*107*		*40 - 200*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C2 6:2 FTS	*98.6*		*40 - 200*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C2 8:2 FTS	*105*		*40 - 300*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C2 PFDoA	*89.7*		*10 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C2 PFTeDA	*67.0*		*10 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C3 HFPO-DA	*89.7*		*40 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C3 PFBS	*103*		*40 - 135*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C3 PFHxS	*87.4*		*40 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C4 PFBA	*102*		*5 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C4 PFHpA	*96.9*		*40 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C5 PFHxA	*99.0*		*40 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C5 PFPeA	*101*		*40 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C6 PFDA	*95.0*		*40 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C7 PFUnA	*89.0*		*30 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C8 FOSA	*90.8*		*40 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C8 PFOA	*92.9*		*40 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C8 PFOS	*96.1*		*40 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
13C9 PFNA	*92.7*		*40 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
d3-NMeFOSAA	*88.9*		*40 - 170*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
d3-NMePFOSA	*76.5*		*10 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
d5-NEtFOSAA	*79.1*		*25 - 135*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
d5-NEtPFOSA	*73.8*		*10 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
d7-N-MeFOSE-M	*56.6*		*10 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*
d9-N-EtFOSE-M	*43.1*		*10 - 130*	*10/08/25 08:36*	*10/08/25 22:56*	*1*

Lab Sample ID: LCS 410-710704/6-A
Matrix: Water
Analysis Batch: 710552

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 710704

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)	37.8	32.7		ng/L		87	55 - 160
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)	76.8	73.7		ng/L		96	60 - 150
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)	75.0	68.0		ng/L		91	70 - 145
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)	76.2	73.4		ng/L		96	65 - 155
3-Perfluoroheptylpropanoic acid (7:3 FTCA)	200	135		ng/L		68	50 - 145
3-Perfluoropentylpropanoic acid (5:3 FTCA)	200	163		ng/L		81	70 - 135
3-Perfluoropropylpropanoic acid (3:3 FTCA)	80.0	64.4		ng/L		81	65 - 130

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LCS 410-710704/6-A
Matrix: Water
Analysis Batch: 710552

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 710704

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	37.8	39.0		ng/L		103	65 - 145
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	37.4	35.5		ng/L		95	70 - 155
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	30.0	28.0		ng/L		93	70 - 140
N-ethylperfluorooctane sulfonamide (NEtFOSA)	40.0	37.7		ng/L		94	65 - 145
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	200	216		ng/L		108	70 - 135
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	40.0	38.7		ng/L		97	70 - 145
N-methylperfluorooctane sulfonamide (NMeFOSA)	40.0	37.5		ng/L		94	60 - 150
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	200	193		ng/L		96	70 - 145
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	40.0	39.3		ng/L		98	50 - 140
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	40.0	31.6		ng/L		79	50 - 150
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	35.7	36.1		ng/L		101	70 - 140
Perfluoro-3-methoxypropanoic acid (PFMPA)	40.0	34.0		ng/L		85	55 - 140
Perfluoro-4-methoxybutanoic acid (PFMBA)	40.0	34.6		ng/L		86	60 - 150
Perfluorobutanesulfonic acid (PFBS)	35.5	32.3		ng/L		91	60 - 145
Perfluorobutanoic acid (PFBA)	80.0	74.9		ng/L		94	70 - 140
Perfluorodecanesulfonic acid (PFDS)	38.6	33.8		ng/L		88	60 - 145
Perfluorodecanoic acid (PFDA)	40.0	35.8		ng/L		90	70 - 140
Perfluorododecanesulfonic acid (PFDoS)	38.8	30.4		ng/L		78	50 - 145
Perfluorododecanoic acid (PFDoA)	40.0	34.8		ng/L		87	70 - 140
Perfluoroheptanesulfonic acid (PFHpS)	38.2	33.8		ng/L		88	70 - 150
Perfluoroheptanoic acid (PFHpA)	40.0	37.5		ng/L		94	70 - 150
Perfluorohexanesulfonic acid (PFHxS)	36.5	39.4		ng/L		108	65 - 145
Perfluorohexanoic acid (PFHxA)	40.0	38.8		ng/L		97	70 - 145
Perfluorononanesulfonic acid (PFNS)	38.5	35.0		ng/L		91	65 - 145
Perfluorononanoic acid (PFNA)	40.0	35.7		ng/L		89	70 - 150
Perfluorooctanesulfonamide (PFOSA)	40.0	39.0		ng/L		97	70 - 145
Perfluorooctanesulfonic acid (PFOS)	37.2	32.6		ng/L		88	55 - 150
Perfluorooctanoic acid (PFOA)	40.0	36.6		ng/L		92	70 - 150
Perfluoropentanesulfonic acid (PFPeS)	37.6	40.5		ng/L		108	65 - 140
Perfluoropentanoic acid (PFPeA)	40.0	35.1		ng/L		88	65 - 135

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LCS 410-710704/6-A
Matrix: Water
Analysis Batch: 710552

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 710704

Analyte			Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits		
Perfluorotetradecanoic acid (PFTeDA)			40.0	35.4		ng/L		89	60 - 140		
Perfluorotridecanoic acid (PFTrDA)			40.0	37.3		ng/L		93	65 - 140		
Perfluoroundecanoic acid (PFUnA)			40.0	35.4		ng/L		88	70 - 145		

Isotope Dilution	*LCS %Recovery*	*LCS Qualifier*	*Limits*
13C2 4:2 FTS	*111*		*40 - 200*
13C2 6:2 FTS	*93.4*		*40 - 200*
13C2 8:2 FTS	*110*		*40 - 300*
13C2 PFDoA	*93.9*		*10 - 130*
13C2 PFTeDA	*79.6*		*10 - 130*
13C3 HFPO-DA	*91.7*		*40 - 130*
13C3 PFBS	*110*		*40 - 135*
13C3 PFHxS	*95.4*		*40 - 130*
13C4 PFBA	*103*		*5 - 130*
13C4 PFHpA	*93.7*		*40 - 130*
13C5 PFHxA	*96.2*		*40 - 130*
13C5 PFPeA	*102*		*40 - 130*
13C6 PFDA	*96.0*		*40 - 130*
13C7 PFUnA	*89.6*		*30 - 130*
13C8 FOSA	*88.0*		*40 - 130*
13C8 PFOA	*93.0*		*40 - 130*
13C8 PFOS	*98.1*		*40 - 130*
13C9 PFNA	*103*		*40 - 130*
d3-NMeFOSAA	*88.0*		*40 - 170*
d3-NMePFOSA	*78.5*		*10 - 130*
d5-NEtFOSAA	*78.6*		*25 - 135*
d5-NEtPFOSA	*78.4*		*10 - 130*
d7-N-MeFOSE-M	*72.7*		*10 - 130*
d9-N-EtFOSE-M	*58.2*		*10 - 130*

Lab Sample ID: LCSD 410-710704/7-A
Matrix: Water
Analysis Batch: 710552

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA
Prep Batch: 710704

Analyte			Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)			37.8	31.5		ng/L		83	55 - 160	4	30
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)			76.8	74.8		ng/L		97	60 - 150	1	30
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)			75.0	64.4		ng/L		86	70 - 145	6	30
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)			76.2	81.7		ng/L		107	65 - 155	11	30
3-Perfluoroheptylpropanoic acid (7:3 FTCA)			200	127		ng/L		64	50 - 145	6	30
3-Perfluoropentylpropanoic acid (5:3 FTCA)			200	158		ng/L		79	70 - 135	3	30

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LCSD 410-710704/7-A
Matrix: Water
Analysis Batch: 710552

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA
Prep Batch: 710704

Analyte			Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
3-Perfluoropropylpropanoic acid (3:3 FTCA)			80.0	63.7		ng/L		80	65 - 130	1	30
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)			37.8	40.4		ng/L		107	65 - 145	4	30
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)			37.4	35.4		ng/L		95	70 - 155	0	30
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)			30.0	29.6		ng/L		99	70 - 140	5	30
N-ethylperfluorooctane sulfonamide (NEtFOSA)			40.0	38.7		ng/L		97	65 - 145	3	30
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)			200	203		ng/L		101	70 - 135	6	30
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)			40.0	37.9		ng/L		95	70 - 145	2	30
N-methylperfluorooctane sulfonamide (NMeFOSA)			40.0	38.8		ng/L		97	60 - 150	3	30
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)			200	197		ng/L		98	70 - 145	2	30
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)			40.0	38.0		ng/L		95	50 - 140	3	30
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)			40.0	30.4		ng/L		76	50 - 150	4	30
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)			35.7	34.5		ng/L		97	70 - 140	5	30
Perfluoro-3-methoxypropanoic acid (PFMPA)			40.0	34.3		ng/L		86	55 - 140	1	30
Perfluoro-4-methoxybutanoic acid (PFMBA)			40.0	35.3		ng/L		88	60 - 150	2	30
Perfluorobutanesulfonic acid (PFBS)			35.5	36.4		ng/L		103	60 - 145	12	30
Perfluorobutanoic acid (PFBA)			80.0	75.8		ng/L		95	70 - 140	1	30
Perfluorodecanesulfonic acid (PFDS)			38.6	34.2		ng/L		89	60 - 145	1	30
Perfluorodecanoic acid (PFDA)			40.0	34.6		ng/L		87	70 - 140	3	30
Perfluorododecanesulfonic acid (PFDoS)			38.8	23.3		ng/L		60	50 - 145	26	30
Perfluorododecanoic acid (PFDoA)			40.0	34.8		ng/L		87	70 - 140	0	30
Perfluoroheptanesulfonic acid (PFHpS)			38.2	35.6		ng/L		93	70 - 150	5	30
Perfluoroheptanoic acid (PFHpA)			40.0	37.2		ng/L		93	70 - 150	1	30
Perfluorohexanesulfonic acid (PFHxS)			36.5	39.8		ng/L		109	65 - 145	1	30
Perfluorohexanoic acid (PFHxA)			40.0	36.9		ng/L		92	70 - 145	5	30
Perfluorononanesulfonic acid (PFNS)			38.5	34.5		ng/L		90	65 - 145	1	30
Perfluorononanoic acid (PFNA)			40.0	38.5		ng/L		96	70 - 150	7	30
Perfluorooctanesulfonamide (PFOSA)			40.0	39.0		ng/L		98	70 - 145	0	30
Perfluorooctanesulfonic acid (PFOS)			37.2	32.1		ng/L		86	55 - 150	1	30
Perfluorooctanoic acid (PFOA)			40.0	37.4		ng/L		94	70 - 150	2	30
Perfluoropentanesulfonic acid (PFPeS)			37.6	41.5		ng/L		110	65 - 140	2	30

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LCSD 410-710704/7-A
Matrix: Water
Analysis Batch: 710552

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA
Prep Batch: 710704

Analyte			Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
Perfluoropentanoic acid (PFPeA)			40.0	36.2		ng/L		91	65 - 135	3	30
Perfluorotetradecanoic acid (PFTeDA)			40.0	30.6		ng/L		77	60 - 140	15	30
Perfluorotridecanoic acid (PFTrDA)			40.0	35.8		ng/L		89	65 - 140	4	30
Perfluoroundecanoic acid (PFUnA)			40.0	37.0		ng/L		93	70 - 145	5	30

Isotope Dilution	*LCSD %Recovery*	*LCSD Qualifier*	*Limits*
13C2 4:2 FTS	*109*		*40 - 200*
13C2 6:2 FTS	*94.3*		*40 - 200*
13C2 8:2 FTS	*103*		*40 - 300*
13C2 PFDoA	*87.9*		*10 - 130*
13C2 PFTeDA	*63.6*		*10 - 130*
13C3 HFPO-DA	*88.6*		*40 - 130*
13C3 PFBS	*100*		*40 - 135*
13C3 PFHxS	*89.7*		*40 - 130*
13C4 PFBA	*97.7*		*5 - 130*
13C4 PFHpA	*92.3*		*40 - 130*
13C5 PFHxA	*100*		*40 - 130*
13C5 PFPeA	*97.4*		*40 - 130*
13C6 PFDA	*90.7*		*40 - 130*
13C7 PFUnA	*86.1*		*30 - 130*
13C8 FOSA	*86.6*		*40 - 130*
13C8 PFOA	*88.4*		*40 - 130*
13C8 PFOS	*95.5*		*40 - 130*
13C9 PFNA	*93.5*		*40 - 130*
d3-NMeFOSAA	*87.5*		*40 - 170*
d3-NMePFOSA	*72.3*		*10 - 130*
d5-NEtFOSAA	*76.6*		*25 - 135*
d5-NEtPFOSA	*71.0*		*10 - 130*
d7-N-MeFOSE-M	*61.1*		*10 - 130*
d9-N-EtFOSE-M	*55.0*		*10 - 130*

Lab Sample ID: LLCS 410-710704/8-A
Matrix: Water
Analysis Batch: 710552

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 710704

Analyte			Spike Added	LLCS Result	LLCS Qualifier	Unit	D	%Rec	%Rec Limits
11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)			3.78	3.03		ng/L		80	55 - 160
1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)			7.68	6.79		ng/L		88	60 - 150
1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)			7.50	6.10		ng/L		81	70 - 145
1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)			7.62	6.58		ng/L		86	65 - 155
3-Perfluoroheptylpropanoic acid (7:3 FTCA)			20.0	12.0		ng/L		60	50 - 145

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LLCS 410-710704/8-A
Matrix: Water
Analysis Batch: 710552

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 710704

Analyte	Spike Added	LLCS Result	LLCS Qualifier	Unit	D	%Rec	%Rec Limits
3-Perfluoropentylpropanoic acid (5:3 FTCA)	20.0	16.8		ng/L		84	70 - 135
3-Perfluoropropylpropanoic acid (3:3 FTCA)	8.00	5.57		ng/L		70	65 - 130
4,8-Dioxa-3H-perfluorononanoic acid (ADONA)	3.78	3.94		ng/L		104	65 - 145
9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid (9Cl-PF3ONS)	3.74	3.43		ng/L		92	70 - 155
Hexafluoropropylene Oxide Dimer Acid (HFPO-DA)	3.00	2.31		ng/L		77	70 - 140
N-ethylperfluorooctane sulfonamide (NEtFOSA)	4.00	3.07		ng/L		77	65 - 145
N-ethylperfluorooctane sulfonamidoethanol (NEtFOSE)	20.0	22.6		ng/L		113	70 - 135
N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)	4.00	3.77		ng/L		94	70 - 145
N-methylperfluorooctane sulfonamide (NMeFOSA)	4.00	3.54		ng/L		89	60 - 150
N-methylperfluorooctane sulfonamidoethanol (NMeFOSE)	20.0	22.5		ng/L		113	70 - 145
N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)	4.00	2.58		ng/L		65	50 - 140
Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)	4.00	3.20		ng/L		80	50 - 150
Perfluoro (2-ethoxyethane) sulfonic acid (PFEESA)	3.57	2.64		ng/L		74	70 - 140
Perfluoro-3-methoxypropanoic acid (PFMPA)	4.00	3.03		ng/L		76	55 - 140
Perfluoro-4-methoxybutanoic acid (PFMBA)	4.00	3.16		ng/L		79	60 - 150
Perfluorobutanesulfonic acid (PFBS)	3.55	3.02		ng/L		85	60 - 145
Perfluorobutanoic acid (PFBA)	8.00	5.92		ng/L		74	70 - 140
Perfluorodecanesulfonic acid (PFDS)	3.86	2.98		ng/L		77	60 - 145
Perfluorodecanoic acid (PFDA)	4.00	3.08		ng/L		77	70 - 140
Perfluorododecanesulfonic acid (PFDoS)	3.88	2.15		ng/L		55	50 - 145
Perfluorododecanoic acid (PFDoA)	4.00	3.05		ng/L		76	70 - 140
Perfluoroheptanesulfonic acid (PFHpS)	3.82	3.10		ng/L		81	70 - 150
Perfluoroheptanoic acid (PFHpA)	4.00	3.60		ng/L		90	70 - 150
Perfluorohexanesulfonic acid (PFHxS)	3.65	3.76		ng/L		103	65 - 145
Perfluorohexanoic acid (PFHxA)	4.00	3.37		ng/L		84	70 - 145
Perfluorononanesulfonic acid (PFNS)	3.85	2.81		ng/L		73	65 - 145
Perfluorononanoic acid (PFNA)	4.00	3.41		ng/L		85	70 - 150
Perfluorooctanesulfonamide (PFOSA)	4.00	3.80		ng/L		95	70 - 145
Perfluorooctanesulfonic acid (PFOS)	3.72	5.56		ng/L		150	55 - 150
Perfluorooctanoic acid (PFOA)	4.00	3.47		ng/L		87	70 - 150

Method: 1633A - Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A (Continued)

Lab Sample ID: LLCS 410-710704/8-A
Matrix: Water
Analysis Batch: 710552

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 710704

Analyte	Spike Added	LLCS Result	LLCS Qualifier	Unit	D	%Rec	%Rec Limits
Perfluoropentanesulfonic acid (PFPeS)	3.76	3.50		ng/L		93	65 - 140
Perfluoropentanoic acid (PFPeA)	4.00	3.04		ng/L		76	65 - 135
Perfluorotetradecanoic acid (PFTeDA)	4.00	2.99		ng/L		75	60 - 140
Perfluorotridecanoic acid (PFTrDA)	4.00	3.21		ng/L		80	65 - 140
Perfluoroundecanoic acid (PFUnA)	4.00	3.33		ng/L		83	70 - 145

Isotope Dilution	*LLCS %Recovery*	*LLCS Qualifier*	*Limits*
13C2 4:2 FTS	*112*		*40 - 200*
13C2 6:2 FTS	*101*		*40 - 200*
13C2 8:2 FTS	*113*		*40 - 300*
13C2 PFDoA	*89.1*		*10 - 130*
13C2 PFTeDA	*73.8*		*10 - 130*
13C3 HFPO-DA	*84.6*		*40 - 130*
13C3 PFBS	*108*		*40 - 135*
13C3 PFHxS	*94.4*		*40 - 130*
13C4 PFBA	*101*		*5 - 130*
13C4 PFHpA	*93.5*		*40 - 130*
13C5 PFHxA	*100*		*40 - 130*
13C5 PFPeA	*101*		*40 - 130*
13C6 PFDA	*91.0*		*40 - 130*
13C7 PFUnA	*90.9*		*30 - 130*
13C8 FOSA	*87.3*		*40 - 130*
13C8 PFOA	*90.3*		*40 - 130*
13C8 PFOS	*99.2*		*40 - 130*
13C9 PFNA	*95.0*		*40 - 130*
d3-NMeFOSAA	*98.0*		*40 - 170*
d3-NMePFOSA	*78.3*		*10 - 130*
d5-NEtFOSAA	*83.9*		*25 - 135*
d5-NEtPFOSA	*77.3*		*10 - 130*
d7-N-MeFOSE-M	*67.4*		*10 - 130*
d9-N-EtFOSE-M	*55.8*		*10 - 130*

Method: 6010D - Metals (ICP)

Lab Sample ID: MB 480-758178/1-A
Matrix: Water
Analysis Batch: 758324

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 758178

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron	ND		0.050	0.019	mg/L		09/29/25 08:35	09/29/25 14:54	1

Method: 6010D - Metals (ICP) (Continued)

Lab Sample ID: LCS 480-758178/2-A
Matrix: Water
Analysis Batch: 758324

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 758178

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
Iron	5.10	5.35		mg/L		105	80 - 120

Lab Sample ID: LCSD 480-758178/3-A
Matrix: Water
Analysis Batch: 758324

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA
Prep Batch: 758178

Analyte	Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
Iron	5.10	5.38		mg/L		105	80 - 120	0	20

Lab Sample ID: MB 480-758180/1-A
Matrix: Water
Analysis Batch: 758350

Client Sample ID: Method Blank
Prep Type: Total Recoverable
Prep Batch: 758180

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Iron, Dissolved	ND		0.050	0.019	mg/L		09/29/25 08:34	09/29/25 23:53	1

Lab Sample ID: LCS 480-758180/2-A
Matrix: Water
Analysis Batch: 758350

Client Sample ID: Lab Control Sample
Prep Type: Total Recoverable
Prep Batch: 758180

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits
Iron, Dissolved	5.10	5.40		mg/L		106	80 - 120

Lab Sample ID: 480-232963-1 MS
Matrix: Water
Analysis Batch: 758350

Client Sample ID: MW1R
Prep Type: Dissolved
Prep Batch: 758180

Analyte	Sample Result	Sample Qualifier	Spike Added	MS Result	MS Qualifier	Unit	D	%Rec	%Rec Limits
Iron, Dissolved	0.21		5.10	5.55		mg/L		105	75 - 125

Lab Sample ID: 480-232963-1 MSD
Matrix: Water
Analysis Batch: 758350

Client Sample ID: MW1R
Prep Type: Dissolved
Prep Batch: 758180

Analyte	Sample Result	Sample Qualifier	Spike Added	MSD Result	MSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
Iron, Dissolved	0.21		5.10	5.59		mg/L		106	75 - 125	1	20

Method: 300.0 - Anions, Ion Chromatography

Lab Sample ID: MB 480-758298/28
Matrix: Water
Analysis Batch: 758298

Client Sample ID: Method Blank
Prep Type: Total/NA

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Sulfate	ND		2.0	0.35	mg/L			09/29/25 21:33	1

Method: 300.0 - Anions, Ion Chromatography (Continued)

Lab Sample ID: LCS 480-758298/29
Matrix: Water
Analysis Batch: 758298

Client Sample ID: Lab Control Sample
Prep Type: Total/NA

Analyte			Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits		
Sulfate			50.1	47.43		mg/L		95	90 - 110		

Lab Sample ID: 480-232963-4 MS
Matrix: Water
Analysis Batch: 758298

Client Sample ID: MW3I
Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MS Result	MS Qualifier	Unit	D	%Rec	%Rec Limits		
Sulfate	79.9		250	284.0		mg/L		82	80 - 120		

Lab Sample ID: MB 480-758402/4
Matrix: Water
Analysis Batch: 758402

Client Sample ID: Method Blank
Prep Type: Total/NA

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Sulfate	ND		2.0	0.35	mg/L			09/30/25 15:24	1

Lab Sample ID: LCS 480-758402/5
Matrix: Water
Analysis Batch: 758402

Client Sample ID: Lab Control Sample
Prep Type: Total/NA

Analyte			Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits		
Sulfate			50.1	47.13		mg/L		94	90 - 110		

