



Excellence Delivered As Promised

December 2, 2021

File #47358.003

Mr. Kevin McKnight
Remediation & Redevelopment Program
Wisconsin Department of Natural Resources
625 East County Road Y, Suite 700
Oshkosh, WI 54901-9731

Re: **Groundwater Monitoring Status Report**
Koeller Center, 1020 – 1142 South Koeller Street
Oshkosh, Wisconsin
BRRTS: 02-71-547941

Dear Mr. McKnight:

On behalf of Koeller One, LLC, Gannett Fleming, Inc. (GF) is submitting this groundwater monitoring status report for the former dry cleaner facility located at 1142 South Koeller Street in the Koeller Center shopping mall site in Oshkosh. Figure 1 is a site location map, and Figure 2 is an aerial photo that shows the area of investigation.

On April 17, 2017, GF submitted a *Closure Request* that summarized site investigation and groundwater monitoring results through November 2016. On June 20, 2017, the WDNR sent a letter to the Livesey Company, the owner of the Koeller Center, denying site closure. In its June 2017 letter, the WDNR indicated that additional groundwater monitoring of MW-1 through MW-4 and MW-6 was necessary to establish whether the tetrachloroethylene (PCE) plume was stable or receding. The WDNR's June 2017 letter also indicated that Koeller should evaluate whether additional investigation or remedial actions are necessary to obtain closure.

PCE concentrations measured in source area wells MW-1 through MW-3 decreased significantly over the last two sampling rounds in July 2020 and August 2021, while PCE concentrations measured in downgradient well MW-6 increased very slightly from 5.25 to 6.0 micrograms per liter ($\mu\text{g}/\ell$). Based on those results and the WDNR's proposed changes to the NR 140 enforcement standard (ES) for PCE from 5.0 to 20 $\mu\text{g}/\ell$, GF does not believe additional investigation or remediation is warranted at this time. GF plans to collect groundwater samples on a semi-annual basis until a stable or receding trend can be documented, and then request the WDNR to reconsider the site for closure.

Mr. Kevin McKnight
Wisconsin Department of Natural Resources
December 2, 2021

-2-

Scope of Work – July 2017 through August 2021

Since the WDNR's June 2017 letter, six rounds of groundwater samples have been collected. Sample collection dates are shown on the table below with an "x" indicating that the well was sampled on that date.

| Sample Date | Well ID | | | | | | |
|-------------|---------|------|------|------|------|------|------|
| | MW-1 | MW-2 | MW-3 | MW-4 | MW-5 | MW-6 | MW-7 |
| 07/06/17 | x | x | x | x | - | x | - |
| 01/08/18 | x | x | x | x | - | x | - |
| 08/01/18 | x | x | x | x | - | x | - |
| 03/12/19 | x | x | x | x | - | x | - |
| 09/18/19 | x | x | x | x | x | x | x |
| 12/13/19 | x | x | x | x | x | x | x |
| 07/08/20 | x | x | x | x | x | x | x |
| 08/18/21 | x | x | x | x | x | x | x |

GF's latest status report was submitted to the WDNR on January 30, 2020, and included the analytical results of groundwater samples collected through December 2019. A summary of groundwater monitoring activities conducted in the months of July 2020 and August 2021 are included in this report.

Periodic reporting of site remediation progress to the WDNR is required pursuant to ss. NR 700.11(1) and 724.13(3), Wisconsin Administrative Code. A completed certification page for the report is included with this report.

Recent Scope of Work (July 2020 – August 2021)

GF conducted groundwater monitoring activities on July 8, 2020, and August 18, 2021, that included:

- Measuring groundwater elevations in all site wells – MW-1 through MW-7.
- Measuring remediation by natural attenuation (RNA) parameters (dissolved oxygen [DO], oxidation-reduction potential [ORP], temperature, pH, and conductivity) in August 2021. RNA parameters were measured using a YSI 550 multi-meter in-situ in each of the wells prior to collecting groundwater samples (i.e., static conditions prior to purging), and then again after the well had been purged and sampled.

Mr. Kevin McKnight
Wisconsin Department of Natural Resources
December 2, 2021

-3-

- Collecting groundwater samples from each monitoring well in July 2020 and August 2020 for analyses of volatile organic compounds (VOCs). Duplicate samples were collected from MW-6 in July 2020 and from MW-1 in August 2021.
- Additional samples were collected from MW-1 in August 2021 for analysis of 1,4-dioxane (1,4-D) and per- and polyfluoroalkyl substances (PFAS).

Groundwater samples collected from each well were placed into laboratory-supplied containers with preservatives, as required for each analysis. The groundwater samples were placed in a cooler with ice and shipped via overnight courier for laboratory analysis of VOCs using EPA Method 8260B. Samples collected in July 2020 and August 2021 were submitted to Pace Analytical Laboratory in Green Bay, Wisconsin, and ALS Laboratory Group in Holland, Michigan, respectively. The August 2021 samples collected from MW-1 were also submitted for analysis of 1,4-D and PFAS using EPA Methods 8260B and 537 Modified, respectively. The laboratory reports for groundwater samples collected in July 2020 and August 2021 are included with this report as Attachment A.

Field Measurements and Analytical Results

Table 1 presents depth to water (DTW) measurements and calculated groundwater elevation data in the site wells through August 2021 and includes previous measurements since 2006. Figures 3 and 4 show the groundwater flow direction based on elevations measured in MW-1 through MW-7 on July 8, 2020, and August 18, 2021, respectively. As shown on Figures 3 and 4, the groundwater flow direction in July 2020 and August 2021 was to the north-northeast between MW-7 and MW-2, then to the northeast further downgradient. The groundwater flow directions measured in 2020 and 2021 were consistent with directions measured since MW-5 through MW-7 were installed in March 2016.

Table 2 summarizes the analytical results of groundwater samples collected through August 2021, including the analytical results of samples previously collected from the monitoring wells.

- Duplicate results are averaged in Table 2 for statistical analysis/plotting, per December 2013 Interstate Technology & Regulatory Council guidance.
- PCE was the only compound measured in July 2020 and August 2021 at concentrations above its NR 140 ES and Preventative Action Limit (PAL) of 5.0 and 0.5 micrograms per liter ($\mu\text{g}/\ell$), respectively. Notable concentrations of PCE measured in MW-1, MW-2, MW-3, and MW-6 in July 2020 and August 2021 are discussed below.

Table 3 presents the results of the RNA parameters measured through August 2021 and includes previous measurements collected from the site wells since 2013. The pH of the groundwater

Gannett Fleming

Mr. Kevin McKnight
Wisconsin Department of Natural Resources
December 2, 2021

-4-

measured in source well MW-1 and downgradient well MW-3 has ranged from 6.7 to 7.9 since RNA parameters were first measured in June 2013. DO concentrations in those wells have ranged from 0.55 to 3.82 milligrams per liter (mg/l) since June 2013. The relatively low DO concentrations and neutral pH values indicate that the aquifer would be conducive to reductive dechlorination if amendments were made to increase the activity of the microbes that facilitate the breakdown of chlorinated ethenes. However, GF does not believe that remediation is warranted at this time.