GC/MS VOA

Analysis Batch: 758407

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-1	MW1R	Total/NA	Water	8260C	
480-232963-2	MW2R	Total/NA	Water	8260C	
480-232963-3	MW3S	Total/NA	Water	8260C	
480-232963-5	MW4S	Total/NA	Water	8260C	
480-232963-6	MW4I	Total/NA	Water	8260C	
480-232963-7	MW5S	Total/NA	Water	8260C	
480-232963-8	MW5I	Total/NA	Water	8260C	
480-232963-9	MW6	Total/NA	Water	8260C	
480-232963-10	MW7	Total/NA	Water	8260C	
480-232963-11	MW8	Total/NA	Water	8260C	
480-232963-12	MW11	Total/NA	Water	8260C	
480-232963-13	MW12S	Total/NA	Water	8260C	
480-232963-14	MW1C	Total/NA	Water	8260C	
480-232963-15	MW1D	Total/NA	Water	8260C	
480-232963-16	MW1E	Total/NA	Water	8260C	
480-232963-17	TB	Total/NA	Water	8260C	
MB 480-758407/8	Method Blank	Total/NA	Water	8260C	
LCS 480-758407/6	Lab Control Sample	Total/NA	Water	8260C	
480-232963-6 MS	MW4I	Total/NA	Water	8260C	
480-232963-6 MSD	MW4I	Total/NA	Water	8260C	

Analysis Batch: 758501

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-4	MW3I	Total/NA	Water	8260C	
MB 480-758501/8	Method Blank	Total/NA	Water	8260C	
LCS 480-758501/6	Lab Control Sample	Total/NA	Water	8260C	

GC VOA

Analysis Batch: 758460

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-1	MW1R	Total/NA	Water	RSK-175	
480-232963-1 - DL	MW1R	Total/NA	Water	RSK-175	
480-232963-2	MW2R	Total/NA	Water	RSK-175	
480-232963-2 - DL	MW2R	Total/NA	Water	RSK-175	
480-232963-3	MW3S	Total/NA	Water	RSK-175	
480-232963-4	MW3I	Total/NA	Water	RSK-175	
480-232963-5	MW4S	Total/NA	Water	RSK-175	
480-232963-5 - DL	MW4S	Total/NA	Water	RSK-175	
480-232963-6	MW4I	Total/NA	Water	RSK-175	
480-232963-6 - DL	MW4I	Total/NA	Water	RSK-175	
MB 480-758460/4	Method Blank	Total/NA	Water	RSK-175	
LCS 480-758460/5	Lab Control Sample	Total/NA	Water	RSK-175	
LCSD 480-758460/6	Lab Control Sample Dup	Total/NA	Water	RSK-175	

LCMS

Analysis Batch: 707164

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-3 - RE	MW3S	Total/NA	Water	1633A	707351
MB 410-707351/16-A	Method Blank	Total/NA	Water	1633A	707351
LCS 410-707351/17-A	Lab Control Sample	Total/NA	Water	1633A	707351

LCMS (Continued)

Analysis Batch: 707164 (Continued)

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
LCSD 410-707351/18-A	Lab Control Sample Dup	Total/NA	Water	1633A	707351

Prep Batch: 707351

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-3 - RE	MW3S	Total/NA	Water	1633A	
MB 410-707351/16-A	Method Blank	Total/NA	Water	1633A	
LCS 410-707351/17-A	Lab Control Sample	Total/NA	Water	1633A	
LCSD 410-707351/18-A	Lab Control Sample Dup	Total/NA	Water	1633A	

Prep Batch: 708885

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-7	MW5S	Total/NA	Water	1633A	
480-232963-8	MW5I	Total/NA	Water	1633A	
MB 410-708885/15-A	Method Blank	Total/NA	Water	1633A	
LCS 410-708885/16-A	Lab Control Sample	Total/NA	Water	1633A	
LCSD 410-708885/17-A	Lab Control Sample Dup	Total/NA	Water	1633A	
LLCS 410-708885/18-A	Lab Control Sample	Total/NA	Water	1633A	

Analysis Batch: 709341

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-7	MW5S	Total/NA	Water	1633A	708885
480-232963-8	MW5I	Total/NA	Water	1633A	708885
MB 410-708885/15-A	Method Blank	Total/NA	Water	1633A	708885
LCS 410-708885/16-A	Lab Control Sample	Total/NA	Water	1633A	708885
LCSD 410-708885/17-A	Lab Control Sample Dup	Total/NA	Water	1633A	708885
LLCS 410-708885/18-A	Lab Control Sample	Total/NA	Water	1633A	708885

Analysis Batch: 710552

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-3	MW3S	Total/NA	Water	1633A	710704
480-232963-4	MW3I	Total/NA	Water	1633A	710704
480-232963-5	MW4S	Total/NA	Water	1633A	710704
480-232963-6	MW4I	Total/NA	Water	1633A	710704
MB 410-710704/5-A	Method Blank	Total/NA	Water	1633A	710704
LCS 410-710704/6-A	Lab Control Sample	Total/NA	Water	1633A	710704
LCSD 410-710704/7-A	Lab Control Sample Dup	Total/NA	Water	1633A	710704
LLCS 410-710704/8-A	Lab Control Sample	Total/NA	Water	1633A	710704

Prep Batch: 710704

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-3	MW3S	Total/NA	Water	1633A	
480-232963-4	MW3I	Total/NA	Water	1633A	
480-232963-5	MW4S	Total/NA	Water	1633A	
480-232963-6	MW4I	Total/NA	Water	1633A	
MB 410-710704/5-A	Method Blank	Total/NA	Water	1633A	
LCS 410-710704/6-A	Lab Control Sample	Total/NA	Water	1633A	
LCSD 410-710704/7-A	Lab Control Sample Dup	Total/NA	Water	1633A	
LLCS 410-710704/8-A	Lab Control Sample	Total/NA	Water	1633A	

Metals

Prep Batch: 758178

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-1	MW1R	Total/NA	Water	3005A	
480-232963-2	MW2R	Total/NA	Water	3005A	
480-232963-3	MW3S	Total/NA	Water	3005A	
480-232963-4	MW3I	Total/NA	Water	3005A	
480-232963-5	MW4S	Total/NA	Water	3005A	
480-232963-6	MW4I	Total/NA	Water	3005A	
MB 480-758178/1-A	Method Blank	Total/NA	Water	3005A	
LCS 480-758178/2-A	Lab Control Sample	Total/NA	Water	3005A	
LCSD 480-758178/3-A	Lab Control Sample Dup	Total/NA	Water	3005A	

Prep Batch: 758180

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-1	MW1R	Dissolved	Water	3005A	
480-232963-2	MW2R	Dissolved	Water	3005A	
480-232963-3	MW3S	Dissolved	Water	3005A	
480-232963-4	MW3I	Dissolved	Water	3005A	
480-232963-5	MW4S	Dissolved	Water	3005A	
480-232963-6	MW4I	Dissolved	Water	3005A	
MB 480-758180/1-A	Method Blank	Total Recoverable	Water	3005A	
LCS 480-758180/2-A	Lab Control Sample	Total Recoverable	Water	3005A	
480-232963-1 MS	MW1R	Dissolved	Water	3005A	
480-232963-1 MSD	MW1R	Dissolved	Water	3005A	

Analysis Batch: 758324

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-1	MW1R	Total/NA	Water	6010D	758178
480-232963-2	MW2R	Total/NA	Water	6010D	758178
480-232963-3	MW3S	Total/NA	Water	6010D	758178
480-232963-4	MW3I	Total/NA	Water	6010D	758178
480-232963-5	MW4S	Total/NA	Water	6010D	758178
480-232963-6	MW4I	Total/NA	Water	6010D	758178
MB 480-758178/1-A	Method Blank	Total/NA	Water	6010D	758178
LCS 480-758178/2-A	Lab Control Sample	Total/NA	Water	6010D	758178
LCSD 480-758178/3-A	Lab Control Sample Dup	Total/NA	Water	6010D	758178

Analysis Batch: 758350

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-1	MW1R	Dissolved	Water	6010D	758180
480-232963-2	MW2R	Dissolved	Water	6010D	758180
480-232963-3	MW3S	Dissolved	Water	6010D	758180
480-232963-4	MW3I	Dissolved	Water	6010D	758180
480-232963-5	MW4S	Dissolved	Water	6010D	758180
480-232963-6	MW4I	Dissolved	Water	6010D	758180
MB 480-758180/1-A	Method Blank	Total Recoverable	Water	6010D	758180
LCS 480-758180/2-A	Lab Control Sample	Total Recoverable	Water	6010D	758180
480-232963-1 MS	MW1R	Dissolved	Water	6010D	758180
480-232963-1 MSD	MW1R	Dissolved	Water	6010D	758180

QC Association Summary

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

General Chemistry

Analysis Batch: 758298

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-1	MW1R	Total/NA	Water	300.0	
480-232963-2	MW2R	Total/NA	Water	300.0	
480-232963-3	MW3S	Total/NA	Water	300.0	
480-232963-4	MW3I	Total/NA	Water	300.0	
MB 480-758298/28	Method Blank	Total/NA	Water	300.0	
LCS 480-758298/29	Lab Control Sample	Total/NA	Water	300.0	
480-232963-4 MS	MW3I	Total/NA	Water	300.0	

Analysis Batch: 758402

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
480-232963-5	MW4S	Total/NA	Water	300.0	
480-232963-6	MW4I	Total/NA	Water	300.0	
MB 480-758402/4	Method Blank	Total/NA	Water	300.0	
LCS 480-758402/5	Lab Control Sample	Total/NA	Water	300.0	

Lab Chronicle

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW1R

Lab Sample ID: 480-232963-1

Date Collected: 09/24/25 14:30

Matrix: Water

Date Received: 09/26/25 10:00

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		5	758407	ATG	EET BUF	10/01/25 01:19
Total/NA	Analysis	RSK-175		1	758460	MAN	EET BUF	10/01/25 10:47
Total/NA	Analysis	RSK-175	DL	22	758460	MAN	EET BUF	10/01/25 12:40
Dissolved	Prep	3005A			758180	EMO	EET BUF	09/29/25 08:34
Dissolved	Analysis	6010D		1	758350	BMB	EET BUF	09/29/25 23:58
Total/NA	Prep	3005A			758178	EMO	EET BUF	09/29/25 08:35
Total/NA	Analysis	6010D		1	758324	MP	EET BUF	09/29/25 15:41
Total/NA	Analysis	300.0		5	758298	AF	EET BUF	09/30/25 01:42

Client Sample ID: MW2R

Lab Sample ID: 480-232963-2

Date Collected: 09/24/25 12:30

Matrix: Water

Date Received: 09/26/25 10:00

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		1	758407	ATG	EET BUF	10/01/25 01:42
Total/NA	Analysis	RSK-175		1	758460	MAN	EET BUF	10/01/25 11:06
Total/NA	Analysis	RSK-175	DL	22	758460	MAN	EET BUF	10/01/25 12:59
Dissolved	Prep	3005A			758180	EMO	EET BUF	09/29/25 08:34
Dissolved	Analysis	6010D		1	758350	BMB	EET BUF	09/30/25 00:15
Total/NA	Prep	3005A			758178	EMO	EET BUF	09/29/25 08:35
Total/NA	Analysis	6010D		1	758324	MP	EET BUF	09/29/25 15:43
Total/NA	Analysis	300.0		5	758298	AF	EET BUF	09/30/25 01:56

Client Sample ID: MW3S

Lab Sample ID: 480-232963-3

Date Collected: 09/24/25 14:15

Matrix: Water

Date Received: 09/26/25 10:00

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		2000	758407	ATG	EET BUF	10/01/25 02:05
Total/NA	Analysis	RSK-175		110	758460	MAN	EET BUF	10/01/25 11:25
Total/NA	Prep	1633A	RE		707351	W2FB	ELLE	10/01/25 08:35
Total/NA	Analysis	1633A	RE	1	707164	RPU6	ELLE	10/02/25 03:58
Total/NA	Prep	1633A			710704	W2FB	ELLE	10/08/25 08:36
Total/NA	Analysis	1633A		1	710552	DX7G	ELLE	10/09/25 03:28
Dissolved	Prep	3005A			758180	EMO	EET BUF	09/29/25 08:34
Dissolved	Analysis	6010D		1	758350	BMB	EET BUF	09/30/25 00:17
Total/NA	Prep	3005A			758178	EMO	EET BUF	09/29/25 08:35
Total/NA	Analysis	6010D		1	758324	MP	EET BUF	09/29/25 15:45
Total/NA	Analysis	300.0		10	758298	AF	EET BUF	09/30/25 02:11

Lab Chronicle

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW3I

Date Collected: 09/24/25 11:50
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-4

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		1	758501	ATG	EET BUF	10/01/25 16:02
Total/NA	Analysis	RSK-175		1	758460	MAN	EET BUF	10/01/25 11:44
Total/NA	Prep	1633A			710704	W2FB	ELLE	10/08/25 08:36
Total/NA	Analysis	1633A		1	710552	DX7G	ELLE	10/09/25 03:41
Dissolved	Prep	3005A			758180	EMO	EET BUF	09/29/25 08:34
Dissolved	Analysis	6010D		1	758350	BMB	EET BUF	09/30/25 00:19
Total/NA	Prep	3005A			758178	EMO	EET BUF	09/29/25 08:35
Total/NA	Analysis	6010D		1	758324	MP	EET BUF	09/29/25 15:47
Total/NA	Analysis	300.0		5	758298	AF	EET BUF	09/30/25 02:26

Client Sample ID: MW4S

Date Collected: 09/24/25 15:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-5

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		2000	758407	ATG	EET BUF	10/01/25 02:51
Total/NA	Analysis	RSK-175		1	758460	MAN	EET BUF	10/01/25 12:03
Total/NA	Analysis	RSK-175	DL	44	758460	MAN	EET BUF	10/01/25 13:18
Total/NA	Prep	1633A			710704	W2FB	ELLE	10/08/25 08:36
Total/NA	Analysis	1633A		1	710552	DX7G	ELLE	10/09/25 03:55
Dissolved	Prep	3005A			758180	EMO	EET BUF	09/29/25 08:34
Dissolved	Analysis	6010D		1	758350	BMB	EET BUF	09/30/25 00:24
Total/NA	Prep	3005A			758178	EMO	EET BUF	09/29/25 08:35
Total/NA	Analysis	6010D		1	758324	MP	EET BUF	09/29/25 15:49
Total/NA	Analysis	300.0		20	758402	AF	EET BUF	09/30/25 16:37

Client Sample ID: MW4I

Date Collected: 09/24/25 14:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-6

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		40	758407	ATG	EET BUF	10/01/25 03:14
Total/NA	Analysis	RSK-175		1	758460	MAN	EET BUF	10/01/25 12:21
Total/NA	Analysis	RSK-175	DL	22	758460	MAN	EET BUF	10/01/25 13:56
Total/NA	Prep	1633A			710704	W2FB	ELLE	10/08/25 08:36
Total/NA	Analysis	1633A		1	710552	DX7G	ELLE	10/09/25 04:08
Dissolved	Prep	3005A			758180	EMO	EET BUF	09/29/25 08:34
Dissolved	Analysis	6010D		1	758350	BMB	EET BUF	09/30/25 00:22
Total/NA	Prep	3005A			758178	EMO	EET BUF	09/29/25 08:35
Total/NA	Analysis	6010D		1	758324	MP	EET BUF	09/29/25 15:52
Total/NA	Analysis	300.0		5	758402	AF	EET BUF	09/30/25 16:52

Lab Chronicle

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW5S
Date Collected: 09/24/25 09:50
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-7
Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		10	758407	ATG	EET BUF	10/01/25 03:38
Total/NA	Prep	1633A			708885	V3FW	ELLE	10/03/25 13:46
Total/NA	Analysis	1633A		1	709341	DX7G	ELLE	10/06/25 06:17

Client Sample ID: MW5I
Date Collected: 09/24/25 10:40
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-8
Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		1	758407	ATG	EET BUF	10/01/25 04:01
Total/NA	Prep	1633A			708885	V3FW	ELLE	10/03/25 13:46
Total/NA	Analysis	1633A		1	709341	DX7G	ELLE	10/06/25 06:31

Client Sample ID: MW6
Date Collected: 09/23/25 14:20
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-9
Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		1	758407	ATG	EET BUF	10/01/25 04:24

Client Sample ID: MW7
Date Collected: 09/23/25 15:35
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-10
Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		1	758407	ATG	EET BUF	10/01/25 04:47

Client Sample ID: MW8
Date Collected: 09/24/25 10:30
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-11
Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		1	758407	ATG	EET BUF	10/01/25 05:10

Client Sample ID: MW11
Date Collected: 09/24/25 16:05
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-12
Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		1	758407	ATG	EET BUF	10/01/25 05:33

Lab Chronicle

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Client Sample ID: MW12S
Date Collected: 09/24/25 16:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-13
Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		1	758407	ATG	EET BUF	10/01/25 05:56

Client Sample ID: MW1C
Date Collected: 09/25/25 12:15
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-14
Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		1	758407	ATG	EET BUF	10/01/25 06:19

Client Sample ID: MW1D
Date Collected: 09/25/25 09:15
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-15
Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		20	758407	ATG	EET BUF	10/01/25 06:43

Client Sample ID: MW1E
Date Collected: 09/25/25 11:05
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-16
Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		100	758407	ATG	EET BUF	10/01/25 07:06

Client Sample ID: TB
Date Collected: 09/25/25 00:00
Date Received: 09/26/25 10:00

Lab Sample ID: 480-232963-17
Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	8260C		1	758407	ATG	EET BUF	10/01/25 07:29

Laboratory References:

EET BUF = Eurofins Buffalo, 10 Hazelwood Drive, Amherst, NY 14228-2298, TEL (716)691-2600

ELLE = Eurofins Lancaster Laboratories Environment Testing, LLC, 2425 New Holland Pike, Lancaster, PA 17601, TEL (717)656-2300

Accreditation/Certification Summary

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Laboratory: Eurofins Buffalo

The accreditations/certifications listed below are applicable to this report.

Authority	Program	Identification Number	Expiration Date
New York	NELAP	10026	10-15-25

Laboratory: Eurofins Lancaster Laboratories Environment Testing, LLC

All accreditations/certifications held by this laboratory are listed. Not all accreditations/certifications are applicable to this report.

Authority	Program	Identification Number	Expiration Date
A2LA	Dept. of Defense ELAP	0001.01	11-30-26
A2LA	Dept. of Energy	0001.01	11-30-26
A2LA	ISO/IEC 17025	0001.01	11-30-26
Alabama	State	43200	01-31-26
Alaska	State	PA00009	06-30-26
Alaska (UST)	State	17-027	12-30-26
Arizona	State	AZ0780	03-12-26
Arkansas DEQ	State	88-00660	08-09-26
California	State	2792	01-31-26
Colorado	State	PA00009	06-30-26
Connecticut	State	PH-0746	06-30-27
DE Haz. Subst. Cleanup Act (HSCA)	State	019-006 (PA cert)	01-31-26
Delaware (DW)	State	N/A	01-31-26
Florida	NELAP	E87997	07-01-26
Georgia (DW)	State	C048	01-31-26
Illinois	NELAP	200027	01-31-26
Iowa	State	361	03-01-26
Kansas	NELAP	E-10151	10-31-25
Kentucky (DW)	State	KY90088	12-31-25
Kentucky (UST)	State	0001.01	11-30-26
Kentucky (WW)	State	KY90088	12-31-25
Louisiana (All)	NELAP	02055	06-30-26
Maine	State	2019012	03-12-27
Maryland	State	100	06-30-26
Massachusetts	State	M-PA009	06-30-26
Michigan	State	9930	01-31-26
Minnesota	NELAP	042-999-487	12-31-25
Mississippi	State	023	01-31-26
Missouri	State	450	01-31-28
Montana (DW)	State	0098	01-01-26
Nebraska	State	NE-OS-32-17	01-31-26
New Hampshire	NELAP	2730	01-10-26
New Jersey	NELAP	PA011	06-30-26
New York	NELAP	10670	10-15-25
North Carolina (DW)	State	42705	07-31-26
North Carolina (WW/SW)	State	521	12-31-25
North Dakota	State	R-205	01-31-24 *
Ohio	State	87787	01-31-26
Oklahoma	NELAP	9804	12-31-25
Oregon	NELAP	PA200001	09-12-26
Pennsylvania	NELAP	36-00037	01-31-26
Quebec Ministry of Environment and Fight against Climate Change	PALA	507	09-16-29
Rhode Island	State	LAO00338	12-30-25
South Carolina	State	89002	01-31-26

* Accreditation/Certification renewal pending - accreditation/certification considered valid.

Accreditation/Certification Summary

Client: GEI Consultants Inc
Job ID: 480-232963-1
Project/Site: Unifirst – Liverpool, NY

Laboratory: Eurofins Lancaster Laboratories Environment Testing, LLC (Continued)

All accreditations/certifications held by this laboratory are listed. Not all accreditations/certifications are applicable to this report.

Authority	Program	Identification Number	Expiration Date
Tennessee	State	02838	01-31-26
Texas	NELAP	T104704194-23-46	08-31-26
USDA	US Federal Programs	525-22-298-19481	10-25-25
Vermont	State	VT - 36037	10-27-25
Virginia	NELAP	460182	06-14-26
Washington	State	C457	04-11-26
West Virginia (DW)	State	9906 C	03-31-26
West Virginia DEP	State	055	07-31-26
Wyoming	State	8TMS-L	01-31-26
Wyoming (UST)	A2LA	0001.01	11-30-26

Method Summary

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Method	Method Description	Protocol	Laboratory
8260C	Volatile Organic Compounds by GC/MS	SW846	EET BUF
RSK-175	Dissolved Gases (GC)	RSK	EET BUF
1633A	Per- and Polyfluoroalkyl Substances by LC/MS/MS, 1633A	EPA	ELLE
6010D	Metals (ICP)	SW846	EET BUF
300.0	Anions, Ion Chromatography	EPA	EET BUF
1633A	Solid-Phase Extraction (SPE)	EPA	ELLE
3005A	Preparation, Total Metals	SW846	EET BUF
3005A	Preparation, Total Recoverable or Dissolved Metals	SW846	EET BUF
5030C	Purge and Trap	SW846	EET BUF

Protocol References:

EPA = US Environmental Protection Agency

RSK = Sample Prep And Calculations For Dissolved Gas Analysis In Water Samples Using A GC Headspace Equilibration Technique, RSKSOP-175, Rev. 0, 8/11/94, USEPA Research Lab

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

Laboratory References:

EET BUF = Eurofins Buffalo, 10 Hazelwood Drive, Amherst, NY 14228-2298, TEL (716)691-2600

ELLE = Eurofins Lancaster Laboratories Environment Testing, LLC, 2425 New Holland Pike, Lancaster, PA 17601, TEL (717)656-2300

Sample Summary

Client: GEI Consultants Inc
Project/Site: Unifirst – Liverpool, NY

Job ID: 480-232963-1

Lab Sample ID	Client Sample ID	Matrix	Collected	Received	Sample Origin
480-232963-1	MW1R	Water	09/24/25 14:30	09/26/25 10:00	New York
480-232963-2	MW2R	Water	09/24/25 12:30	09/26/25 10:00	New York
480-232963-3	MW3S	Water	09/24/25 14:15	09/26/25 10:00	New York
480-232963-4	MW3I	Water	09/24/25 11:50	09/26/25 10:00	New York
480-232963-5	MW4S	Water	09/24/25 15:00	09/26/25 10:00	New York
480-232963-6	MW4I	Water	09/24/25 14:00	09/26/25 10:00	New York
480-232963-7	MW5S	Water	09/24/25 09:50	09/26/25 10:00	New York
480-232963-8	MW5I	Water	09/24/25 10:40	09/26/25 10:00	New York
480-232963-9	MW6	Water	09/23/25 14:20	09/26/25 10:00	New York
480-232963-10	MW7	Water	09/23/25 15:35	09/26/25 10:00	New York
480-232963-11	MW8	Water	09/24/25 10:30	09/26/25 10:00	New York
480-232963-12	MW11	Water	09/24/25 16:05	09/26/25 10:00	New York
480-232963-13	MW12S	Water	09/24/25 16:00	09/26/25 10:00	New York
480-232963-14	MW1C	Water	09/25/25 12:15	09/26/25 10:00	New York
480-232963-15	MW1D	Water	09/25/25 09:15	09/26/25 10:00	New York
480-232963-16	MW1E	Water	09/25/25 11:05	09/26/25 10:00	New York
480-232963-17	TB	Water	09/25/25 00:00	09/26/25 10:00	New York

Amherst, NY 14228-2298
Phone: 716-691-2600 Fax: 716-691-7991

Chain of Custody Record

eurofins Environment Testing

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Client Information	Sampler: J. DesRosiers	Lab PM: Schove, John R	Carrier Tracking No(s):	COC No: 480-207985-42102.1
Client Contact: Ms. Abigal Jock	Phone: 607-216-8985	E-Mail: John.Schove@et.eurofinsus.com	State of Origin: NY	Page: Page 1 of 2
Company: GEI Consultants Inc	PWSID:	Analysis Requested		Job #
Address: 950 Danby Road Suite 201-F	Due Date Requested: Standard			Preservation Codes: A - HCL, N - None, D - HNO3
City: Ithaca	TAT Requested (days):			
State, Zip: NY, 14850	Compliance Project: Δ Yes Δ No			
Phone: 607-216-8983(Tel)	PO #: Purchase Order Requested			
Email: AJock@geiconsultants.com NMoore@geiconsultants.com	WO #: 2001642			
Project Name: Unifirst – Liverpool, NY	Project #: 48029086			
Site: UniFirst Liverpool	SSOW#:			Other:

480-232963 Chain of Custody

3 coolers total

Sample Identification	Sample Date	Sample Time	Sample Type (C=Comp, G=grab)	Matrix (W=water, S=solid, O=wasteoil, BT=Tissue, A=Air, DW=Drinking Water)	Field Filtered Sample (Yes or No)	Perform MS/MSD (Yes or No)	8260C - TCL VOCs	RSK_175 - Methane, Ethane, Ethene	300.0_28D - Sulfate	6010D - Metals - Iron (Fe), Total	6010D - Metals - Iron (Fe), Dissolved	1633A - PFAS - NY Standard 40 Analytes	Total Number of containe	Special Instructions/Note:
Preservation Code:							A	A	N	D	D	N		
MW1R	9/24/25	14:30		Water	Y	N	X	X	X	X	X		9	
MW2R	9/24/25	12:30		Water	Y	N	X	X	X	X	X		9	
MW3S	9/24/25	14:15		Water	Y	N	X	X	X	X	X	X	11	
MW3I	9/24/25	11:50		Water	Y	N	X	X	X	X	X	X	11	
MW4S	9/24/25	16:00		Water	Y	N	X	X	X	X	X	X	11	
MW4I	9/24/25	14:00		Water	Y	N	X	X	X	X	X	X	11	
MW5S	9/24/25	9:50		Water	N	N	X					X	5	
MW5I	9/24/25	10:40		Water	N	N	X					X	5	
MW6	9/23/25	14:20		Water	N	N	X						3	
MW7	9/23/25	15:35		Water	N	N	X						3	
MW8	9/24/25	10:30		Water	N	N	X						3	

Possible Hazard Identification
☐ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown ☐ Radiological

Sample Disposal (A fee may be assessed if samples are retained longer than 1 month)
☐ Return To Client ☒ Disposal By Lab ☐ Archive For _____ Months

Deliverable Requested: I, II, III, IV, Other (specify)

Special Instructions/QC Requirements:

Empty Kit Relinquished by:	Date:	Time:	Method of Shipment:

	Relinquished by	Date/Time	Company	Received by	Date/Time	Company
1	[signature]	9/25/25 14:50	GEI	[signature]	9/25/25 1450	
2	[signature]	9/25/25 1600		[signature]	9/26/25 1000	EUROFINS
3						

Custody Seals Intact: Δ Yes Δ No	Custody Seal No.:	Cooler Temperature(s) °C and Other Remarks: 2.9, 3.2 IR #SC

Ver: 10/10/2024

Amherst, NY 14228-2298
Phone: 716-691-2600 Fax: 716-691-7991

Chain of Custody Record — Syracuse # 225

eurofins | Environment Testing

Client Information	Sampler: J. DesRosiers	Lab PM: Schove, John R	Carrier Tracking No(s):	COC No: 480-207985-42102.2
Client Contact: Ms. Abigal Jock	Phone: 607-216-8985	E-Mail: John.Schove@et.eurofinsus.com	State of Origin: NY	Page: Page 2 of 2
Company: GEI Consultants Inc	PWSID:	Analysis Requested		Job #:
Address: 950 Danby Road Suite 201-F	Due Date Requested: Standard			Preservation Codes: A - HCL, N - None, D - HNO3
City: Ithaca	TAT Requested (days):			
State, Zip: NY, 14850	Compliance Project: Δ Yes Δ No			
Phone: 607-216-8983(Tel)	PO #: Purchase Order Requested			
Email: AJock@geiconsultants.com WMoore@geiconsultants.com	WO #: 2001642			
Project Name: Unifirst – Liverpool, NY	Project #: 48029086			Other:
Site: Uni First Liverpool	SSOW#:			

Sample Identification	Sample Date	Sample Time	Sample Type (C=comp, G=grab)	Matrix (W=water, S=solid, O=waste/oil, BT=Tissue, A=Air, DW=Drinking Water)	Field Filtered Sample (Yes or No)	Perform MS/MSD (Yes or No)	8260C - TCL VOCs	RSK_175 - Methane, Ethane, Ethene	300.0_28D - Sulfate	6010D - Metals - Iron (Fe), Total	6010D - Metals - Iron (Fe), Dissolved	1633A - PFAS - NY Standard 40 Analytes	Total Number of containers	Special Instructions/Note:
			Preservation Code:				A	A	N	D	D	N		3 coolers total
MW11	9/24/25	16:05		Water	N	N	X						2 (MS/MSD)	
MW12S	9/24/25	16:00		Water	N	N	X							
MW1C	9/25/25	12:15		Water	N	N	X							
MW1D	9/25/25	9:15		Water	N	N	X							
MW1E	9/25/25	11:05		Water	N	N	X							
TB	9/25/25			Water			X						2	
				~~Water~~										
~~MW1R~~				~~Water~~										
~~MW2R~~				~~Water~~										
~~MW3S~~				~~Water~~										
~~MW3I~~				~~Water~~										

Possible Hazard Identification
☐ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown ☐ Radiological

Sample Disposal (A fee may be assessed if samples are retained longer than 1 month)
☐ Return To Client ☒ Disposal By Lab ☐ Archive For _____ Months

Deliverable Requested: I, II, III, IV, Other (specify)

Special Instructions/QC Requirements:

Empty Kit Relinquished by:	Date:	Time:	Method of Shipment:

	Relinquished by	Date/Time	Company	Received by	Date/Time	Company
1	Jordan DesRosiers	9/25/25 14:50			9/25/25 1450	
2		9/25/25 1900			9/26/25 1000	GURO
3		Date/Time:	Company	Received by:	Date/Time:	Company

Custody Seals Intact: Δ Yes Δ No | Custody Seal No.:

Cooler Temperature(s) °C and Other Remarks:

Ver: 10/10/2024

Eurofins Buffalo
10 Hazelwood Drive
Amherst, NY 14228-2298
Phone: 716-691-2600 Fax: 716-691-7991

Chain of Custody Record

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Client Information (Sub Contract Lab)	Sampler: N/A	Lab PM: Schove, John R	Carrier Tracking No(s): N/A	COC No: 480-95087.1
Client Contact: Shipping/Receiving	Phone: N/A	E-Mail: John.Schove@et eurofinsus com	State of Origin: New York	Page: Page 1 of 1
Company: Eurofins Lancaster Laboratories Environm		Accreditations Required (See note): NELAP - New York		Job #: 480-232963-1
Address: 2425 New Holland Pike, ,	Due Date Requested: 10/16/2025			
City: Lancaster	TAT Requested (days): N/A			
State, Zip: PA, 17601				
Phone: 717-656-2300(Tel)	PO #: N/A			
Email: N/A	WO #: N/A			
Project Name: Unifirst – Liverpool, NY	Project #: 48029086			
Site: N/A	SSOW#: N/A			

Analysis Requested

Preservation Codes:
-

Other:
N/A

Sample Identification - Client ID (Lab ID)	Sample Date	Sample Time	Sample Type (C=comp, G=grab)	Matrix (W=water, S=solid, O=waste/oil, BT=Tissue, A=Air)	Field Filtered Sample (Yes or No)	Perform MS/MSD (Yes or No)	1633A/1633A_SPEPFAS - NY Standard 40 Analytes	Total Number of containers	Special Instructions/Note:
			Preservation Code:						
MW3S (480-232963-3)	9/24/25	14:15 Eastern	G	Water			X	2	
MW3I (480-232963-4)	9/24/25	11:50 Eastern	G	Water			X	2	
MW4S (480-232963-5)	9/24/25	15:00 Eastern	G	Water			X	2	
MW4I (480-232963-6)	9/24/25	14:00 Eastern	G	Water			X	2	
MW5S (480-232963-7)	9/24/25	09:50 Eastern	G	Water			X	2	
MW5I (480-232963-8)	9/24/25	10:40 Eastern	G	Water			X	2	

Note: Since laboratory accreditations are subject to change, Eurofins Environment Testing Northeast, LLC places the ownership of method, analyte & accreditation compliance upon our subcontract laboratories. This sample shipment is forwarded under chain-of-custody. If the laboratory does not currently maintain accreditation in the State of Origin listed above for analysis/tests/matrix being analyzed, the samples must be shipped back to the Eurofins Environment Testing Northeast, LLC laboratory or other instructions will be provided. Any changes to accreditation status should be brought to Eurofins Environment Testing Northeast, LLC attention immediately. If all requested accreditations are current to date, return the signed Chain of Custody attesting to said compliance to Eurofins Environment Testing Northeast, LLC

Possible Hazard Identification	Sample Disposal (A fee may be assessed if samples are retained longer than 1 month)
Unconfirmed	☐ Return To Client ☐ Disposal By Lab ☐ Archive For _____ Months
Deliverable Requested: I, II, III, IV, Other (specify) Primary Deliverable Rank: 2	Special Instructions/QC Requirements:

Empty Kit Relinquished by:	Date:	Time:	Method of Shipment:

Relinquished by	Date/Time	Company	Received by	Date/Time	Company
VHS	9-26-25 1600				
Relinquished by	Date/Time	Company	Received by	Date/Time	Company
Relinquished by	Date/Time	Company	Received by	Date/Time	Company
				09/27/25 1100	

Custody Seals Intact: Δ Yes Δ No	Custody Seal No.:	Cooler Temperature(s) °C and Other Remarks: R:2.0 C:1.9

Ver: 10/10/2024

Login Sample Receipt Checklist

Client: GEI Consultants Inc

Job Number: 480-232963-1

Login Number: 232963
List Number: 1
Creator: Yeager, Brian A

List Source: Eurofins Buffalo

Question	Answer	Comment
Radioactivity either was not measured or, if measured, is at or below background	True	
The cooler's custody seal, if present, is intact.	True	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	2.9,3.2 ICE IR# SC
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
Is the Field Sampler's name present on COC?	True	
There are no discrepancies between the sample IDs on the containers and the COC.	True	
Samples are received within Holding Time (Excluding tests with immediate HTs)..	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
Sample Preservation Verified	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
VOA sample vials do not have headspace or bubble is <6mm (1/4") in diameter.	True	
If necessary, staff have been informed of any short hold time or quick TAT needs	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	
Sampling Company provided.	True	GEI
Samples received within 48 hours of sampling.	False	
Samples requiring field filtration have been filtered in the field.	True	
Chlorine Residual checked.	N/A	

Login Sample Receipt Checklist

Client: GEI Consultants Inc

Job Number: 480-232963-1

Login Number: 232963
List Number: 2
Creator: Yeager, Brian A

List Source: Eurofins Buffalo

Question	Answer	Comment
Radioactivity wasn't checked or is </= background as measured by a survey meter.	True	
The cooler's custody seal, if present, is intact.	True	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
Is the Field Sampler's name present on COC?	True	
There are no discrepancies between the sample IDs on the containers and the COC.	True	
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
Sample Preservation Verified	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
Containers requiring zero headspace have no headspace or bubble is <6mm (1/4").	True	
If necessary, staff have been informed of any short hold time or quick TAT needs	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	N/A	

Login Sample Receipt Checklist

Client: GEI Consultants Inc

Job Number: 480-232963-1

Login Number: 232963
List Number: 3
Creator: Reynolds, Benjamin

List Source: Eurofins Lancaster Laboratories Environment Testing, LLC
List Creation: 09/27/25 05:03 PM

Question	Answer	Comment
The cooler's custody seal is intact.	True	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature acceptable,where thermal pres is required(</=6C, not frozen).	True	
Cooler Temperature is recorded.	True	
WV:Container Temp acceptable,where thermal pres is required (</=6C, not frozen).	N/A	
WV: Container Temperature is recorded.	N/A	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
There are no discrepancies between the containers received and the COC.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
There is sufficient vol. for all requested analyses.	True	
Is the Field Sampler's name present on COC?	False	Received project as a subcontract.
Sample custody seals are intact.	N/A	
VOA sample vials do not have headspace >6mm in diameter (none, if from WV)?	N/A	

Appendix C Microbial Insights QuantArray Chlor Report

10515 Research Drive
Knoxville, TN 37932
Phone: 865.573.8188
Fax: 865.573.8133
Web: www.microbe.com

SITE LOGIC Report

QuantArray®-Chlor Study

Contact:	Wendy Moore	Phone:	
Address:	GEI Consultants 950 Danby Road Suite 201-F Ithaca, NY 14850	Email:	wmoore@geiconsultants.com

MI Identifier: 073WI

Report Date: 09/30/2025

Project: UniFirst-Liverpool, NY

Reviewed By: *Sarah Keys*

Comments:

The QuantArray®-Chlor Approach

Quantification *Dehalococcoides*, the only known bacterial group capable of complete reductive dechlorination of PCE and TCE to ethene, has become an indispensable component of assessment, remedy selection, and performance monitoring at sites impacted by chlorinated solvents. While undeniably a key group of halorespiring bacteria, *Dehalococcoides* are not the only bacteria of interest in the subsurface because reductive dechlorination is not the only potential biodegradation pathway operative at contaminated sites, and chlorinated ethenes are not always the primary contaminants of concern. The QuantArray®-Chlor not only includes a variety of halorespiring bacteria (*Dehalococcoides*, *Dehalobacter*, *Dehalogenimonas*, etc.) to assess the potential for reductive dechlorination of chloroethenes, chloroethanes, chlorobenzenes, chlorophenols, and chloroform, but also provides quantification of functional genes involved in aerobic (co)metabolic pathways for biodegradation of chlorinated solvents and even competing biological processes. Thus, the QuantArray®-Chlor will give site managers the ability to simultaneously yet economically evaluate the potential for biodegradation of a spectrum of common chlorinated contaminants through a multitude of anaerobic and aerobic (co)metabolic pathways to give a much more clear and comprehensive view of contaminant biodegradation.

The QuantArray®-Chlor is used to quantify specific microorganisms and functional genes to evaluate the following:

Anaerobic Reductive Dechlorination	Quantification of important halorespiring bacteria (e.g. *Dehalococcoides*, *Dehalobacter*, *Dehalogenimonas*, *Desulfitobacterium* spp.) and key functional genes (e.g. vinyl chloride reductases, TCE reductase, chloroform reductase) responsible for reductive dechlorination of a broad spectrum of chlorinated solvents.
Aerobic Cometabolism	Several different types of bacteria including methanotrophs and some toluene/phenol utilizing bacteria can co-oxidize TCE, DCE, and vinyl chloride. The QuantArray®-Chlor quantifies functional genes like soluble methane monooxygenase encoding enzymes capable of co-oxidation of chlorinated ethenes.
Aerobic (Co)metabolism of Vinyl Chloride	Ethene oxidizing bacteria are capable of cometabolism of vinyl chloride. In some cases, ethenotrophs can also utilize vinyl chloride as a growth supporting substrate. The QuantArray®-Chlor targets key functional genes in ethene metabolism.

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Knoxville, TN 37932
Phone: 865.573.8188
Fax: 865.573.8133
Web: www.microbe.com

Results

Table 1: Summary of the QuantArray®-Chlor results obtained for samples MW3S, MW3I, MW4S, and MW4I

Sample Name Sample Date	MW3S 09/24/2025	MW3I 09/24/2025	MW4S 09/24/2025	MW4I 09/24/2025
Reductive Dechlorination	cells/mL	cells/mL	cells/mL	cells/mL
Dehalococcoides (DHC)	**3.10E+06**	**4.80E+02**	**1.00E+02**	**2.80E+03**
tceA Reductase (TCE)	**1.60E+06**	**2.60E+01**	**9.30E+00**	**1.50E-01 (J)**
BAV1 Vinyl Chloride Reductase (BVC)	**2.40E+05**	**1.70E+02**	**1.10E+01**	**2.00E+03**
Vinyl Chloride Reductase (VCR)	**2.40E+05**	**1.80E+02**	**1.50E+01**	**4.80E+02**
Dehalobacter spp. (DHBt)	**2.70E+04**	**2.10E+02**	**2.40E+03**	**3.80E+02**
Dehalobacter DCM (DCM)	**1.10E+03**	<2.50E-01	<4.50E-01	<2.50E-01
Dehalogenimonas spp. (DHG)	**1.10E+06**	**3.00E+03**	**2.00E+03**	**1.60E+04**
cerA Reductase (CER)	**3.10E+03**	<2.50E-01	<4.50E-01	**1.60E+01**
trans-1,2-DCE Reductase (TDR)	<8.30E-01	<2.50E-01	<4.50E-01	<2.50E-01
Desulfitobacterium spp. (DSB)	**3.50E+04**	**2.40E+02**	**9.20E+03**	**1.20E+03**
Dehalobium chlorocoercia (DECO)	**6.10E+02**	**1.40E+02**	**5.20E+03**	**4.60E+02**
Desulfuromonas spp. (DSM)	**8.80E+00**	**2.20E+01**	**2.90E+01**	<2.50E-01
PCE Reductase (PCE-1)	<8.30E-01	<2.50E-01	<4.50E-01	<2.50E-01
PCE Reductase (PCE-2)	<8.30E-01	<2.50E-01	<4.50E-01	<2.50E-01
Chloroform Reductase (CFR)	<8.30E-01	<2.50E-01	<4.50E-01	<2.50E-01
1,1 DCA Reductase (DCA)	<8.30E-01	<2.50E-01	<4.50E-01	<2.50E-01
1,2 DCA Reductase (DCAR)	<8.30E-01	<2.50E-01	<4.50E-01	<2.50E-01
Aerobic (Co)Metabolic				
Soluble Methane Monooxygenase (SMMO)	<8.30E-01	**6.60E+01**	**1.80E+02**	<2.50E-01
Toluene Dioxygenase (TOD)	<8.30E-01	**2.80E+00**	<4.50E-01	**9.90E+01**
Phenol Hydroxylase (PHE)	**6.40E+03**	<2.50E-01	**4.30E+03**	**4.10E+01**
Trichlorobenzene Dioxygenase (TCBO)	<8.30E-01	<2.50E-01	<4.50E-01	<2.50E-01
Toluene Monooxygenase 2 (RDEG)	**2.00E+03**	<2.50E-01	**6.30E+02**	<2.50E-01
Toluene Monooxygenase (RMO)	**3.80E+04**	<2.50E-01	<4.50E-01	**2.00E+03**
Ethene Monooxygenase (EtnC)	**2.50E+04**	<2.50E-01	**4.60E+00**	<2.50E-01
Epoxyalkane Transferase (EtnE)	**3.00E+05**	<2.50E-01	**2.70E+03**	<2.50E-01
Dichloromethane Dehalogenase (DCMA)	<8.30E-01	<2.50E-01	<4.50E-01	<2.50E-01
Other				
Total Eubacteria (EBAC)	**1.03E+07**	**7.50E+04**	**1.50E+06**	**2.30E+05**
Sulfate Reducing Bacteria (APS)	**7.20E+04**	**6.40E+04**	**2.10E+05**	**5.60E+04**
Methanogens (MGN)	**3.20E+03**	**3.90E+01**	**3.70E+02**	**2.90E+01**

Legend:

NA = Not Analyzed
NS = Not Sampled
J = Estimated Gene Copies Below PQL but Above LQL
I = Inhibited
< = Result Not Detected

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Fax: 865.573.8133
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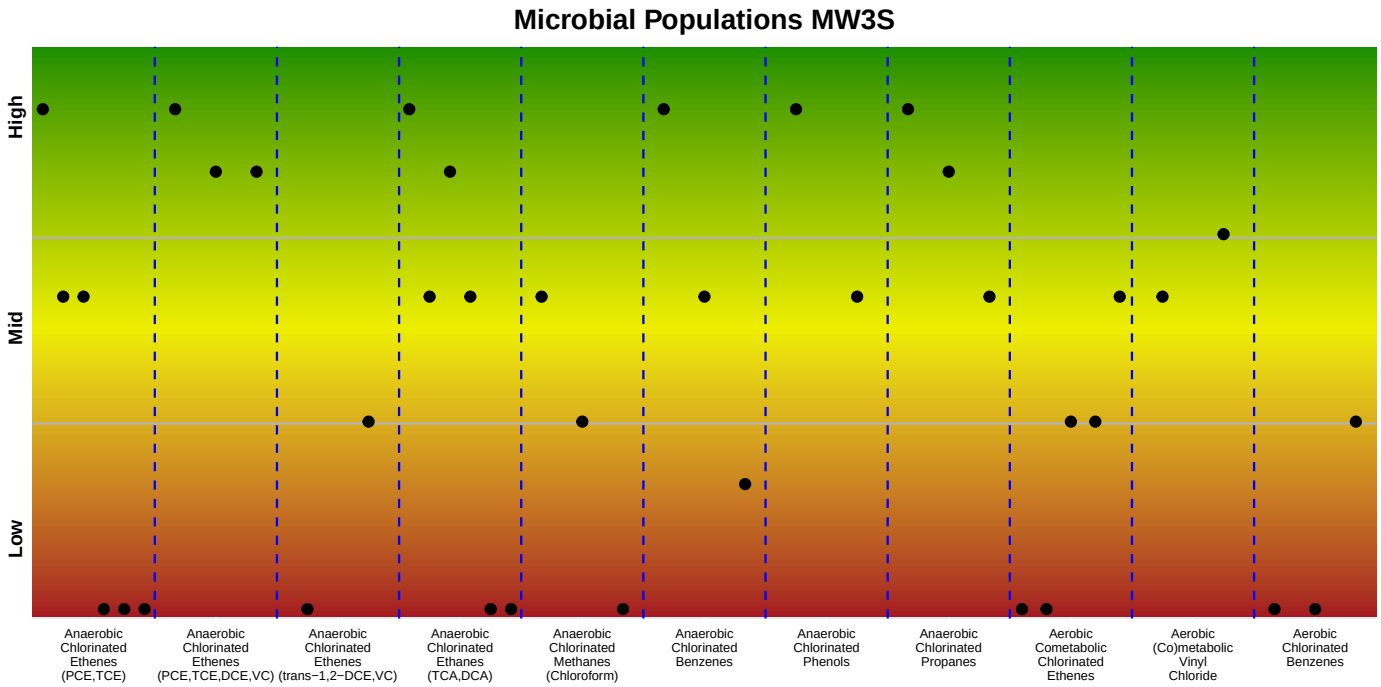

Figure 1: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants. See below to understand the correlation between contaminants and their associated qPCR assays.