As discussed above, additional samples were collected from MW-1 in August 2021 for 1,4-D and PFAS analyses. 1,4-D was not detected ($<0.44 \mu\text{g/l}$) in the groundwater sample collected from MW-1. Eight PFAS compounds were detected in the sample collected from MW-1. See Table 4 for specific PFAS compounds and concentrations measured in the groundwater sample collected from MW-1 in August 2021. These results are further discussed below.

Discussion

PCE Plume

Since groundwater monitoring began in September 2006, PCE has been the only VOC measured at concentrations above its NR 140 ES of $5.0 \mu\text{g/l}$, with source area well MW-1 historically containing the highest PCE concentrations. As shown in Table 2, PCE was measured at concentrations above its NR 140 ES in MW-1 and MW-6 in July 2020 and August 2021 and in MW-2 and MW-3 in July 2020 only.

Since reaching its maximum concentration of $275 \mu\text{g/l}$ (the average of 255 and $295 \mu\text{g/l}$, measured in two duplicate samples) in July 2017, PCE concentrations in MW-1 have fluctuated slightly with an overall decreasing trend. The most recent sample and duplicate sample collected from MW-1 in August 2021 both contained $120 \mu\text{g/l}$ of PCE. Likewise, PCE concentrations measured in MW-2 and MW-3 were the highest in July 2020 ($5.6 \mu\text{g/l}$) and September 2019 ($81.4 \mu\text{g/l}$), respectively, but then decreased to 2.7 and $4.2 \mu\text{g/l}$ in August 2021. Water levels rose about 3.4 to 4.7 feet since January 2018, when the groundwater elevation was at a relative minimum, with the elevations measured in December 2019, July 2020, and August 2021 near historical maximum elevations. A chart showing the groundwater elevations and PCE concentrations measured in MW-1 since 2006 is included with this report as Attachment B. Figure 5 shows the estimated extent of PCE in the groundwater at concentrations at or above its NR 140 ES/PAL of $5/0.5 \mu\text{g/l}$ on August 18, 2021.

GF believes the elevated PCE concentrations measured in the groundwater samples collected from MW-1 and other wells in September and December 2019 were due to the high water table coming into contact with PCE-impacted soil. However, as stated in our January 2020 status report, we believe the overall increase in dissolved-phase PCE concentrations was limited and that the residual mass of PCE in the soil was/is relatively small. That belief is supported by the

Mr. Kevin McKnight
Wisconsin Department of Natural Resources
December 2, 2021

-5-

overall decreasing trend in PCE concentrations measured in source area wells MW-1 through MW-3 between December 2019 and August 2021, even though the water table remained at/near its historical maximum. Additionally, the hydraulic conductivity of the aquifer is relatively low (2.8×10^{-5} cm/sec) and inhibits the migration of the PCE in the groundwater, as evidenced by the relatively limited extent of PCE at concentrations above the NR 140 ES in the groundwater (about 70 feet) since the dry-cleaning facility associated with this property closed in 1994. There are no potable or municipal wells within 1,200 feet of the site, and the downgradient edge of the PCE plume is over 150 feet from the property line. Based on PCE concentrations measured in the soil and groundwater, GF does not believe there is enough mass of PCE in the source area to cause it to migrate offsite without being diluted to concentrations below the NR 140 PALs.

PCE concentrations in downgradient well MW-6 have increased slightly since it was first sampled in April 2016, with the sample collected in August 2021 containing $6.0 \mu\text{g/l}$. However, the increase in PCE concentrations in MW-6 has been relatively slow, and we believe the increasing trend will reverse itself as groundwater with lower PCE concentrations, as measured in MW-2 and MW-3 in August 2021, migrates further downgradient. Additionally, the WDNR is in the process of revising the NR 140 ES and PAL for PCE from $5.0/0.5 \mu\text{g/l}$ to $20/2.0 \mu\text{g/l}$. After that revision is promulgated, the footprint of the PCE plume at concentrations above the NR 140 ES/PAL will dramatically decrease in size, as shown on Figure 6. For those reasons, GF does not believe additional investigation or remediation of the PCE plume is warranted at this time.

PFAS and 1,4-Dioxane Results

The PFAS and 1,4-D samples were collected from MW-1 because it has historically contained the highest concentrations of PCE, indicating it is in the source area where hazardous compounds associated with the former dry cleaner were released.

1,4-D was not detected in the groundwater sample collected from MW-1 in August 2021. Of the eight PFAS compounds detected in the sample from MW-1, only two compounds were measured at concentrations above their proposed NR 140 PAL recommended by the Wisconsin Department of Health Services: PFOA and PFOS were at 5.8 and 9.7 nanograms per liter (ng/l) above their combined proposed NR 140 PAL of 2.0 ng/l but below their combined proposed NR 140 ES of 20 ng/l. All other PFAS compounds were measured at concentrations far below their proposed NR 140 ES and PALs. See Table 4 for the PFAS compounds and concentrations measured in the groundwater sample collected from MW-1 in August 2021. Based on the absence of 1,4-D and the relatively low concentrations of PFAS measured in source area well MW-1, GF does not believe that additional sampling for 1,4-D or PFAS is warranted.

Gannett Fleming

Mr. Kevin McKnight
Wisconsin Department of Natural Resources
December 2, 2021

-6-

Closing

Because of the low groundwater velocity and the relatively wide fluctuations in the water table elevation over the past 10 years, GF proposes to monitor PCE concentrations in the groundwater on a semiannual basis going forward. Groundwater monitoring activities will continue until stable/receding trends in PCE concentrations have been established. The next groundwater sampling events are scheduled for spring and fall 2022. A groundwater monitoring report will be submitted to the WDNR after the fall 2022 sample results have been received and evaluated. In the meantime, please let us know if you have any questions or need additional information to complete your review of this report.

Sincerely,

GANNETT FLEMING, INC.