Anaerobic - Reductive Dechlorination or Dichloroelimination	
Contaminants	**qPCR Assays**
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER
Chlorinated Ethanes (TCA and 1,2- DCA)	DHC, DHBt, DHG, DSB[1], DCA, DCAR
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR
Chlorinated Benzenes	DHC, DHBt[2], DECO
Chlorinated Phenols	DHC, DSB
Chlorinated Propanes	DHC, DHG, DSB[1]
Aerobic - (Co)metabolism	
Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Benzenes	TOD, TCBO, PHE

[1] *Desulfitobacterium dichloroeliminans* DCA1. [2] Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

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Phone: 865.573.8188
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Web: www.microbe.com

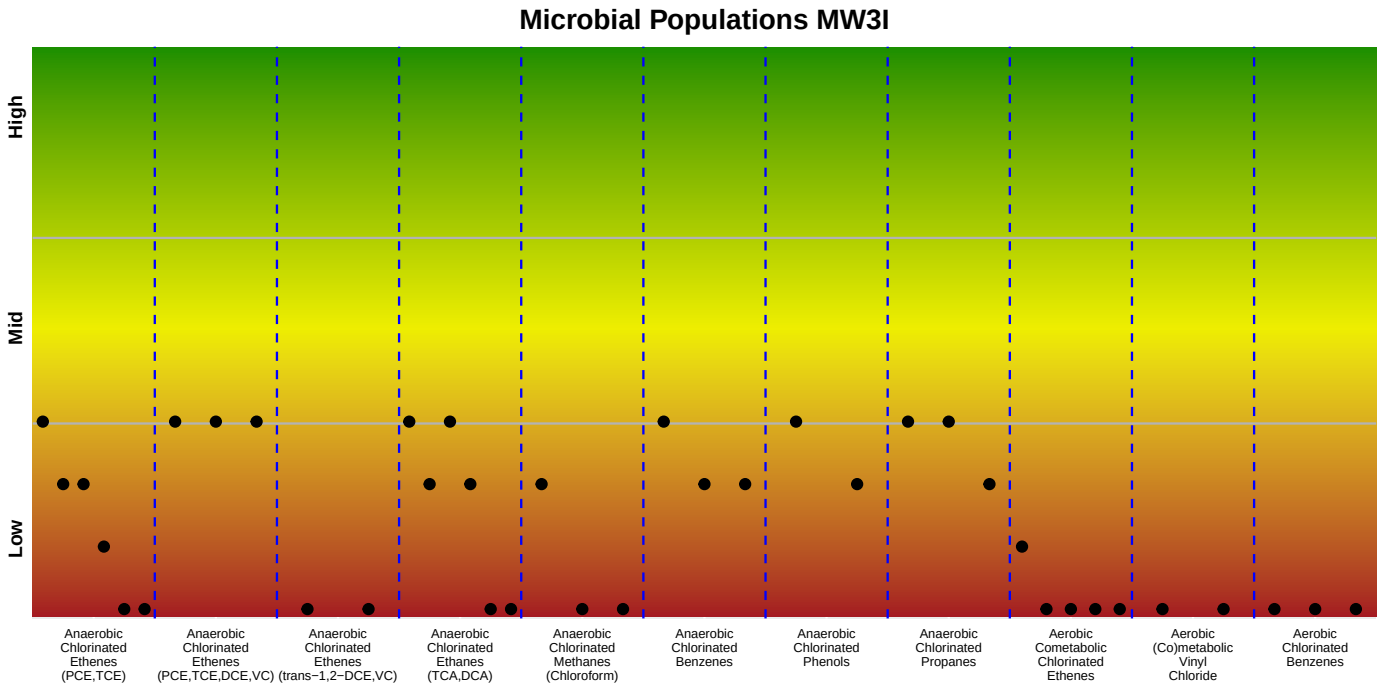

Figure 2: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants. See below to understand the correlation between contaminants and their associated qPCR assays.

Anaerobic - Reductive Dechlorination or Dichloroelimination	
Contaminants	**qPCR Assays**
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER
Chlorinated Ethanes (TCA and 1,2- DCA)	DHC, DHBt, DHG, DSB[1], DCA, DCAR
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR
Chlorinated Benzenes	DHC, DHBt[2], DECO
Chlorinated Phenols	DHC, DSB
Chlorinated Propanes	DHC, DHG, DSB[1]
Aerobic - (Co)metabolism	
Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Benzenes	TOD, TCBO, PHE

[1] *Desulfitobacterium dichloroeliminans* DCA1. [2] Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

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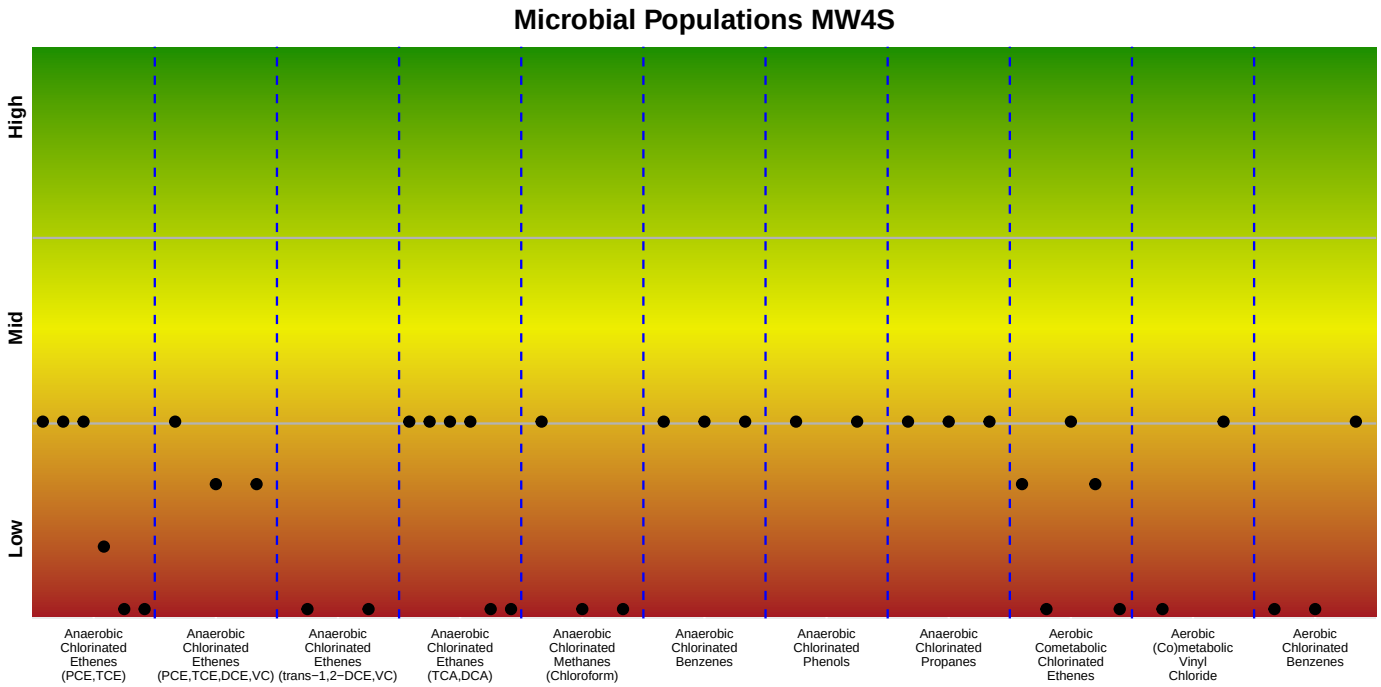

Figure 3: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants. See below to understand the correlation between contaminants and their associated qPCR assays.

Anaerobic - Reductive Dechlorination or Dichloroelimination	
Contaminants	**qPCR Assays**
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER
Chlorinated Ethanes (TCA and 1,2- DCA)	DHC, DHBt, DHG, DSB[1], DCA, DCAR
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR
Chlorinated Benzenes	DHC, DHBt[2], DECO
Chlorinated Phenols	DHC, DSB
Chlorinated Propanes	DHC, DHG, DSB[1]
Aerobic - (Co)metabolism	
Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Benzenes	TOD, TCBO, PHE

[1] *Desulfitobacterium dichloroeliminans* DCA1. [2] Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

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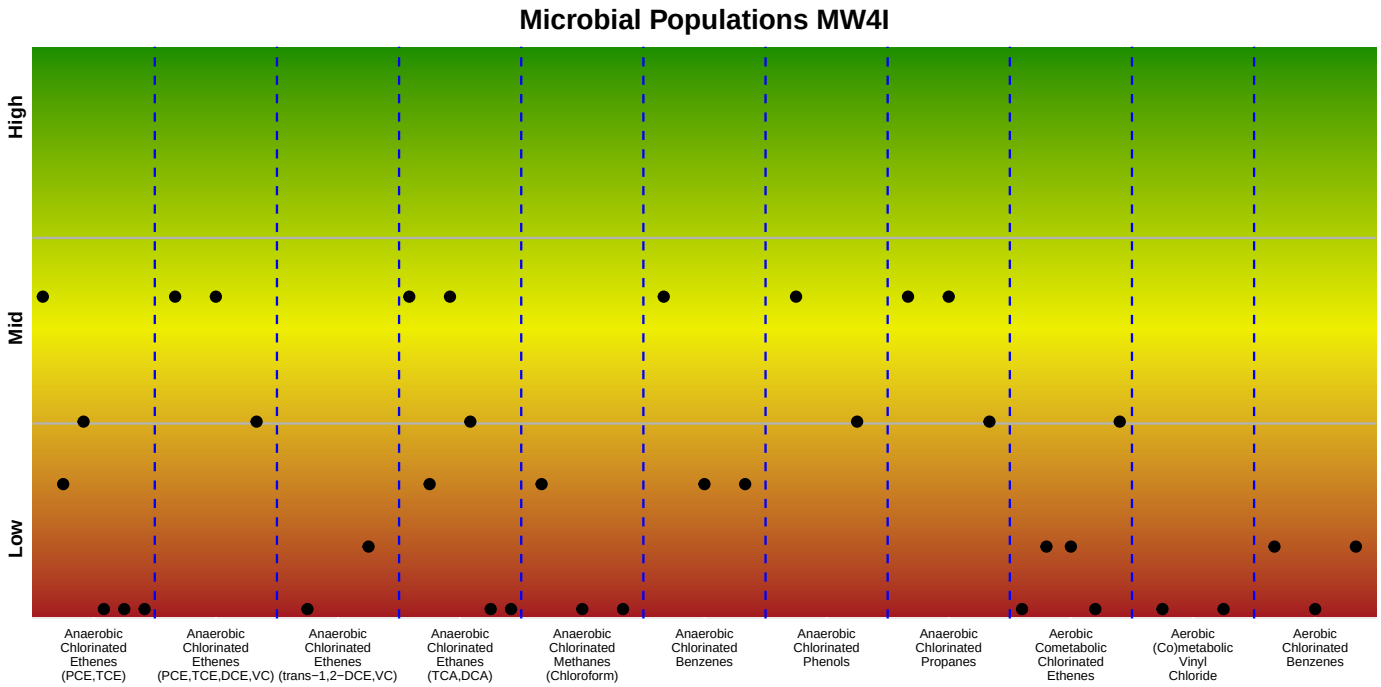

Figure 4: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants. See below to understand the correlation between contaminants and their associated qPCR assays.

Anaerobic - Reductive Dechlorination or Dichloroelimination	
Contaminants	**qPCR Assays**
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER
Chlorinated Ethanes (TCA and 1,2- DCA)	DHC, DHBt, DHG, DSB[1], DCA, DCAR
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR
Chlorinated Benzenes	DHC, DHBt[2], DECO
Chlorinated Phenols	DHC, DSB
Chlorinated Propanes	DHC, DHG, DSB[1]
Aerobic - (Co)metabolism	
Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Benzenes	TOD, TCBO, PHE

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Interpretation

The overall purpose of the QuantArray®-Chlor is to give site managers the ability to simultaneously yet economically evaluate the potential for biodegradation of a spectrum of common chlorinated contaminants through a multitude of anaerobic and aerobic (co)metabolic pathways in order to provide a clearer and more comprehensive view of contaminant biodegradation. The following discussion describes the interpretation of results in general terms and is meant to serve as a guide.

Reductive Dechlorination - Chlorinated Ethenes: While a number of bacterial cultures including *Dehalococcoides*, *Dehalobacter*, *Desulfitobacterium*, *Desulfuromonas* spp. capable, of utilizing PCE and TCE as growth-supporting electron acceptors have been isolated [1–5], *Dehalococcoides* may be the most important because they are the only bacterial group that has been isolated to date which is capable of complete reductive dechlorination of PCE to ethene [6]. In fact, the presence of *Dehalococcoides* has been associated with complete reductive dechlorination to ethene at sites across North America and Europe. Hendrickson and Lu have proposed using a *Dehalococcoides* concentration of 1 x 10^4 cells/mL as a screening criterion to identify sites where biological reductive dechlorination is predicted to proceed at "generally useful" rates [7,8].

At chlorinated ethene sites, any "stall" leading to the accumulation of daughter products, especially vinyl chloride, would be a substantial concern. While *Dehalococcoides* concentrations greater than 1 x 10^4 cells/mL correspond to ethene production and useful rates of dechlorination, the range of chlorinated ethenes degraded varies by strain within the *Dehalococcoides* genus [6,9] and the presence of co-contaminants and competitors can have complex impacts on the halorespiring microbial community [10–15]. Therefore, QuantArray®-Chlor also provides quantification of a suite of reductive dehalogenase genes (PCE, TCE, BVC, VCR, CER, and TDR) to more definitively confirm the potential for reductive dechlorination of all chlorinated ethene compounds including vinyl chloride.

Perhaps most importantly, QuantArray®-Chlor quantifies TCE reductase (TCE) and both known vinyl chloride reductase genes (BVC, VCR) from *Dehalococcoides* to conclusively evaluate the potential for complete reductive dechlorination of chlorinated ethenes to non-toxic ethene [16–18]. In addition, the analysis also includes quantification of reductive dehalogenase genes from *Dehalogenimonas* spp. capable of reductive dechlorination of chlorinated ethenes. More specifically, these are the trans-1,2-DCE dehalogenase gene (TDR) from strain WBC-2 [19] and the vinyl chloride reductase gene (CER) from GP, the only known organisms other than *Dehalococcoides* capable of vinyl chloride reduction [20]. Finally, PCE reductase genes responsible for sequential reductive dechlorination of PCE to *cis*-DCE by *Sulfurospirillum* and *Geobacter* spp. are also quantified. In mixed cultures, evidence increasingly suggests that partial dechlorinators like *Sulfurospirillum}* and *Geobacter* may be responsible for the majority of reductive dechlorination of PCE to TCE and *cis*-DCE while *Dehalococcoides* functions more as *cis*-DCE and vinyl chloride reducing specialists [10,21].

Reductive Dechlorination - Chlorinated Ethanes: Under anaerobic conditions, chlorinated ethanes are susceptible to reductive dechlorination by several groups of halorespiring bacteria including *Dehalobacter*, *Dehalogenimonas*, and *Dehalococcoides*. While the reported range of chlorinated ethanes utilized varies by genus, species, and sometimes at the strain level, several general observations can be made regarding biodegradation pathways and daughter product formation. *Dehalobacter* spp. have been isolated that are capable of sequential reductive dechlorination of 1,1,1-TCA through 1,1-DCA to chloroethane [13]. Biodegradation of 1,1,2-TCA by several halorespiring bacteria including *Dehalobacter* and *Dehalogenimonas* spp. proceeds via dichloroelimination producing vinyl chloride [22–24]. Similarly, 1,2-DCA biodegradation by *Dehalobacter*, *Dehalogenimonas*, and *Dehalococcoides* occurs via dichloroelimination producing ethene. While not utilized by many *Desulfitobacterium* isolates, at least one strain, *Desulfitobacterium dichloroeliminans* strain DCA1, is also capable of dichloroelimination of 1,2-DCA [25]. The 1,2-dichloroethane reductive dehalogenase gene (DCAR) from members of *Desulfitobacterium* and *Dehalobacter* is known to dechlorinate 1,2-DCA to ethene, while the 1,1-dichloroethane reductive dehalogenase (DCA) targets the gene responsible for 1,1-DCA dechlorination in some strains of *Dehalobacter*. In addition to chloroform, chloroform reductase (CFR) has also been shown to be responsible for reductive dechlorination of 1,1,1-TCA [26].

Reductive Dechlorination - Chlorinated Methanes: Chloroform is a common co-contaminant at chlorinated solvent sites and can inhibit reductive dechlorination of chlorinated ethenes. Grostern et al. demonstrated that a *Dehalobacter* population was capable of reductive dechlorination of chloroform to produce dichloromethane [27]. The *cfrA* gene encodes the reductase which catalyzes this initial step in chloroform biodegradation [26]. Justicia-Leon et al. have since shown that dichloromethane can support growth of a distinct group of *Dehalobacter* strains via fermentation [28]. The *Dehalobacter* DCM assay targets the 16S rRNA gene of these strains.

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Reductive Dechlorination - Chlorinated Benzenes: Chlorinated benzenes are an important class of industrial solvents and chemical intermediates in the production of drugs, dyes, herbicides, and insecticides. The physical-chemical properties of chlorinated benzenes as well as susceptibility to biodegradation are functions of their degree of chlorination and the positions of chlorine substituents. Under anaerobic conditions, reductive dechlorination of higher chlorinated benzenes including hexachlorobenzene (HCB), pentachlorobenzene (PeCB), tetrachlorobenzene (TeCB) isomers, and trichlorobenzene (TCB) isomers has been well documented [29], although biodegradation of individual compounds and isomers varies between isolates. For example, *Dehalococcoides* strain CBDB1 reductively dechlorinats HCB, PeCB, all three TeCB isomers, 1,2,3-TCB, and 1,2,4-TCB [9]. *Dehalobium chlorocoercia* DF-1 has been shown to be capable of reductive dechlorination of HCB, PeCB, and 1,2,3,5-TeCB [31]. The dichlorobenzene (DCB) isomers and chlorobenzene (CB) were considered relatively recalcitrant under anaerobic conditions. However, new evidence has demonstrated reductive dechlorination of DCBs to CB and CB to benzene [32] with corresponding increases in concentrations of *Dehalobacter* spp. [33].

Reductive Dechlorination - Chlorinated Phenols: Pentachlorophenol (PCP) was one of the most widely used biocides in the U.S. and despite residential use restrictions, is still extensively used industrially as a wood preservative. Along with PCP, the tetrachlorophenol and trichlorophenol isomers were also used as fungicides in wood preserving formulations. 2,4-Dichlorophenol and 2,4,5-TCP were used as chemical intermediates in herbicide production (e.g. 2,4-D) and chlorophenols are known byproducts of chlorine bleaching in the pulp and paper industry. While the range of compounds utilized varies by strain, some *Dehalococcoides* isolates are capable of reductive dechlorination of PCP and other chlorinated phenols. For example, *Dehalococcoides* strain CBDB1 is capable of utilizing PCP, all three tetrach lorophenol (TeCP) congeners, all six trichlorophenol (TCP) congeners, and 2,3-dichlorophenol (2,3-DCP). PCP dechlorination by strain CBDB1 produces a mixture of 3,5-DCP, 3,4-DCP, 2,4-DCP, 3-CP, and 4-CP [34]. In the same study, however, *Dehalococcoides* strain 195 dechlorinated a more narrow spectrum of chlorophenols which included 2,3-DCP, 2,3,4-TCP, and 2,3,6-TCP, but no other TCPs or PCP. Similar to *Dehalococcoides*, some species and strains of *Desulfitobacterium* are capable of utilizing PCP and other chlorinated phenols. *Desulfitobacterium hafniense* PCP-1 is capable of reductive dechlorination of PCP to 3-CP [35]. However, the ability to biodegrade PCP is not universal among *Desulfitobacterium* isolates. *Desulfitobacterium* spp. strain PCE1 and *D. chlororespirans* strain Co23, for example, can utilize some TCP and DCP isomers, but not PCP for growth [2,36].