Chelsea Payne

Chelsea J. Payne, P.G.
Project Geologist

Anthony W. Miller

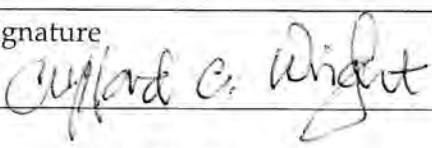
Anthony W. Miller, P.S.S.
Senior Project Manager

AWM/jec
Enc.

ecc: Ryan Eley (Livesey Company, LLC)

ENGINEERING AND HYDROGEOLOGIST CERTIFICATIONS

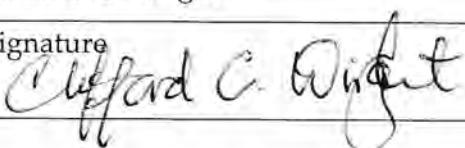
I hereby certify that I am a registered professional engineer in the State of Wisconsin, registered in accordance with the requirements of ch. A-E 4, Wis. Adm. Code; that this document has been prepared in accordance with the rules of Professional Conduct in ch. A-E 8, Wis. Adm. Code; and that, to the best of my knowledge, all information contained in this document is correct and the document was prepared in compliance with all applicable requirements in chs. NR 700 to 726, Wis. Adm. Code.

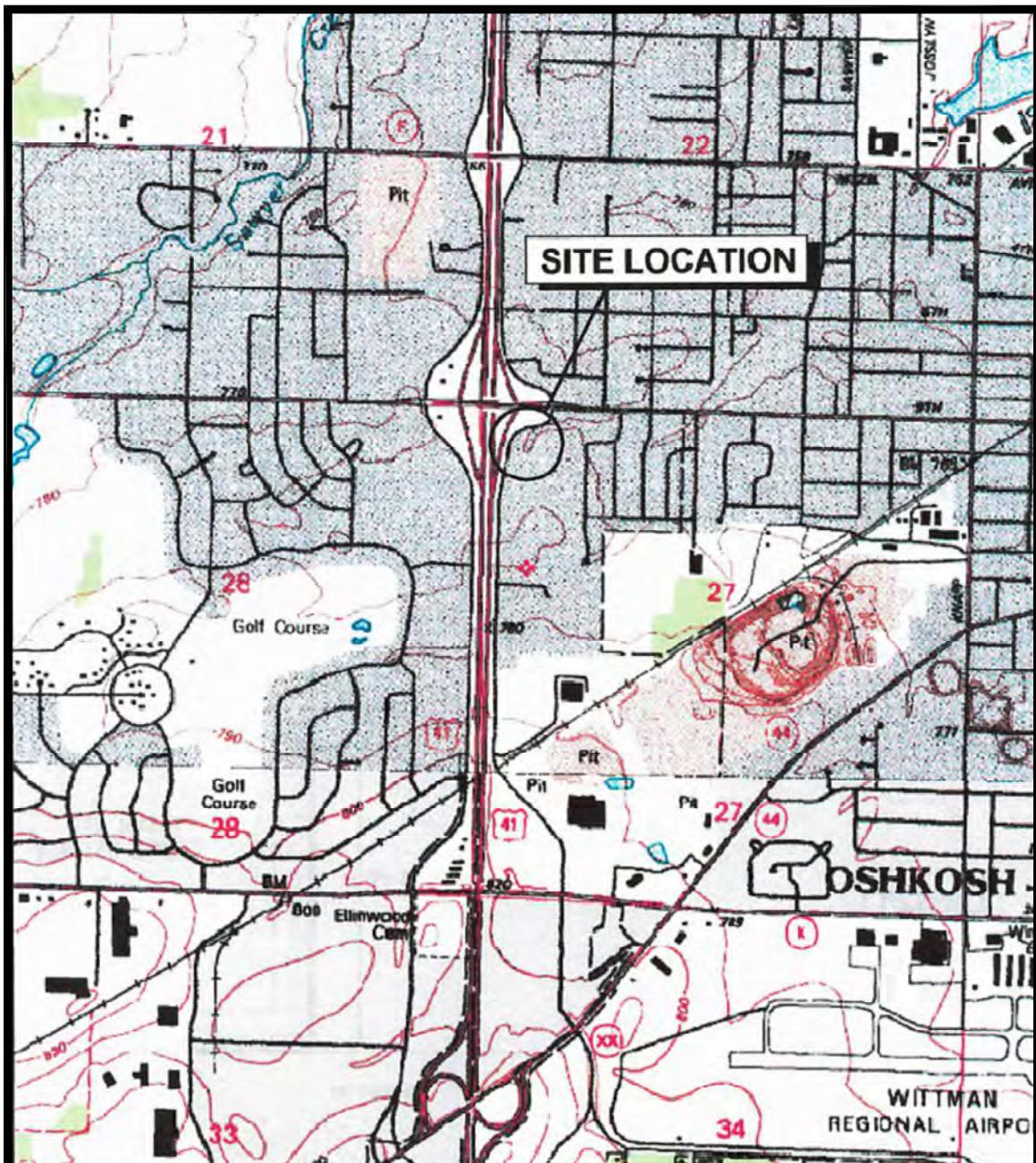
| | |
|--|-------------------------------------|
| Print Name Clifford C. Wright | Title Project Engineer/Geologist |
| Signature  | Date 12-1-2021 |

P.E. Seal for E-31265:



I hereby certify that I am a hydrogeologist as that term is defined in s. NR 712.03(1), Wis. Adm. Code, am registered in accordance with the requirements of ch. GHSS 2, Wis. Adm. Code, or licensed in accordance with the requirements of ch. GHSS 3, Wis. Adm. Code, and that, to the best of my knowledge, all information contained in this document is correct and the document was prepared in compliance with all applicable requirements in chs. NR 700 to 726, Wis. Adm. Code.

| | |
|--|-------------------------------------|
| Print Name Clifford C. Wright | Title Project Engineer/Geologist |
| Signature  | Date 12-1-2021 |