Reductive Dechlorination - Chlorinated Propanes: *Dehalogenimonas* is a recently described bacterial genus of the phylum Chloroflexi which also includes the well-known chloroethene-respiring *Dehalococcoides* [23]. The *Dehalogenimonas* isolates characterized to date are also halorespiring bacteria, but utilize a rather unique range of chlorinated compounds as electron acceptors including chlorinated propanes (1,2,3-TCP and 1,2-DCP) and a variety of other vicinally chlorinated alkanes including 1,1,2,2-tetrachloroethane, 1,1,2-trichloroethane, and 1,2-dichloroethane [23].

Aerobic - Chlorinated Ethene Cometabolism: Under aerobic conditions, several different types of bacteria including methane-oxidizing bacteria (methanotrophs), and many benzene, toluene, ethylbenzene, xylene, and (BTEX)-utilizing bacteria can cometabolize or co-oxidize TCE, DCE, and vinyl chloride [37]. In general, cometabolism of chlorinated ethenes is mediated by monooxygenase enzymes with "relaxed" specificity that oxidize a primary (growth supporting) substrate (*e.g.* methane) and co-oxidize the chlorinated compound (*e.g.* TCE). QuantArray®-Chlor provides quantification of a suite of genes encoding oxygenase enzymes capable of co-oxidation of chlorinated ethenes including soluble methane monooxygenase (sMMO). Soluble methane monooxygenases co-oxidize a broad range of chlorinated compounds [38–41] including TCE, *cis*-DCE, and vinyl chloride. Furthermore, soluble methane monooxygenases are generally believed to support greater rates of aerobic cometabolism [40]. QuantArray®-Chlor also quantifies aromatic oxygenase genes encoding ring hydroxylating toluene monooxygenase genes (RMO, RDEG), toluene dioxygenase (TOD) and phenol hydroxylases (PHE) capable of TCE co-oxidation [42–46]. TCE or a degradation product has been shown to induce expression of toluene monooxygenases in some laboratory studies [43,47] raising the possibility of TCE cometabolism with an alternative (non-aromatic) growth substrate. Moreover, while a number of additional factors must be considered, recent research under ESTCP Project 201584 has shown positive correlations between concentrations of monooxygenase genes (soluble methane monooxygenase, ring hydroxylating monooxygenases, and phenol hydroxylase) and the rate of TCE degradation [48].

Aerobic - Chlorinated Ethane Cometabolism: While less widely studied than cometabolism of chlorinated ethenes, some chlorinated ethanes are also susceptible to co-oxidation. As mentioned previously, soluble methane monooxygenases (sMMO) exhibit very relaxed specificity. In laboratory studies, sMMO has been shown to co-oxidize a number of chlorinated ethanes including 1,1,1-TCA and 1,2-DCA [38,40].

Aerobic - Vinyl Chloride Cometabolism: Beginning in the early 1990s, numerous microcosm studies demonstrated

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aerobic oxidation of vinyl chloride under MNA conditions without the addition of exogenous primary substrates. Since then, strains of *Mycobacterium*, *Nocardioides*, *Pseudomonas*, *Ochrobactrum*, and *Ralstonia* species have been isolated which are capable of aerobic growth on both ethene and vinyl chloride (see Mattes et al. [49] for a review). The initial steps in the pathway are the monooxygenase (*etn*ABCD) catalyzed conversion of ethene and vinyl chloride to their respective epoxyalkanes (epoxyethane and chlorooxirane), followed by epoxyalkane:CoM transferase (*etn*E) mediated conjugation and breaking of the epoxide [50].

Aerobic - Chlorinated Benzenes: In general, chlorobenzenes with four or less chlorine groups are susceptible to aerobic biodegradation and can serve as growth-supporting substrates. Toluene dioxygenase (TOD) has a relatively relaxed substrate specificity and mediates the incorporation of both atoms of oxygen into the aromatic ring of benzene and substituted benzenes (toluene and chlorobenzene). Comparison of TOD levels in background and source zone samples from a CB-impacted site suggested that CBs promoted growth of TOD-containing bacteria [51]. In addition, aerobic biodegradation of some trichlorobenzene and even tetrachlorobenzene isomers is initiated by a group of related trichlorobenzene dioxygenase genes (TCBO). Finally, phenol hydroxylases catalyze the continued oxidation and in some cases, the initial oxidation of a variety of monoaromatic compounds. In an independent study, significant increases in numbers of bacteria containing PHE genes corresponded to increases in biodegradation of DCB isomers [51].

Aerobic - Chlorinated Methanes: Many aerobic methylotrophic bacteria, belonging to diverse genera (*Hyphomicrobium*, *Methylobacterium*, *Methylophilus*, *Pseudomonas*, *Paracoccus*, and *Alibacter*) have been isolated which are capable of utilizing dichloromethane (DCM) as a growth substrate. The DCM metabolic pathway in methylotrophic bacteria is initiated by a dichloromethane dehalogenase (DCMA) gene. DCMA is responsible for aerobic biodegradation of dichloromethane by methylotrophs by first producing formaldehyde which is then further oxidized [52].

As discussed in previous sections, soluble methane monooxygenase (sMMO) exhibits relaxed specificity and co-oxidizes a broad spectrum of chlorinated hydrocarbons. In addition to chlorinated ethenes, sMMO has been shown to co-oxidize chloroform in laboratory studies [38,41].

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Appendix D Trend Monitoring Graphs

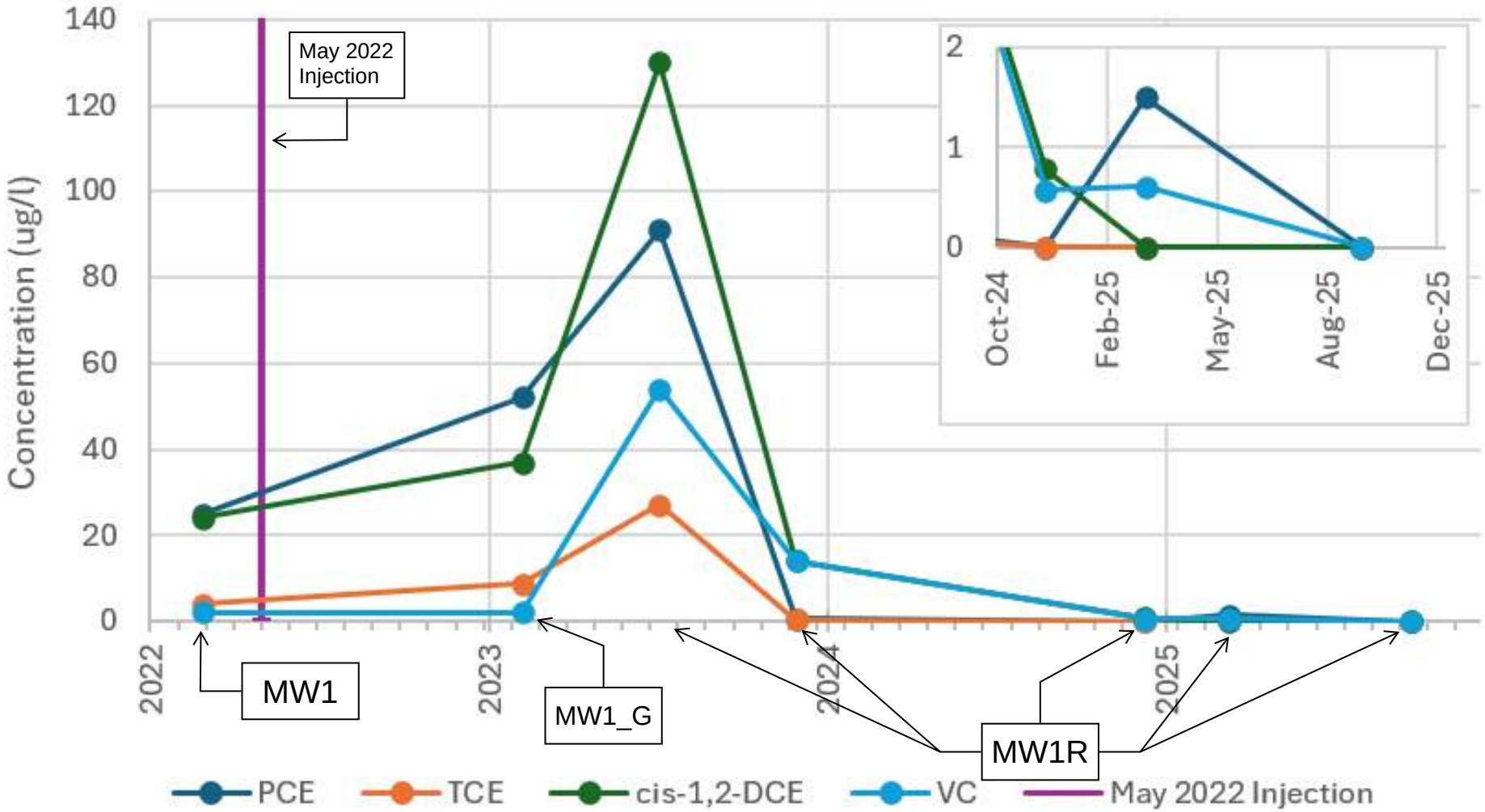
Appendix D-1. CVOC Trend Monitoring Graphs

Appendix D-2. PFAS Trend Monitoring Graphs

Appendix D-3. MNA Parameters Trend Monitoring Graphs

Appendix D-1. CVOC Trend Monitoring Graphs

MW1/1R CVOC Concentrations

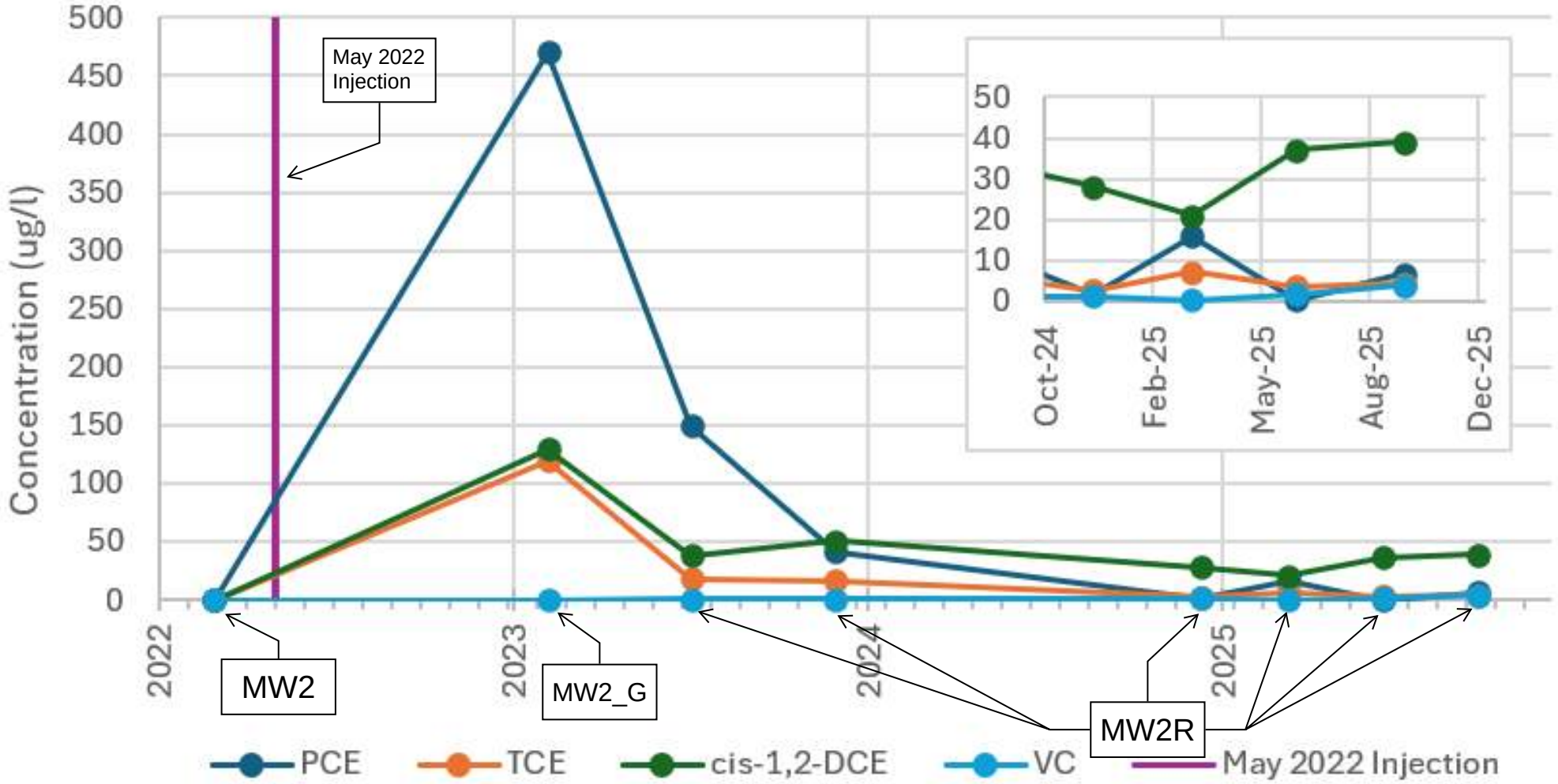


A Hydropunch sampler was used to collect groundwater from a temporary boring at the approximate location of MW1R in February 2023 because a permanent well could not be installed at that time due to ongoing facility construction. This sample is designated with a "_G" suffix.

Permanent well MW1R was installed at the approximate location of former permanent well MW1 and, as such, data for these wells are presented together.

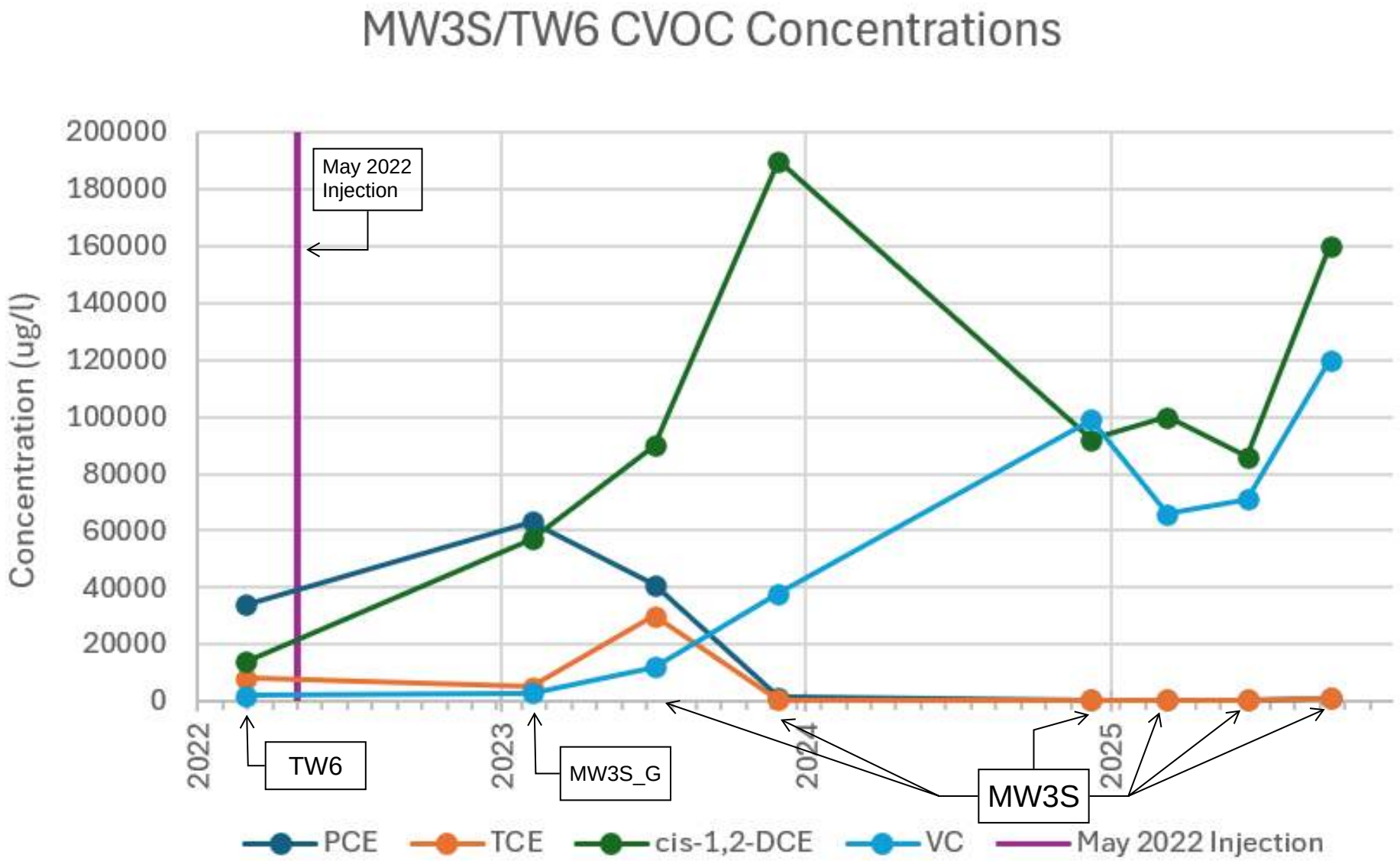
There was not a sample collected from MW1R in June 2025 due to the presence of NAPL.

MW2/2R CVOC Concentrations



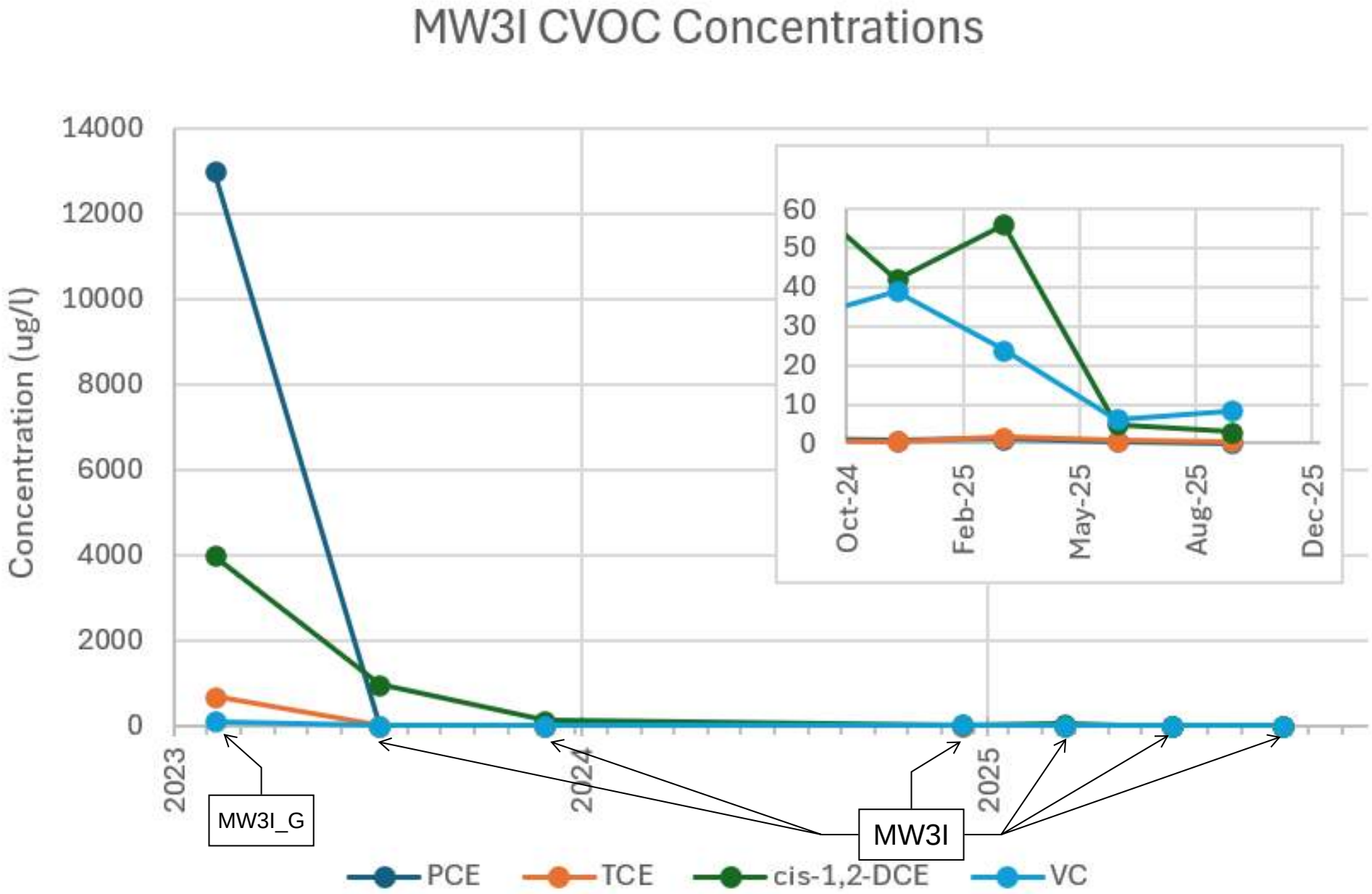
A Hydropunch sampler was used to collect groundwater from temporary boring at the approximate location of MW2R in February 2023 because a permanent well could not be installed at that time due to ongoing facility construction. This sample is designated with a "_G" suffix.

Permanent well MW2R was installed at the approximate location of former permanent well MW2 and, as such, data for these wells are presented together.