SCALE: 1 INCH ~ 1800 FEET

7.5 MIN TOPOGRAPHIC MAP
OSHKOSH, WISCONSIN
1992



LOCATION MAP
KOELLER CENTER—OSHKOSH
KOELLER ONE, LLC
OSHKOSH, WISCONSIN



1 INCH ~ 375 FEET

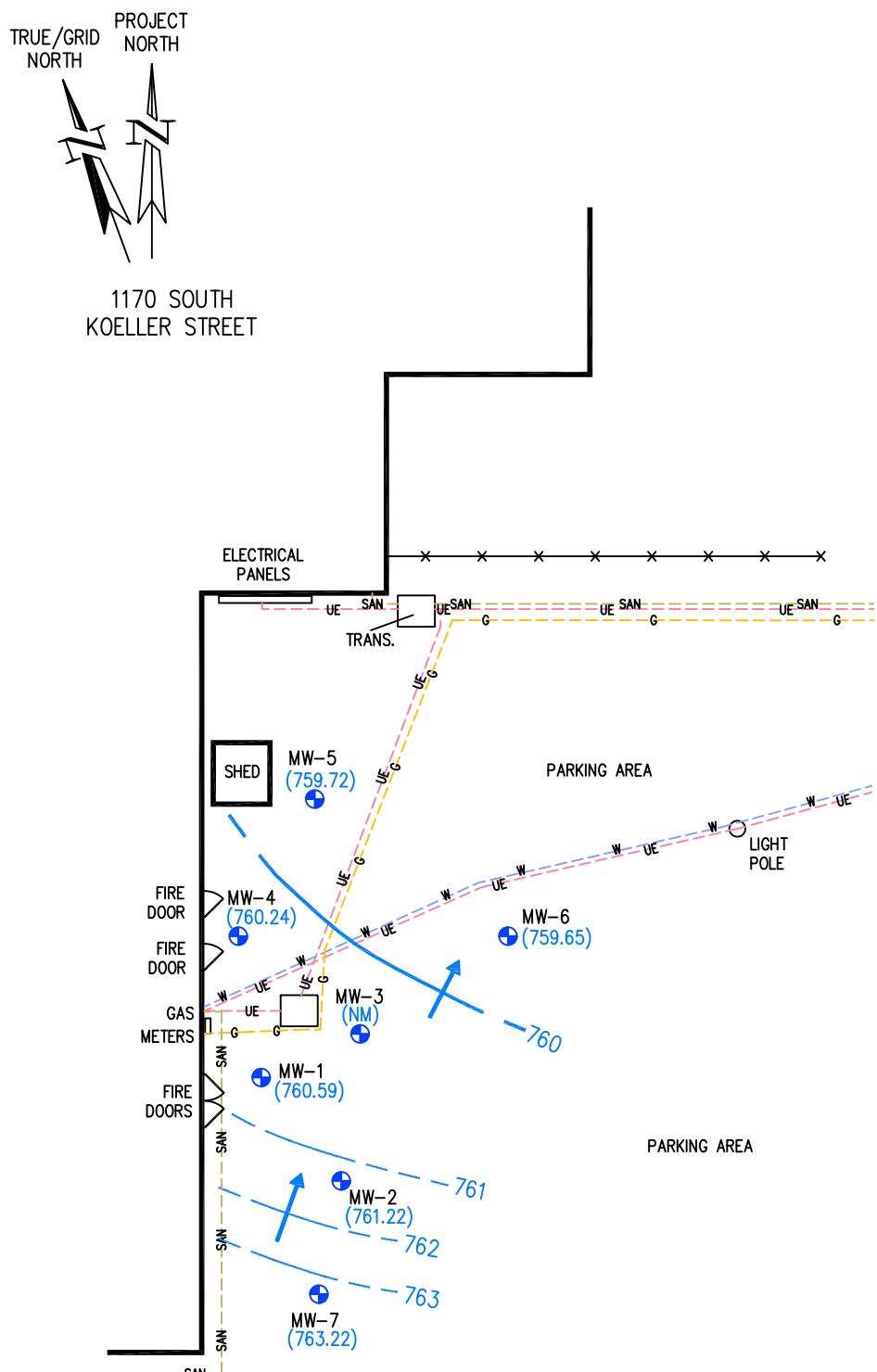
GOOGLE EARTH - 06/15



AERIAL PHOTO OF
PARCEL AND AREA
OF INVESTIGATION
KOELLER CENTER—OSHKOSH
KOELLER ONE, LLC
OSHKOSH, WISCONSIN

Gannett Fleming

FIGURE 3



LEGEND

- This legend identifies various features and infrastructure elements shown in the map:

 - 760 → GROUNDWATER CONTOUR (DASHED WHERE INFERRED)
 - ← GROUNDWATER FLOW DIRECTION (07/08/20)
 - MONITORING WELL
 - *— FENCE
 - G— UNDERGROUND NATURAL GAS
 - UE— UNDERGROUND ELECTRICAL
 - W— UNDERGROUND WATER LINE
 - SAN— UNDERGROUND SANITARY SEWER

NOTES

- 1. LOCATIONS OF UNDERGROUND UTILITIES ARE APPROXIMATE.**

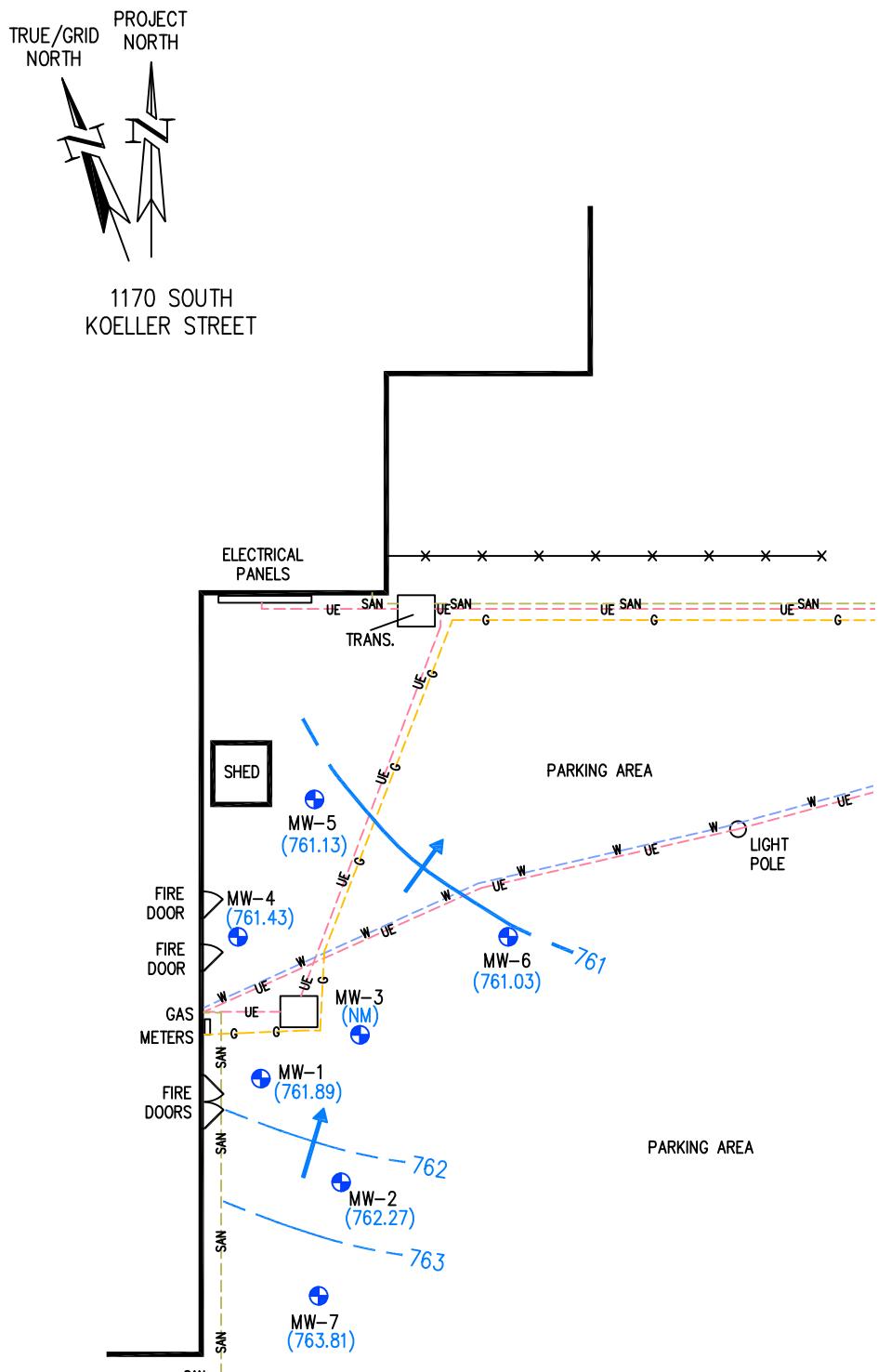
GROUNDWATER CONTOUR MAP (JULY 2020)

KOELLER CENTER—OSHKOSH
KOELLER ONE, LLC.
OSHKOSH, WISCONSIN

A horizontal scale bar with tick marks at 0 and 40. The text "Approximate Scale In Feet" is written below it.

Gannett Fleming

FIGURE 4



LEGEND

- This legend provides key information for interpreting the map:

 - 760** — GROUNDWATER CONTOUR (DASHED WHERE INFERRED)
 - ←** — GROUNDWATER FLOW DIRECTION (08/18/2021)
 - MONITORING WELL** (represented by a circle with a cross)
 - FENCE** (represented by a line with an asterisk)
 - G** — UNDERGROUND NATURAL GAS
 - UE** — UNDERGROUND ELECTRICAL
 - W** — UNDERGROUND WATER LINE
 - SAN** — UNDERGROUND SANITARY SEWER

NOTES

- ## 1. LOCATIONS OF UNDERGROUND UTILITIES ARE APPROXIMATE.

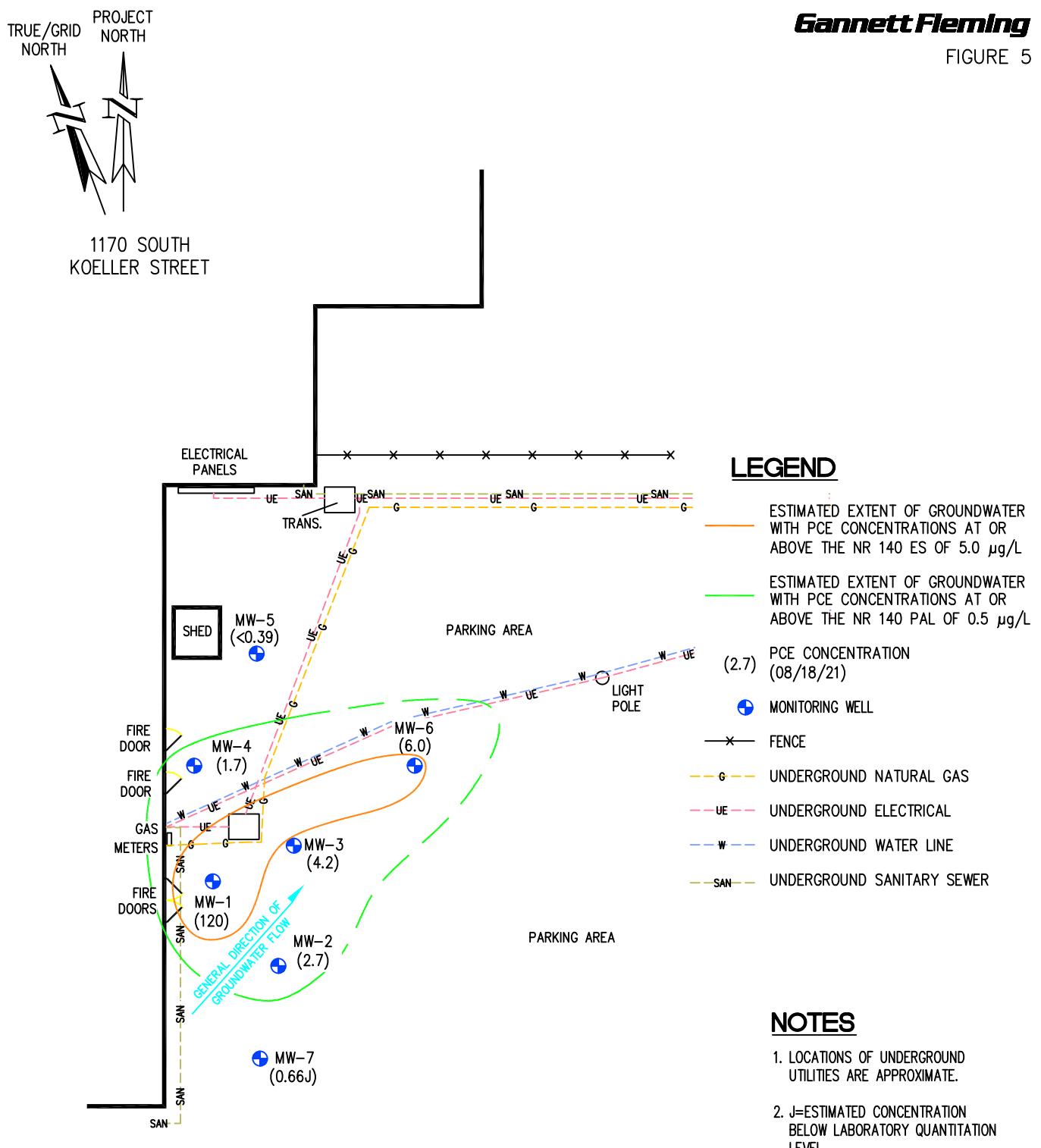
A horizontal scale bar with tick marks at 0 and 40. The segment between them is divided into four equal parts by three internal tick marks. Below the scale bar, the text "Approximate Scale In Feet" is centered.