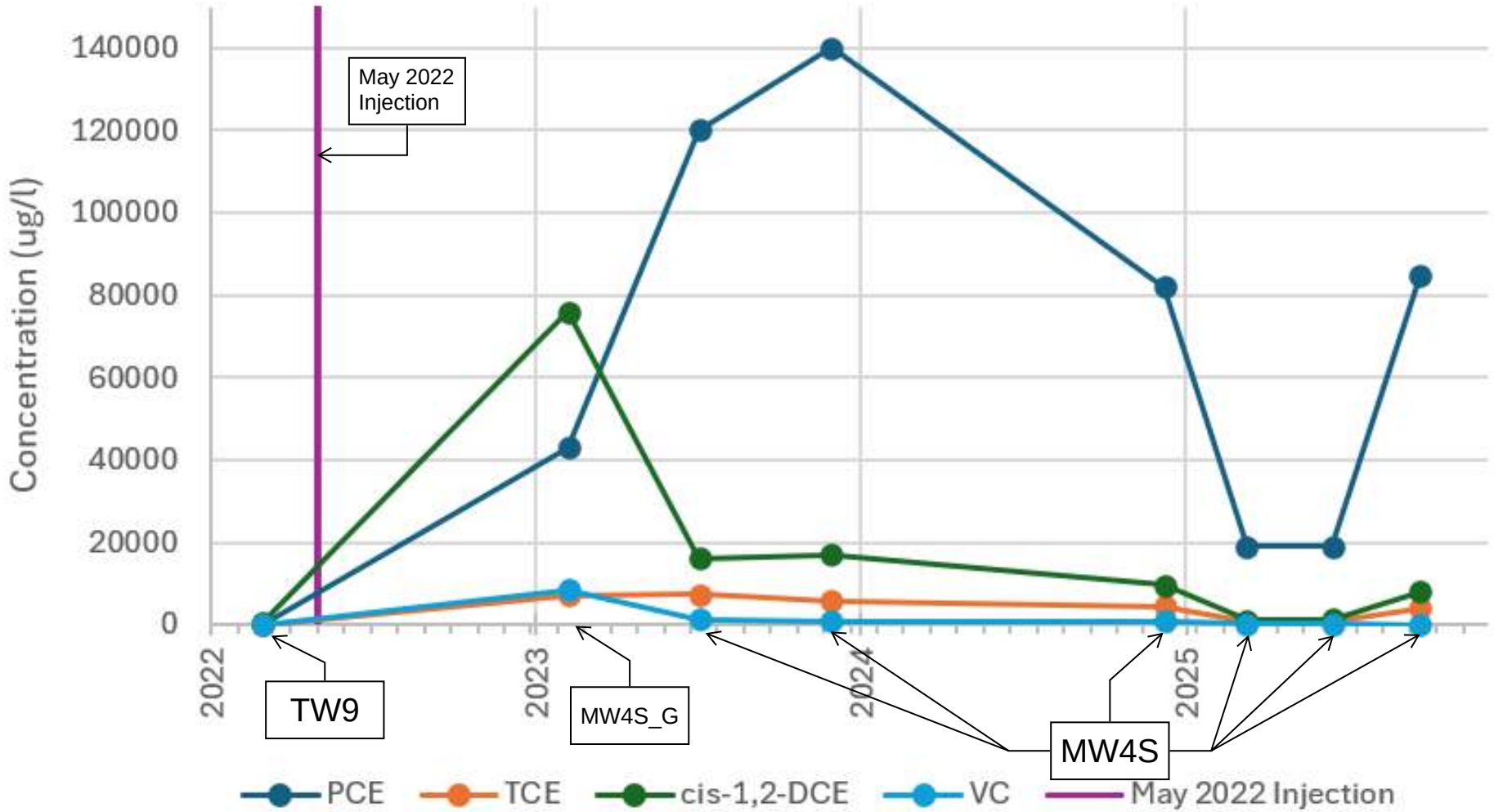
A Hydropunch sampler was used to collect groundwater from a temporary boring at the approximate location of MW3S/I in February 2023 because permanent wells could not be installed at that time due to ongoing facility construction. This sample is designated with a "_G" suffix.

Permanent well MW3S was installed at the approximate location of former temporary well TW6 and, as such, data for these co-located wells are presented together.



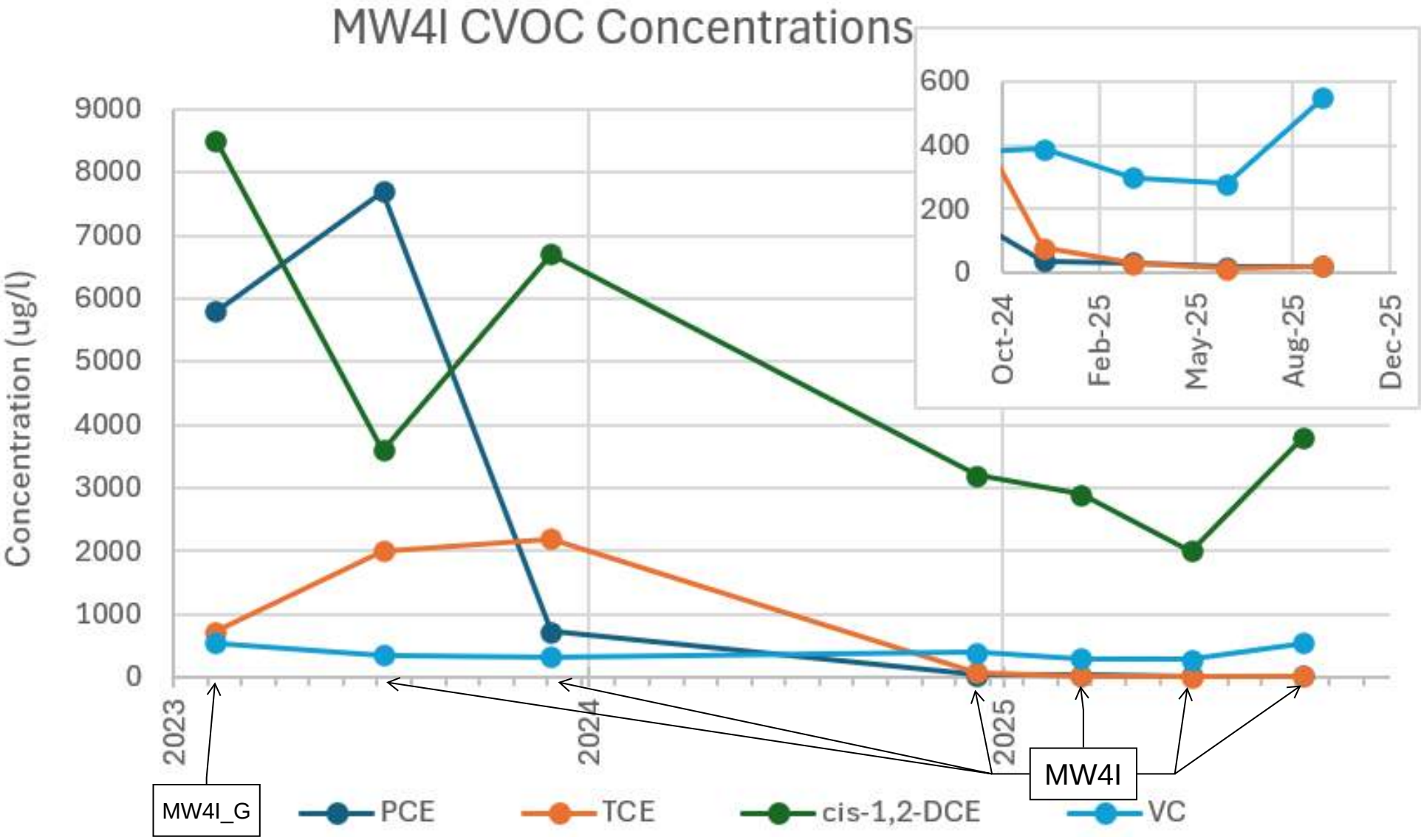
A Hydropunch sampler was used to collect groundwater from a temporary boring at the approximate location of MW3S/I in February 2023 because permanent wells could not be installed at that time due to ongoing facility construction. This sample is designated with a "_G" suffix.

MW4S/TW9 CVOC Concentrations



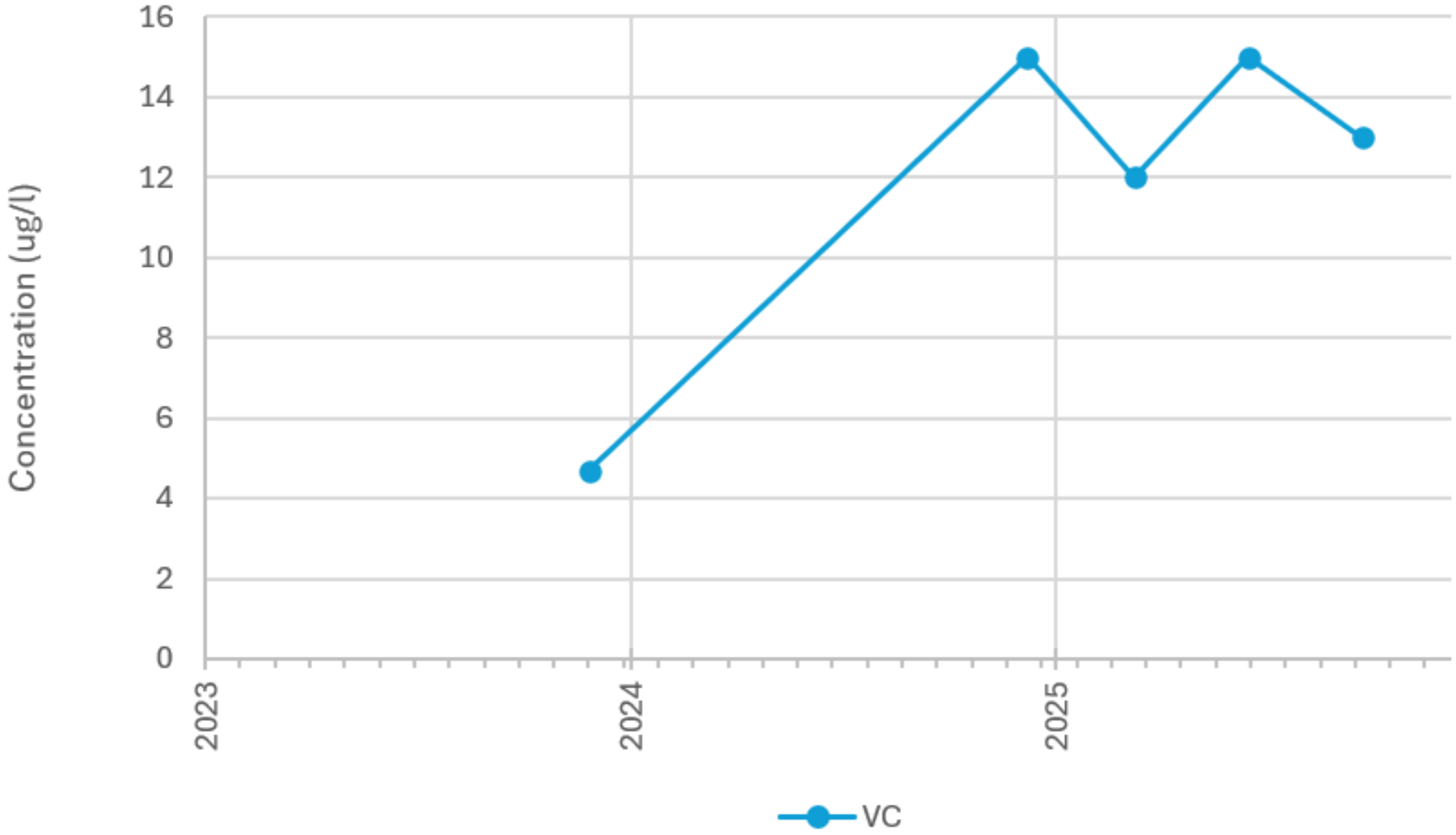
A Hydropunch sampler was used to collect groundwater from a temporary boring at the approximate location of MW4S/I in February 2023 because permanent wells could not be installed at that time due to ongoing facility construction. This sample is designated with a "_G" suffix.

Permanent well MW4S was installed at the approximate location of former temporary well TW9 and, as such, data for these co-located wells are presented together.

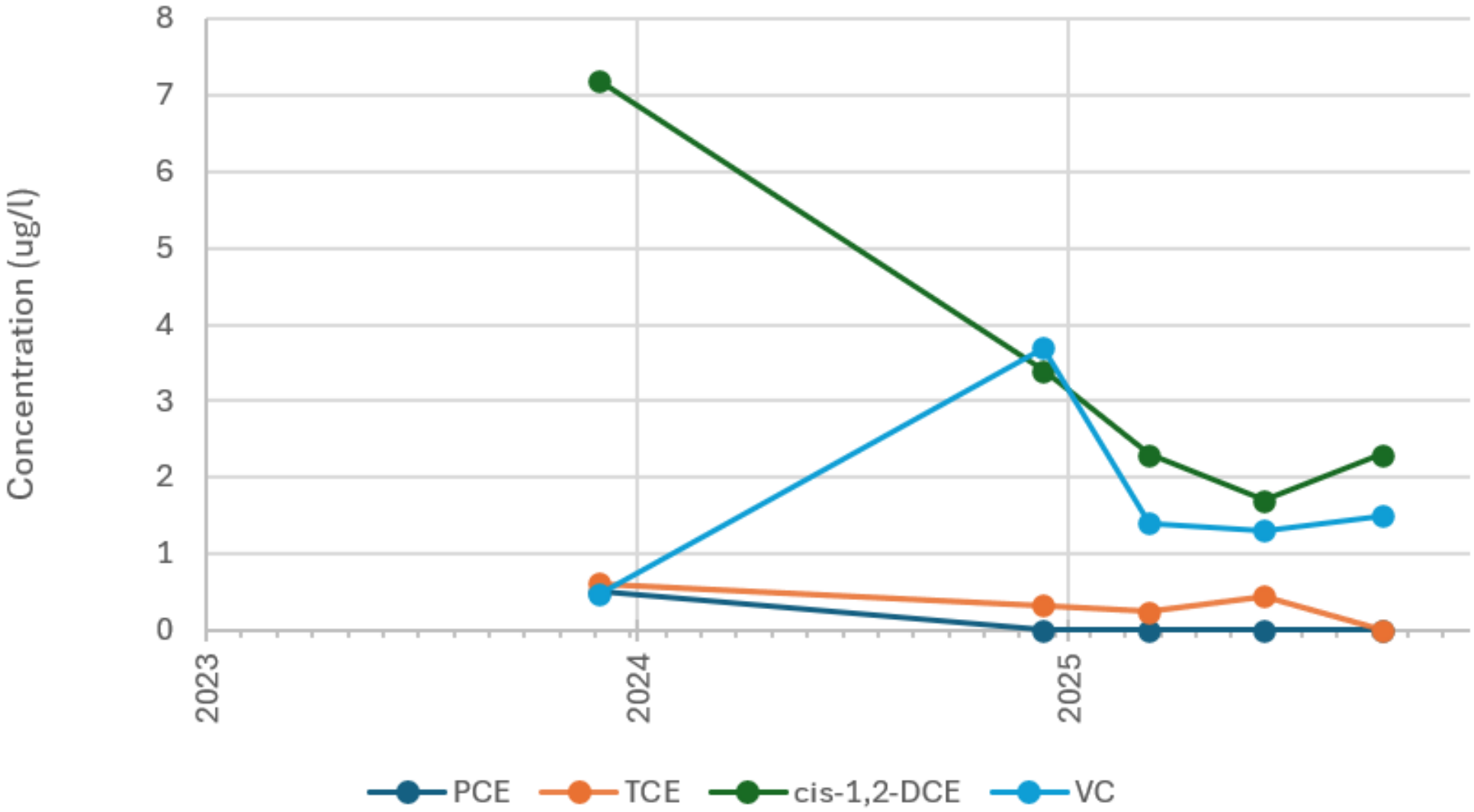


A Hydropunch sampler was used to collect groundwater from a temporary boring at the approximate locations of MW4S/I in February 2023 because permanent wells could not be installed at that time due to ongoing facility construction. This sample is designated with a "_G" suffix.