GROUNDWATER
CONTOUR MAP
(AUGUST 2021)

KOELLER CENTER—OSHKOSH
KOELLER ONE, LLC.
OSHKOSH, WISCONSIN

Gannett Fleming

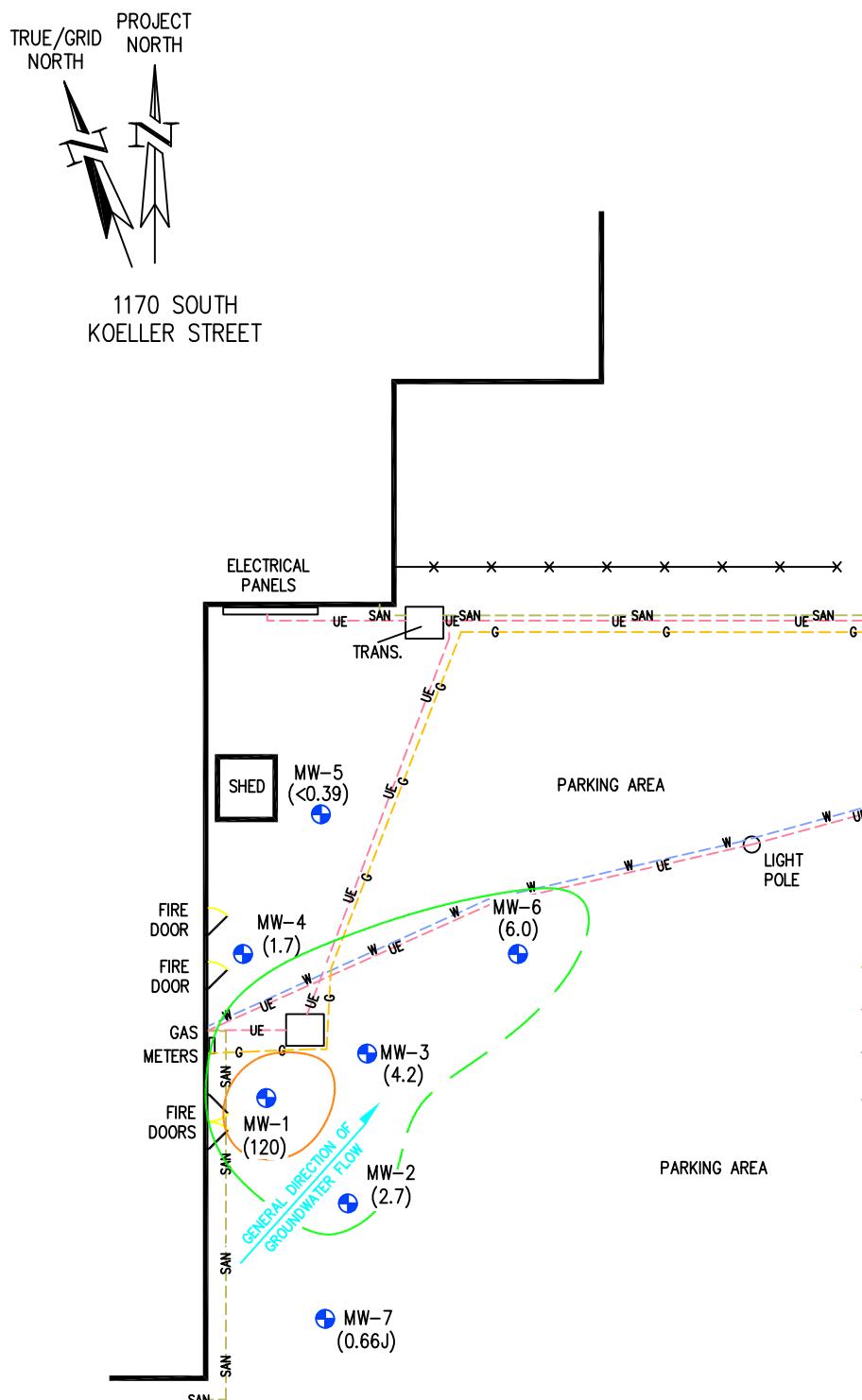
FIGURE 5



PCE CONCENTRATIONS IN GROUNDWATER (AUGUST 2021)

KOELLER CENTER—OSHKOSH
KOELLER ONE, LLC.
OSHKOSH, WISCONSIN

092021
awm_R47358-003_2021_01MR_F05

**LEGEND**

- ESTIMATED EXTENT OF GROUNDWATER WITH PCE CONCENTRATIONS AT OR ABOVE THE PROPOSED NR 140 ES OF 20 µg/L
- ESTIMATED EXTENT OF GROUNDWATER WITH PCE CONCENTRATIONS AT OR ABOVE THE PROPOSED NR 140 PAL OF 2.0 µg/L
- (2.7) PCE CONCENTRATION (08/18/21)
- MONITORING WELL
- FENCE
- UNDERGROUND NATURAL GAS
- UNDERGROUND ELECTRICAL
- UNDERGROUND WATER LINE
- UNDERGROUND SANITARY SEWER

NOTES

- LOCATIONS OF UNDERGROUND UTILITIES ARE APPROXIMATE.
- J=ESTIMATED CONCENTRATION BELOW LABORATORY QUANTITATION LEVEL.
- PROPOSED NR 140 ES AND PAL OF 20 AND 2.0 µg/L RESPECTIVELY, WERE RECOMMENDED BY THE WISCONSIN DEPARTMENT OF HEALTH IN JUNE 2019.

ESTIMATED EXTENT OF PCE USING PROPOSED NR 140 ES & PAL (AUGUST 2021)

KOELLER CENTER-OSHKOSH
KOELLER ONE, LLC.
OSHKOSH, WISCONSIN

0 40
Approximate Scale In Feet

KOELLER ONE, LLC
KOELLER SHOPPING CENTER
OSHKOSH, WISCONSIN

TABLE 1

WATER LEVEL ELEVATION DATA (MW-1 THROUGH MW-7)