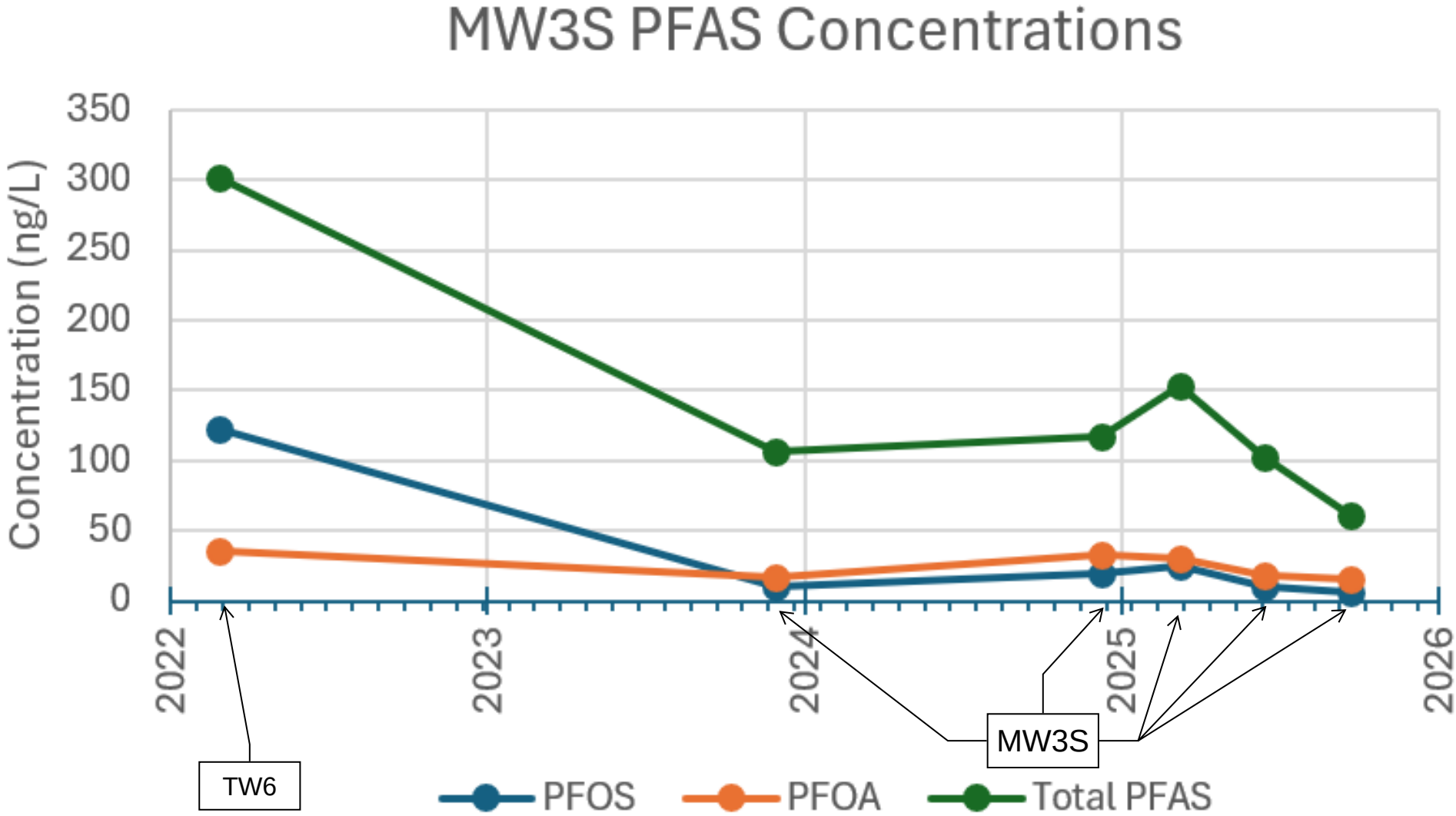
MW7 CVOC Concentrations



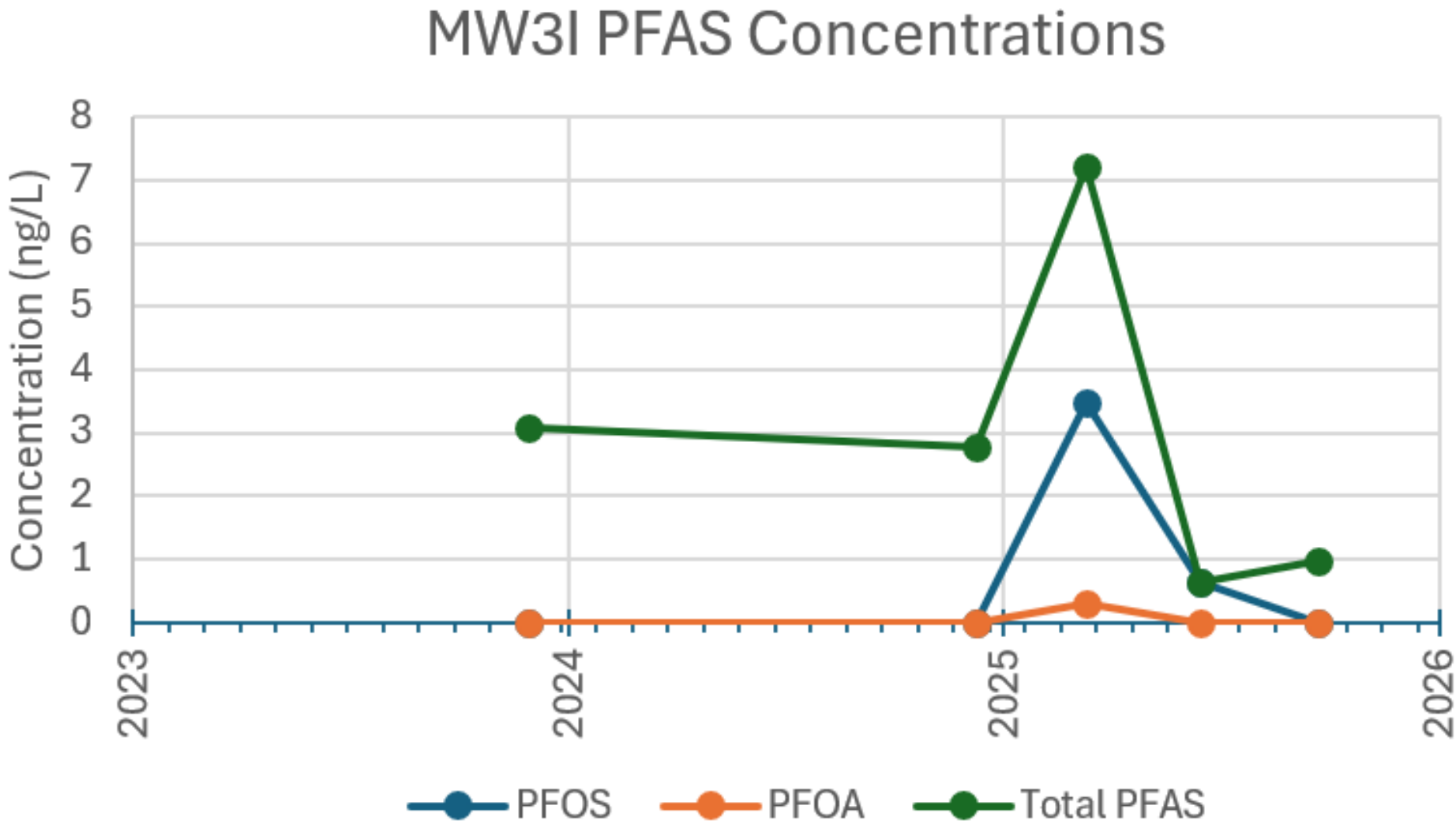
MW12S CVOC Concentrations

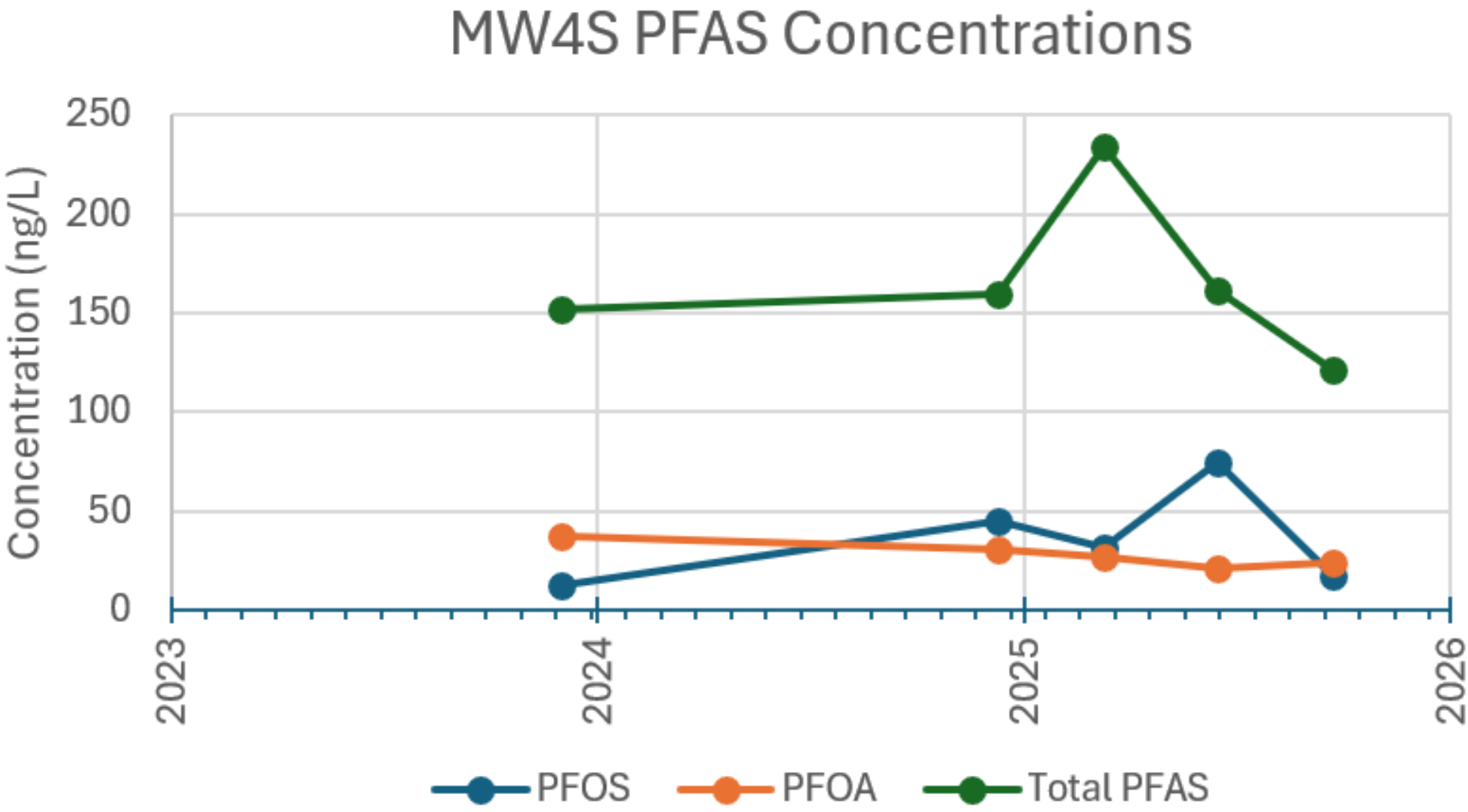


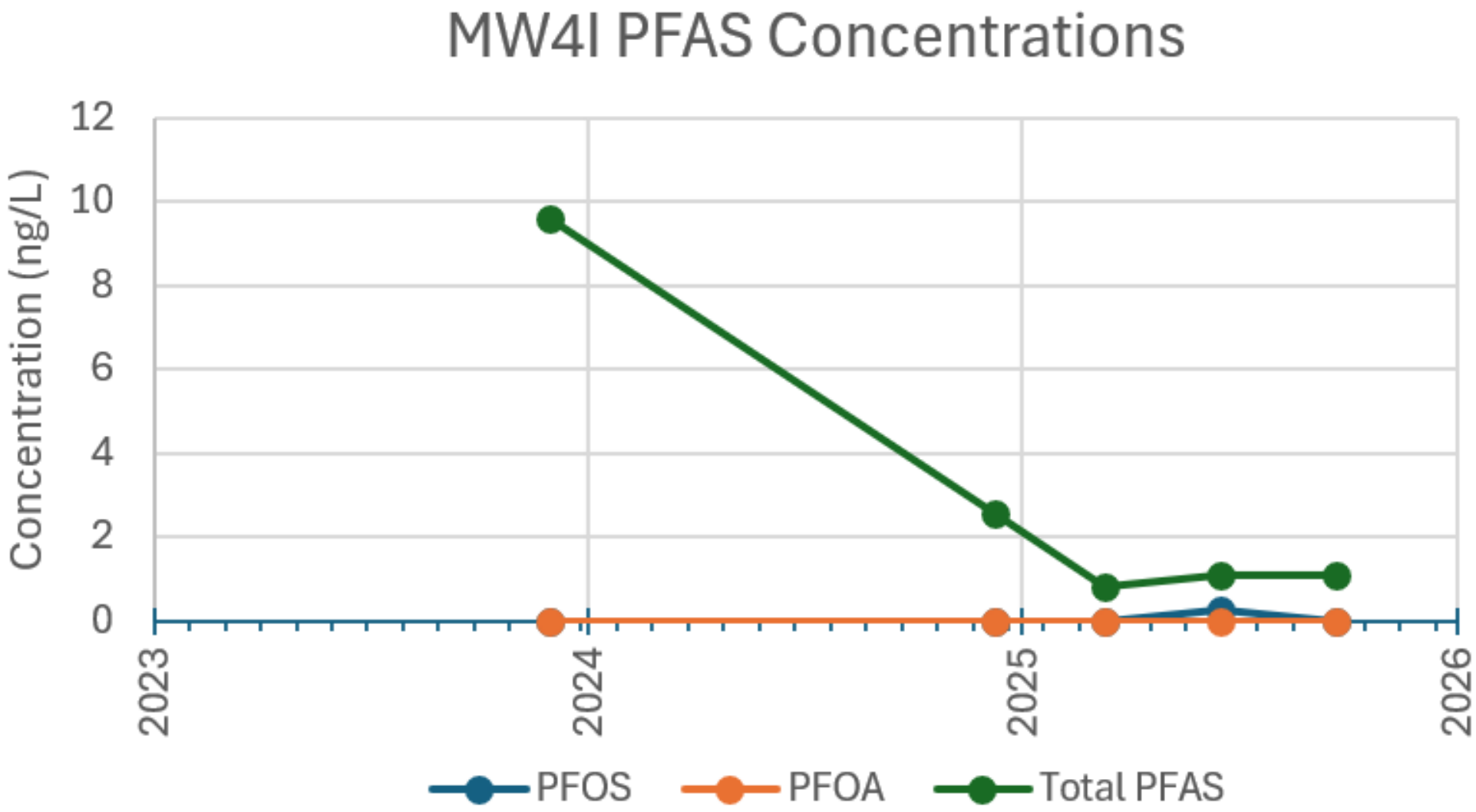
Appendix D-2. PFAS Trend Monitoring Graphs

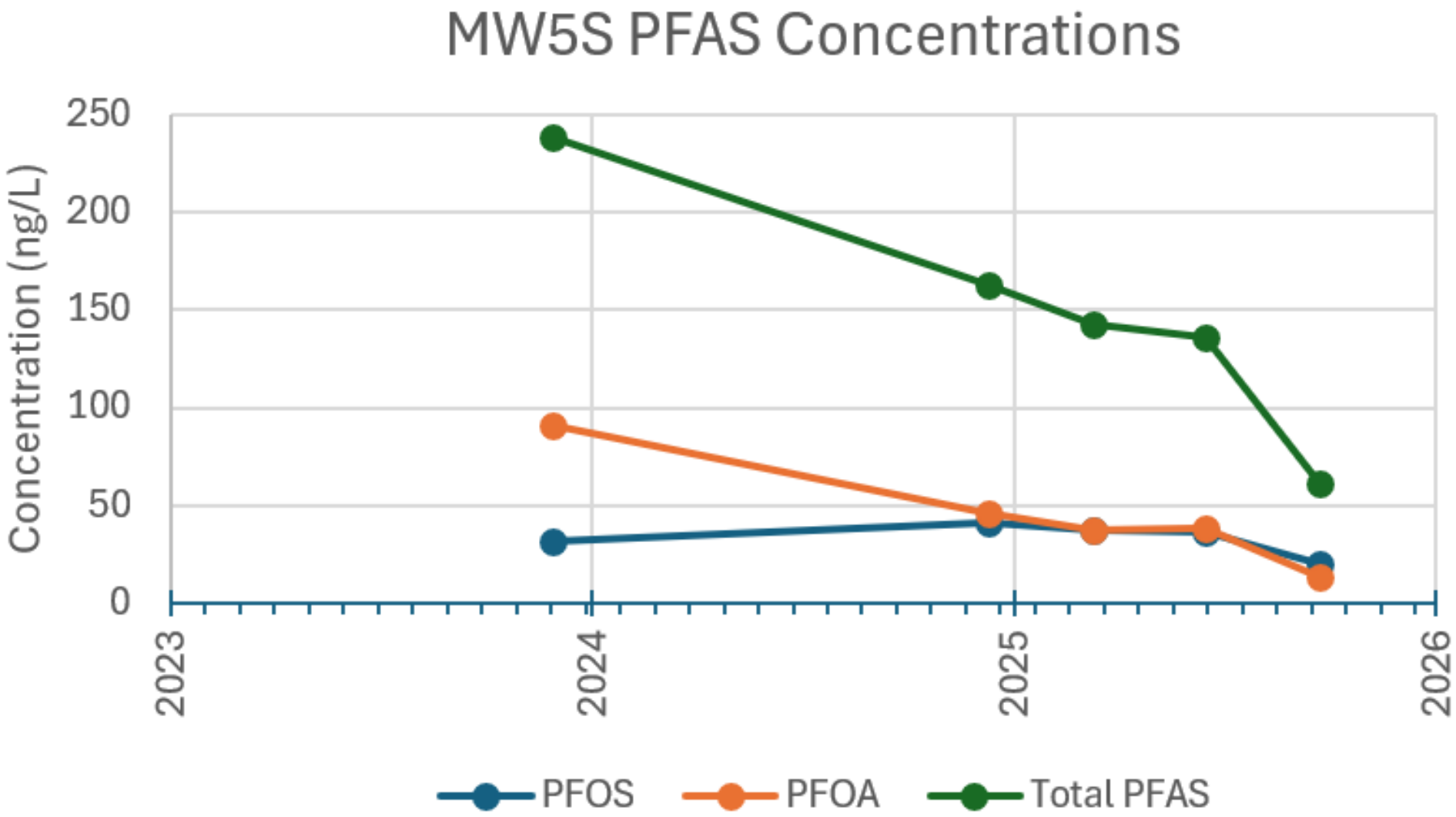


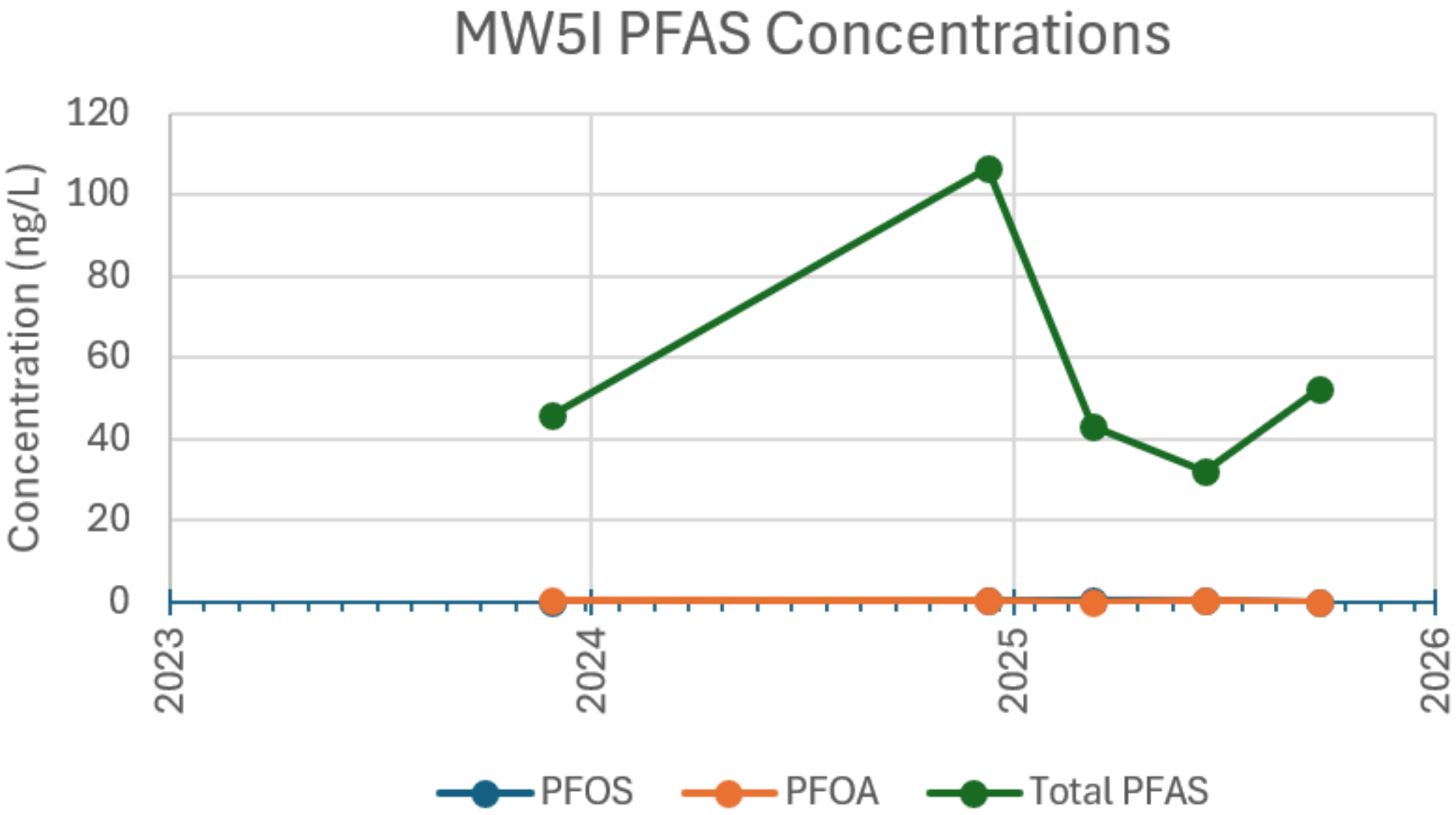
Permanent well MW3S was installed at the approximate location of former temporary well TW6 and, as such, data for these co-located wells are presented together.


MW3I PFAS Concentrations
Concentration (ng/L)
8
7
6
5
4
3
2
1
0
2023
2024
2025
2026
PFOS
PFOA
Total PFAS

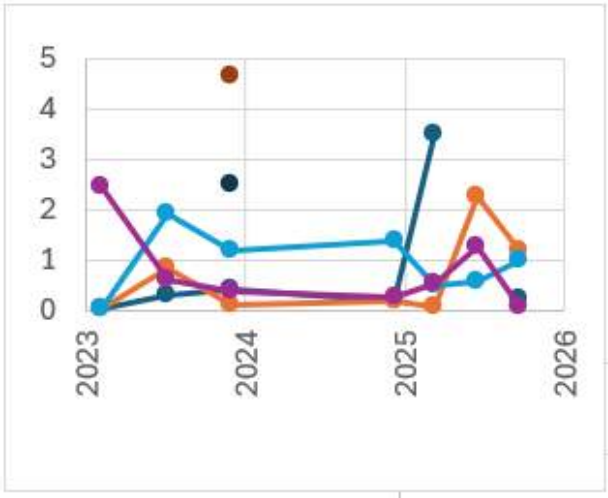

MW4S PFAS Concentrations
Concentration (ng/L)
250
200
150
100
50
0
2023
2024
2025
2026
PFOS
PFOA
Total PFAS


MW4I PFAS Concentrations
Concentration (ng/L)
12
10
8
6
4
2
0
2023
2024
2025
2026
PFOS
PFOA
Total PFAS

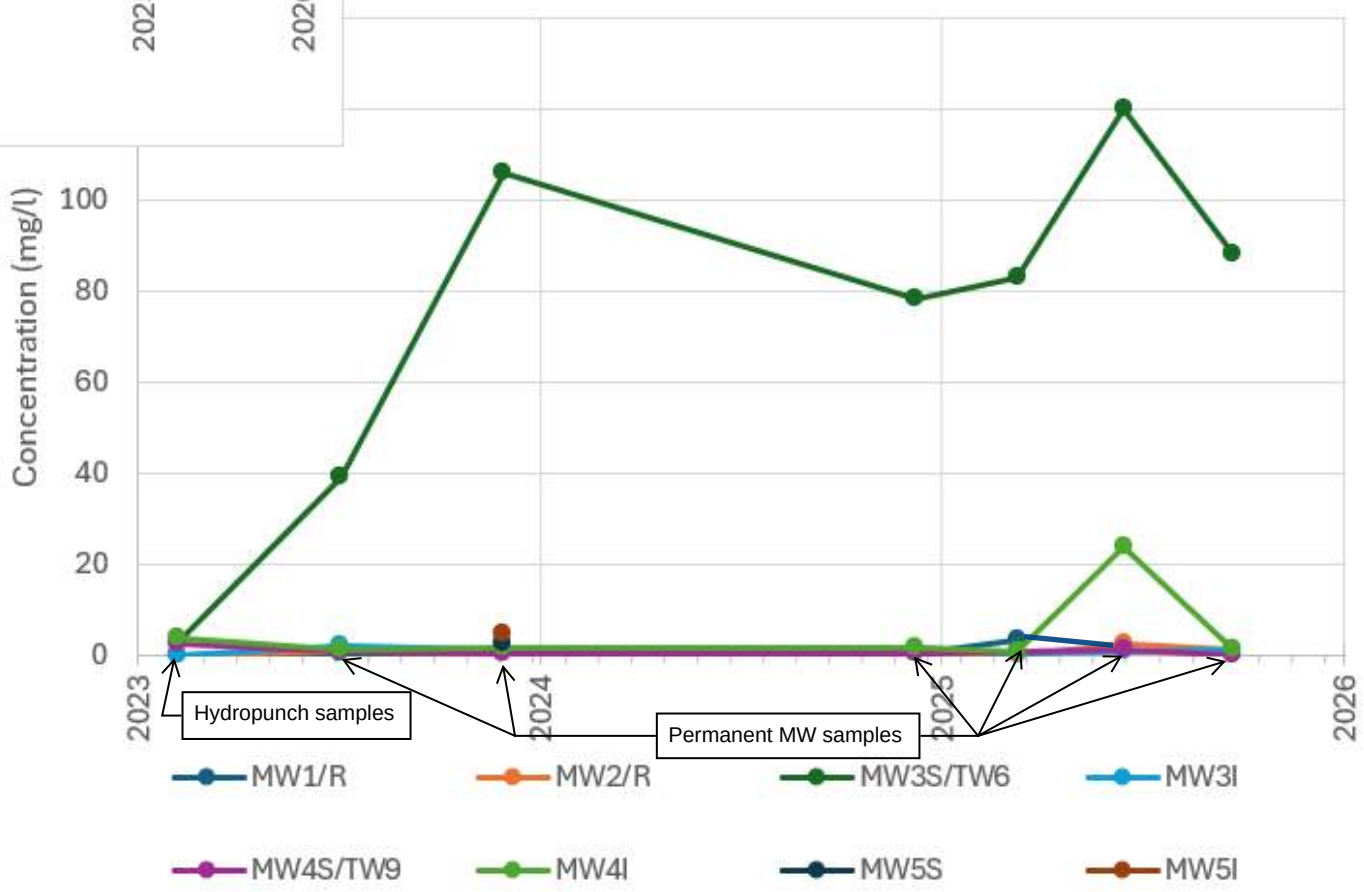

MW5S PFAS Concentrations
Concentration (ng/L)
250
200
150
100
50
0
2023
2024
2025
2026
PFOS
PFOA
Total PFAS


MW5I PFAS Concentrations
Concentration (ng/L)
120
100
80
60
40
20
0
2023
2024
2025
2026
PFOS
PFOA
Total PFAS