| Well ID | Monitoring Well ID and Reference Information | | | | | | |
|---|--|---------|-------------------|---------|---------|---------|---------|
| | MW-1 | MW-2 | MW-3 | MW-4 | MW-5 | MW-6 | MW-7 |
| Date Well Installed | 9/7/06 | 8/14/08 | 8/14/08 | 5/23/13 | 3/28/16 | 3/28/16 | 3/28/16 |
| Top of Casing Elevation (ft MSL) ^(1,2) | 774.19 | 773.87 | 774.17 | 774.25 | 774.10 | 773.57 | 774.04 |
| Top of Screen Elevation (ft MSL) | 757.45 | 762.13 | 762.43 | 762.51 | 765.8 | 764.18 | 763.91 |
| Bottom of Screen Elevation (ft MSL) | 747.45 | 747.13 | 747.43 | 747.51 | 750.8 | 749.18 | 748.91 |
| Measurement Date | Depth to Water (feet below top of casing) | | | | | | |
| 10/31/06 | 16.44 | NI | NI | NI | NI | NI | NI |
| 01/10/07 | 15.82 | NI | NI | NI | NI | NI | NI |
| 05/17/07 | 14.62 | NI | NI | NI | NI | NI | NI |
| 08/14/08 | 14.45 | NI | NI | NI | NI | NI | NI |
| 08/21/08 | 15.20 | 14.42 | 14.97 | NI | NI | NI | NI |
| 01/28/09 | 17.71 | 17.10 | 17.56 | NI | NI | NI | NI |
| 04/08/09 | 12.71 | 11.81 | 12.36 | NI | NI | NI | NI |
| 06/14/13 ⁽³⁾ | 16.78 | 11.80 | 12.56 | 13.34 | NI | NI | NI |
| 09/12/13 | 15.11 | 14.55 | 15.25 | 15.35 | NI | NI | NI |
| 03/14/14 | 16.37 | 16.05 | 16.60 | 16.56 | NI | NI | NI |
| 07/10/14 | 11.84 | 11.10 | 11.62 | 12.20 | NI | NI | NI |
| 11/19/14 | 14.80 | 14.27 | 14.84 | 15.12 | NI | NI | NI |
| 12/04/14 | 14.59 | 13.95 | 14.61 | 14.88 | NI | NI | NI |
| 03/28/16 | 11.05 | 10.45 | 10.78 | 11.41 | 11.51 | 11.06 | 13.51 |
| 04/12/16 | 11.12 | 10.33 | 10.54 | 11.48 | 11.62 | 11.09 | 9.21 |
| 06/03/16 | 13.04 | 12.10 | NM | 13.36 | 13.52 | 13.01 | 10.79 |
| 06/22/16 | 13.39 | 12.50 | 13.23 | 13.76 | 13.95 | 13.45 | 11.10 |
| 07/14/16 | 14.28 | 13.37 | 14.18 | 14.61 | 14.78 | 14.31 | 12.26 |
| 08/01/16 | 13.85 | 12.95 | 13.76 | 14.39 | 14.48 | 13.98 | 11.47 |
| 09/01/16 | 14.89 | 13.82 | 14.76 | 14.98 | 15.41 | 14.93 | 12.18 |
| 10/05/16 | 13.58 | 12.83 | 13.52 | 14.47 | 14.23 | 13.77 | 11.15 |
| 11/04/16 | 14.47 | 13.70 | 14.55 | 14.79 | 14.94 | 14.43 | 12.68 |
| 11/11/16 | 14.90 | 14.09 | 14.89 | 15.18 | 15.21 | 14.79 | 13.39 |
| 07/06/17 | 12.47 | 11.58 | 12.30 | 12.82 | 13.01 | 12.56 | 10.42 |
| 01/08/18 | 16.61 | 16.07 | 16.61 | 16.77 | 16.72 | 16.33 | 15.75 |
| 08/01/18 | 14.21 | 13.37 | 14.05 | 14.55 | 14.79 | 14.35 | 11.95 |
| 03/12/19 | 15.10 | 14.39 | NM ⁽⁴⁾ | 15.50 | 15.63 | 15.22 | 13.74 |
| 09/18/19 | 13.65 | 12.80 | NM ⁽⁵⁾ | 14.37 | 14.42 | 13.95 | 11.10 |
| 12/13/19 | 12.54 | 11.59 | NM ⁽⁶⁾ | 13.08 | 13.31 | 12.83 | 10.08 |
| 07/08/20 | 13.60 | 12.65 | NM ⁽⁶⁾ | 14.01 | 14.38 | 13.92 | 10.82 |
| 08/18/21 | 12.30 | 11.60 | NM ⁽⁶⁾ | 12.82 | 12.97 | 12.54 | 10.23 |

TABLE 1

WATER LEVEL ELEVATION DATA (MW-1 THROUGH MW-7)

| Well ID | Monitoring Well ID and Reference Information | | | | | | |
|---|--|---------------|-------------------|---------------|---------------|---------------|---------------|
| | MW-1 | MW-2 | MW-3 | MW-4 | MW-5 | MW-6 | MW-7 |
| Date Well Installed | 9/7/06 | 8/14/08 | 8/14/08 | 5/23/13 | 3/28/16 | 3/28/16 | 3/28/16 |
| Top of Casing Elevation (ft MSL)^(1,2) | 774.19 | 773.87 | 774.17 | 774.25 | 774.10 | 773.57 | 774.04 |
| Top of Screen Elevation (ft MSL) | 757.45 | 762.13 | 762.43 | 762.51 | 765.8 | 764.18 | 763.91 |
| Bottom of Screen Elevation (ft MSL) | 747.45 | 747.13 | 747.43 | 747.51 | 750.8 | 749.18 | 748.91 |
| Measurement Date | Water Elevation (ft MSL) | | | | | | |
| 10/31/06 | 757.75 | NI | NI | NI | NI | NI | NI |
| 01/10/07 | 758.37 | NI | NI | NI | NI | NI | NI |
| 05/17/07 | 759.57 | NI | NI | NI | NI | NI | NI |
| 08/14/08 | 759.74 | NI | NI | NI | NI | NI | NI |
| 08/21/08 | 758.99 | 759.45 | 759.20 | NI | NI | NI | NI |
| 01/28/09 | 756.48 | 756.77 | 756.61 | NI | NI | NI | NI |
| 04/08/09 | 761.48 | 762.06 | 761.81 | NI | NI | NI | NI |
| 06/14/13 ⁽³⁾ | 757.41 | 762.07 | 761.61 | 760.91 | NI | NI | NI |
| 09/12/13 | 759.08 | 759.32 | 758.92 | 758.90 | NI | NI | NI |
| 03/14/14 | 757.82 | 757.82 | 757.57 | 757.69 | NI | NI | NI |
| 07/10/14 | 762.35 | 762.77 | 762.55 | 762.05 | NI | NI | NI |
| 11/19/14 | 759.39 | 759.60 | 759.33 | 759.13 | NI | NI | NI |
| 12/04/14 | 759.60 | 759.92 | 759.56 | 759.37 | NI | NI | NI |
| 03/28/16 | 763.14 | 763.42 | 763.39 | 762.84 | 762.59 | 762.51 | 760.53 |
| 04/12/16 | 763.07 | 763.54 | 763.63 | 762.77 | 762.48 | 762.48 | 764.83 |
| 06/03/16 | 761.15 | 761.77 | NM | 760.89 | 760.58 | 760.56 | 763.25 |
| 06/22/16 | 760.80 | 761.37 | 760.94 | 760.49 | 760.15 | 760.12 | 762.94 |
| 07/14/16 | 759.91 | 760.50 | 759.99 | 759.64 | 759.32 | 759.26 | 761.78 |
| 08/01/16 | 760.34 | 760.92 | 760.41 | 759.86 | 759.62 | 759.59 | 762.57 |
| 09/01/16 | 759.30 | 760.05 | 759.41 | 759.27 | 758.69 | 758.64 | 761.86 |
| 10/05/16 | 760.61 | 761.04 | 760.65 | 759.78 | 759.87 | 759.80 | 762.89 |
| 11/04/16 | 759.72 | 760.17 | 759.62 | 759.46 | 759.16 | 759.14 | 761.36 |
| 11/11/16 | 759.29 | 759.78 | 759.28 | 759.07 | 758.89 | 758.78 | 760.65 |
| 07/06/17 | 761.72 | 762.29 | 761.87 | 761.43 | 761.09 | 761.01 | 763.62 |
| 01/08/18 | 757.58 | 757.80 | 757.56 | 757.48 | 757.38 | 757.24 | 758.29 |
| 08/01/18 | 759.98 | 760.50 | 760.12 | 759.70 | 759.31 | 759.22 | 762.09 |
| 03/12/19 | 759.09 | 759.48 | NM ⁽⁴⁾ | 758.75 | 758.47 | 758.35 | 760.30 |
| 09/18/19 | 760.54 | 761.07 | NM ⁽⁵⁾ | 759.88 | 759.68 | 759.62 | 762.94 |
| 12/13/19 | 761.65 | 762.28 | NM ⁽⁶⁾ | 761.17 | 760.79 | 760.74 | 763.96 |
| 07/08/20 | 760.59 | 761.22 | NM ⁽⁶⁾ | 760.24 | 759.72 | 759.65 | 763.22 |
| 08/18/21 | 761.89 | 762.27 | NM(6) | 761.43 | 761.13 | 761.03 | 763.81 |