Appendix D-3. MNA Parameters Trend Monitoring Graphs



Dissolved Iron

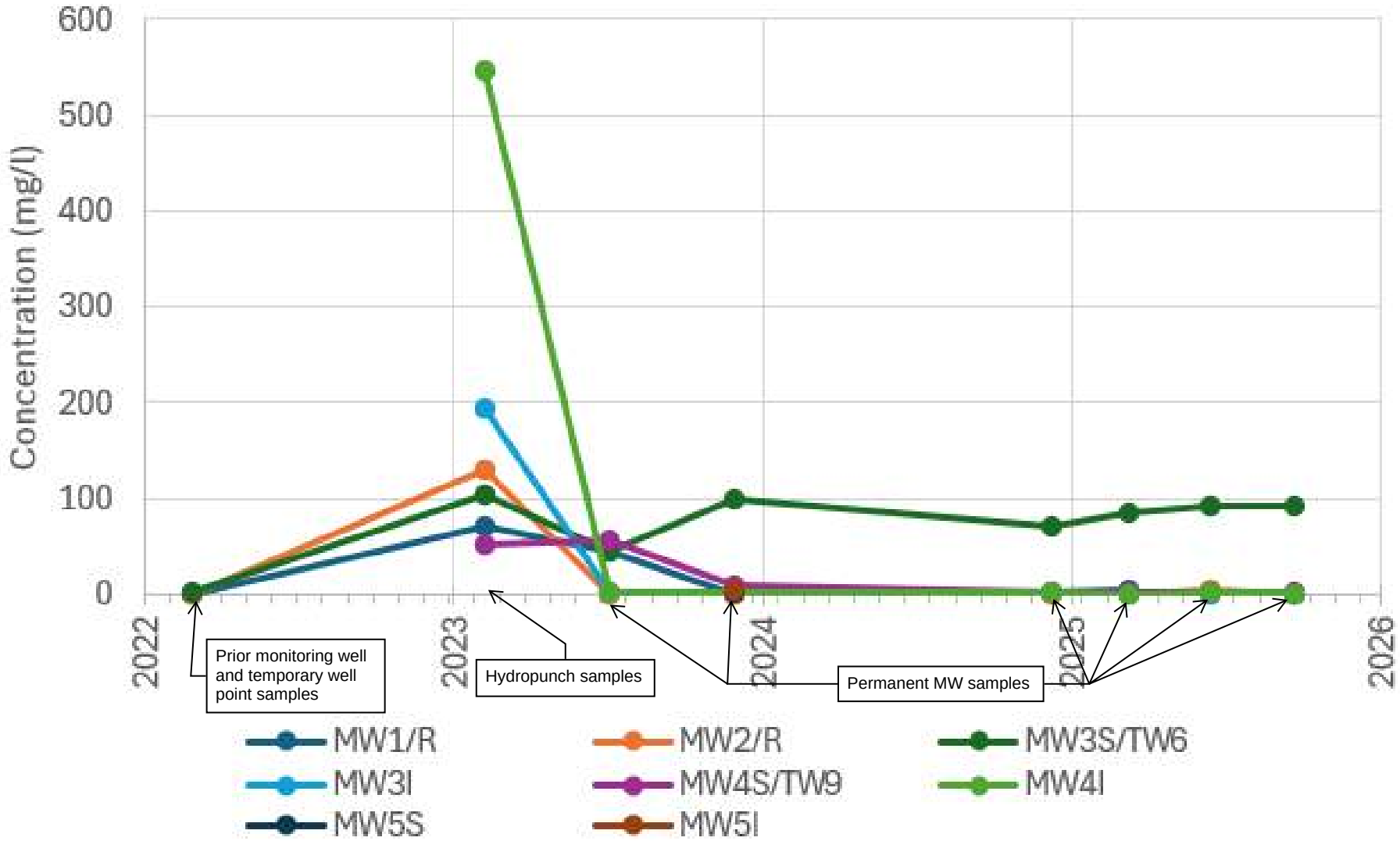


A Hydropunch sampler was used to collect groundwater from temporary borings at the approximate locations of MW1R MW2R, MW3S/I, and MW4S/I in February 2023 because permanent wells could not be installed at that time due to ongoing facility construction. These samples are designated with "_G" suffixes. One exception: samples from MW5S and MW5I in February 2023 were collected from permanent monitoring wells.

Permanent wells MW1R, MW2R, MW3S/I, and MW4S/I were installed at the approximate locations of former associated samples and, as such, data for these wells are presented together.

There was not a sample collected from MW1R in June 2025 due to the presence of NAPL.

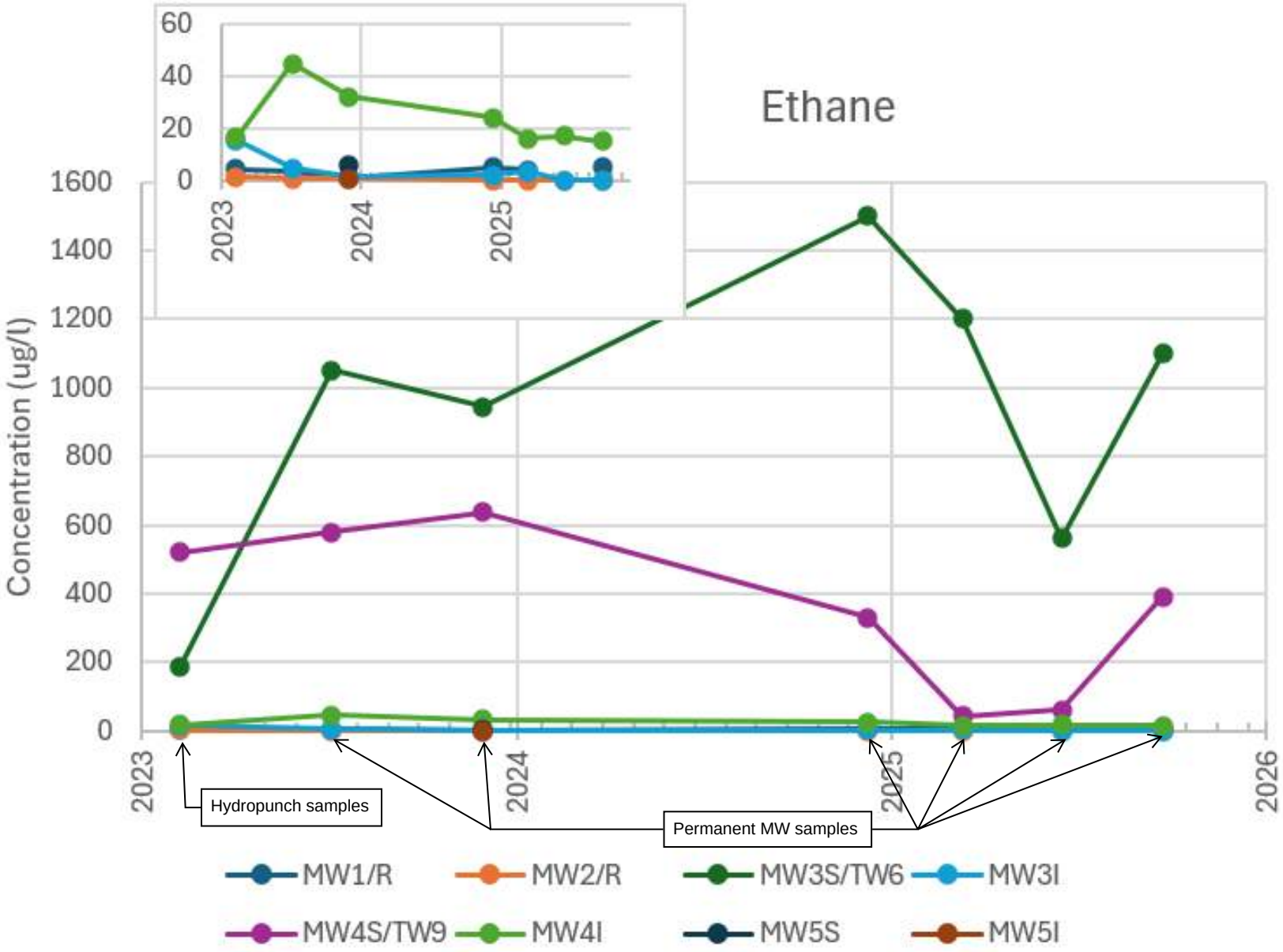
Total Iron



Samples collected in 2022 were taken from the prior monitoring wells MW1 and MW2 and temporary well points TW6 and TW9. A Hydropunch sampler was used to collect groundwater from temporary borings at the approximate locations of MW1R MW2R, MW3S/I, and MW4S/I in February 2023 because permanent wells could not be installed at that time due to ongoing facility construction. These samples are designated with "_G" suffixes. One exception: samples from MW5S and MW5I in February 2023 were collected from permanent monitoring wells.

Permanent wells MW1R, MW2R, MW3S/I, and MW4S/I were installed at the approximate locations of former associated samples and, as such, data for these wells are presented together.

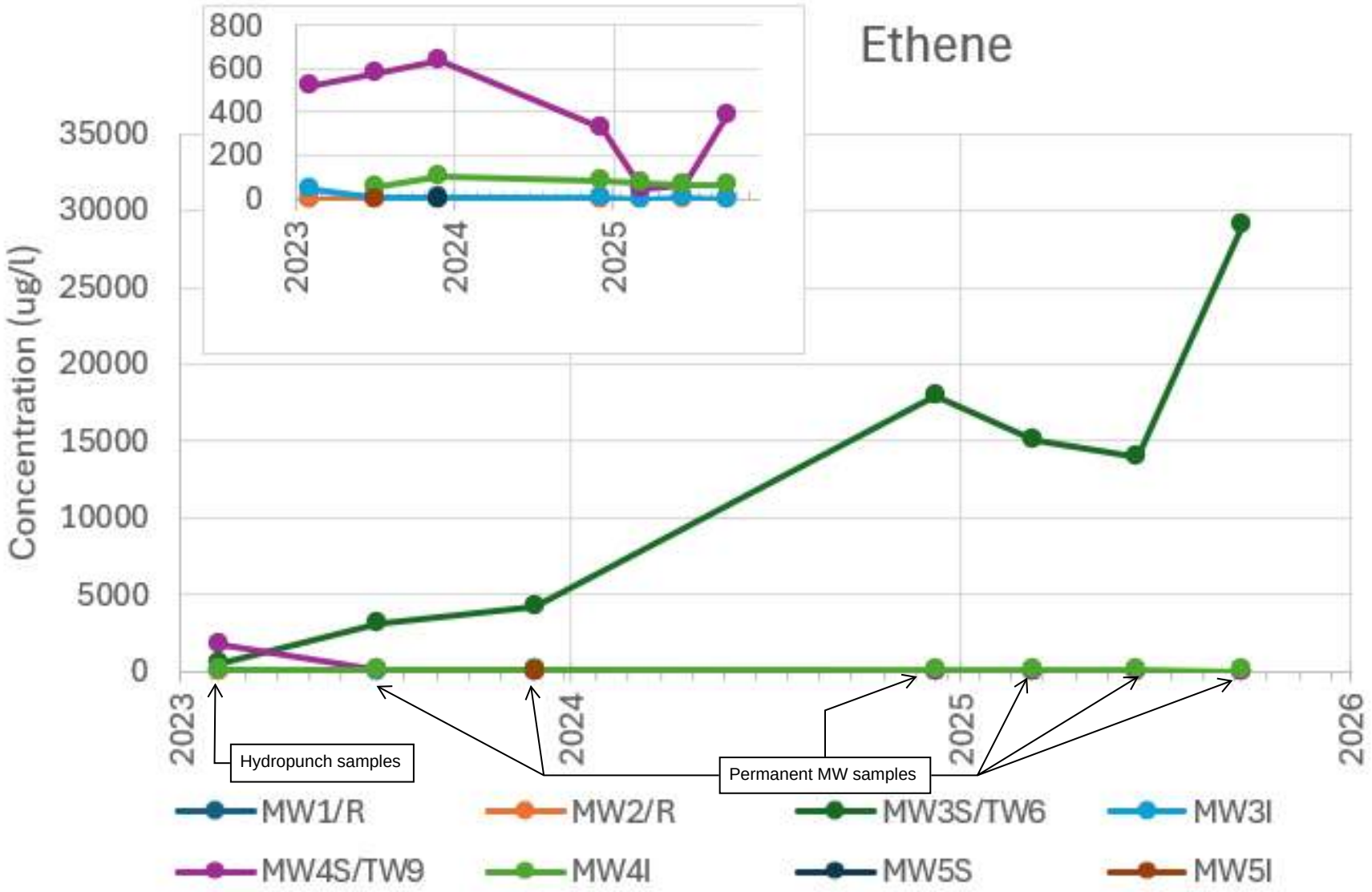
There was not a sample collected from MW1R in June 2025 due to the presence of NAPL.



A Hydropunch sampler was used to collect groundwater from temporary borings at the approximate locations of MW1R MW2R, MW3S/I, and MW4S/I in February 2023 because permanent wells could not be installed at that time due to ongoing facility construction. These samples are designated with "_G" suffixes. One exception: samples from MW5S and MW5I in February 2023 were collected from permanent monitoring wells.

Permanent wells MW1R, MW2R, MW3S/I, and MW4S/I were installed at the approximate locations of former associated samples and, as such, data for these wells are presented together.

There was not a sample collected from MW1R in June 2025 due to the presence of NAPL.

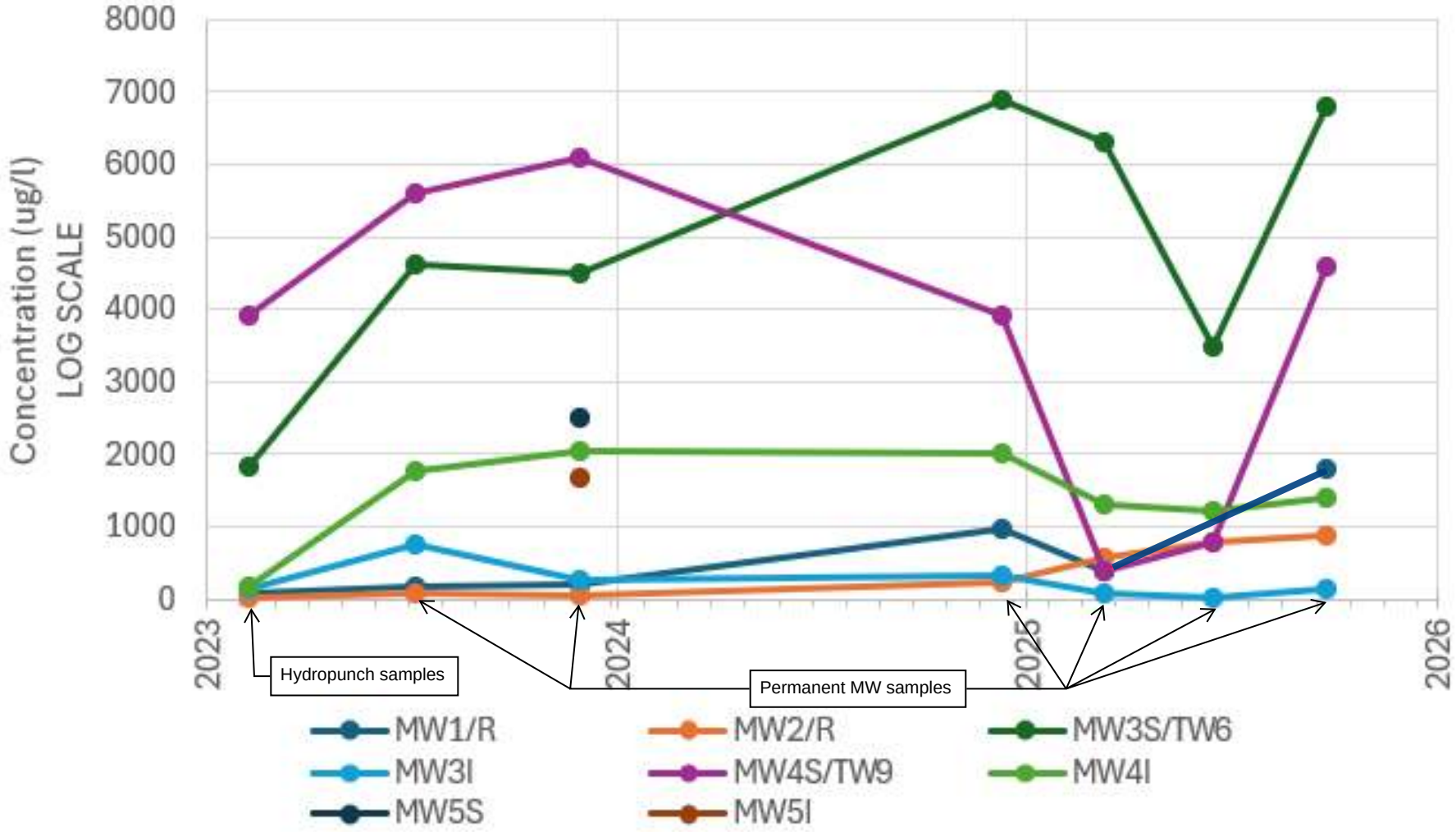


A Hydropunch sampler was used to collect groundwater from temporary borings at the approximate locations of MW1R MW2R, MW3S/I, and MW4S/I in February 2023 because permanent wells could not be installed at that time due to ongoing facility construction. These samples are designated with "_G" suffixes. One exception: samples from MW5S and MW5I in February 2023 were collected from permanent monitoring wells.

Permanent wells MW1R, MW2R, MW3S/I, and MW4S/I were installed at the approximate locations of former associated samples and, as such, data for these wells are presented together.

There was not a sample collected from MW1R in June 2025 due to the presence of NAPL.

Methane

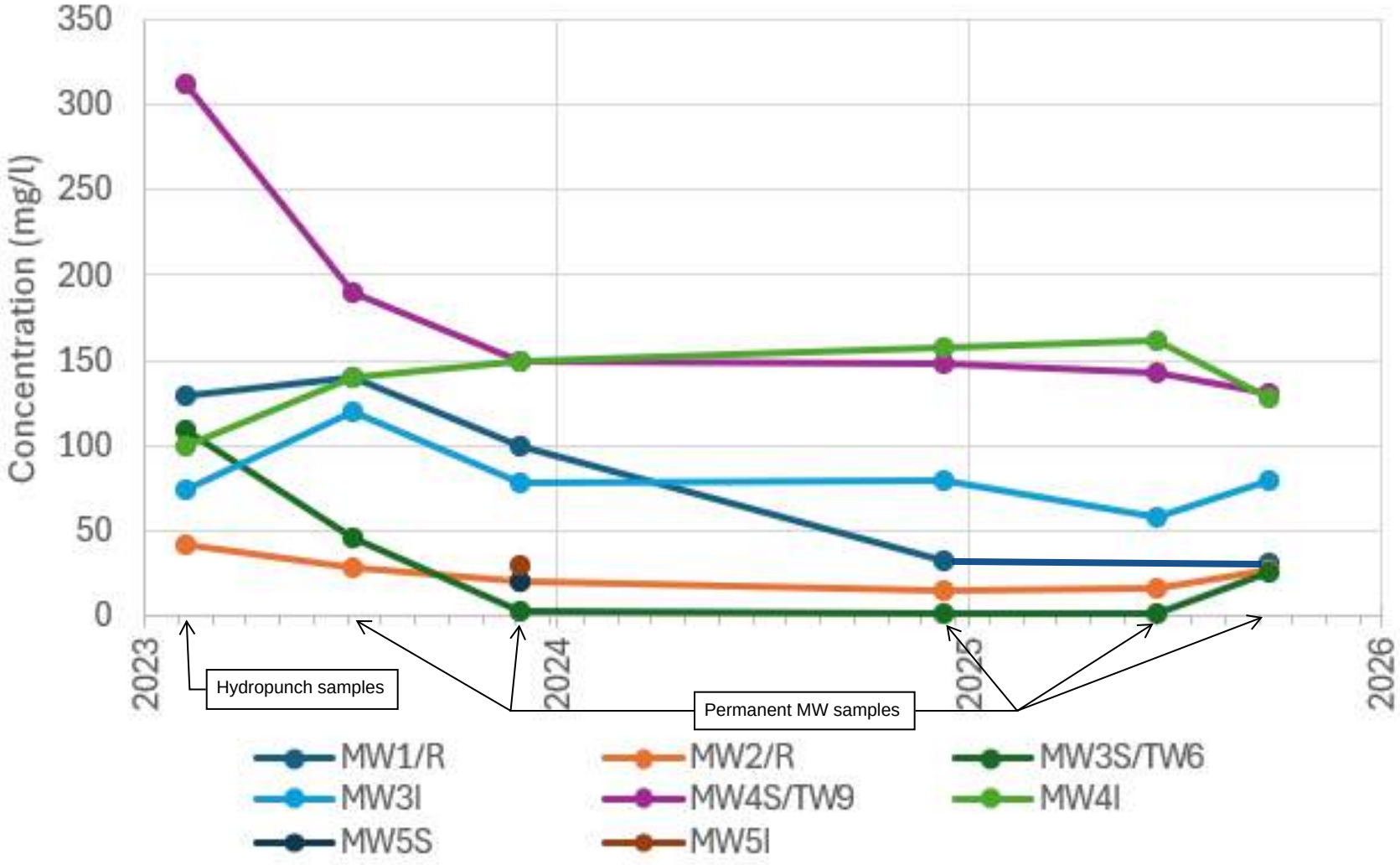


A Hydropunch sampler was used to collect groundwater from temporary borings at the approximate locations of MW1R MW2R, MW3S/I, and MW4S/I in February 2023 because permanent wells could not be installed at that time due to ongoing facility construction. These samples are designated with "_G" suffixes. One exception: samples from MW5S and MW5I in February 2023 were collected from permanent monitoring wells.

Permanent wells MW1R, MW2R, MW3S/I, and MW4S/I were installed at the approximate locations of former associated samples and, as such, data for these wells are presented together.

There was not a sample collected from MW1R in June 2025 due to the presence of NAPL.

Sulfate



A Hydropunch sampler was used to collect groundwater from temporary borings at the approximate locations of MW1R MW2R, MW3S/I, and MW4S/I in February 2023 because permanent wells could not be installed at that time due to ongoing facility construction. These samples are designated with "_G" suffixes. One exception: samples from MW5S and MW5I in February 2023 were collected from permanent monitoring wells.

Permanent wells MW1R, MW2R, MW3S/I, and MW4S/I were installed at the approximate locations of former associated samples and, as such, data for these wells are presented together.

There was not a sample collected from MW1R in June 2025 due to the presence of NAPL.