TABLE 1

WATER LEVEL ELEVATION DATA (MW-1 THROUGH MW-7)

NOTES:

Site datum = feet above mean sea level (ft MSL).

USGS Registered Benchmark = 776.04 ft MSL, top of nut of fire hydrant in front of Mitchell Insurance at 1746 W. 9th Ave. (Source City of Oshkosh Engineering Department, 920-236-5065.)

Local Benchmark = 776.11 ft MSL, top of nut of fire hydrant, 240 feet east of MW-1.

NI = Not installed.

NM = Not measured.

FOOTNOTES:

(1) Top of casing elevations for MW-1 through MW-4 based on 03/14/14 survey using MW-1= 774.19 ft MSL as benchmark.

(2) MW-5 through MW-7 surveyed on 3/28/16 using MW-1 top of casing as benchmark. The top nut of the local benchmark fire hydrant was measured at 776.10 ft MSL on 3/28/16 and 4/12/16.

(3) The relatively deep groundwater elevation measured in MW-1 on 06/14/13 was likely due to field error caused by not allowing the water table to rise and stabilize after removing the air-tight cap. This phenomenon was discussed in more detail in Gannett Fleming's November 2013 *Site Investigation Report*.

(4) On 03/12/19, the MW-3 top of casing was bent so that a water level probe could not reach past 8 inches below top of casing.

(5) On 09/18/19, the MW-3 top of casing was filled with sand and debris so DTW was not measured.

(6) The top of casing is damaged and therefore an accurate DTW could not be measured.

KOELLER ONE, LLC
KOELLER SHOPPING CENTER
OSHKOSH, WISCONSIN

TABLE 2

SUMMARY OF GROUNDWATER MONITORING ANALYTICAL RESULTS

| Well ID | Concentration ($\mu\text{g/l}$) and Results Qualifier(s) for Detected Volatile Organic Compounds (VOCs) | | | | | | | | | Comments Footnotes |
|-------------|---|-------------------|-------------------------|------------------------|---------|--------------|---------|---------|---------|--------------------|
| | Tetrachloroethylene | Trichloroethylene | Dichlorodifluoromethane | 1,2,4-Trimethylbenzene | Benzene | Ethylbenzene | Xylenes | Styrene | Toluene | |
| Sample Date | | | | | | | | | | |
| NR 140 PAL | 0.5 | 0.5 | 200 | 96 | 0.5 | 140 | 1,000 | 10 | 200 | |
| NR 140 ES | 5.0 | 5.0 | 1,000 | 480 | 5.0 | 700 | 10,000 | 100 | 1,000 | |
| MW-1 | | | | | | | | | | |
| 09/06/06 | <0.50 | <0.50 | NA | NA | <0.50 | <0.50 | <1.0 | NA | <0.50 | |
| 10/31/06 | 2.95 | 0.27 | 0.90 | <0.15 | <0.15 | <0.10 | <0.50 | <0.10 | <0.40 | |
| 01/10/07 | 14.8 | <0.20 | 1.50 | <0.15 | <0.15 | 0.13 J | <0.50 | <0.10 | <0.40 | |
| 05/17/07 | 12.0 | <0.20 | 1.30 | <0.20 | <0.20 | <0.10 | <0.60 | <0.10 | <0.40 | |
| 08/21/08 | 54.4 | <0.40 | 1.95 | <0.20 | <0.20 | <0.20 | <0.60 | <0.10 | <0.40 | |
| 01/28/09 | 36.4 | <0.40 | 0.82 J | <0.20 | <0.20 | <0.20 | <0.60 | <0.10 | <0.40 | (1) |
| 04/08/09 | 28.7 | <0.40 | 1.22 | <0.20 | <0.20 | <0.20 | <0.60 | <0.10 | <0.40 | |
| 06/14/13 | 89.4 | <0.43 | <0.40 | <0.57 | <0.50 | <0.50 | <0.82 | <0.35 | <0.44 | |
| 09/12/13 | 94.4 | <0.43 | <0.40 | <0.57 | <0.50 | <0.50 | <0.82 | <0.35 | <0.44 | |
| 03/13/14 | 50.1 | <0.36 | <0.40 | <0.50 | <0.50 | <0.50 | <0.82 | <0.35 | <0.44 | |
| 07/10/14 | 92.3 | <0.33 | <0.20 | <0.50 | <0.50 | <0.50 | <0.82 | <0.50 | <0.50 | |
| 12/04/14 | 110 | <0.33 | <0.20 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 04/12/16 | 185 | <0.33 | <0.20 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 11/11/16 | 213 | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 07/06/17 | 275 | 0.58 U | 0.39 U | 0.85 U | 0.85 U | 0.85 U | 2.60 U | 0.85 U | 0.85 U | Dup |
| 01/08/18 | 192 | <0.83 | <0.56 | <1.2 | <1.2 | <1.2 | <3.7 | <1.2 | <1.2 | |
| 08/01/18 | 162 | <0.64 | <1.2 | <2.1 | <0.62 | <0.55 | <1.85 | <1.2 | <0.43 | |
| 03/12/19 | 125 | <0.64 | <1.2 | <2.1 | <0.62 | <0.55 | <1.85 | <1.2 | <0.43 | Dup |
| 09/18/19 | 192 | <0.64 | <1.2 | <2.1 | <0.62 | <0.55 | <1.85 | <1.2 | <0.43 | |
| 12/13/19 | 184 | <0.64 | <1.2 | <2.1 | <0.62 | <0.55 | <1.85 | <1.2 | <0.43 | |
| 07/08/20 | 181 | <0.64 | <1.2 | <2.1 | <0.62 | <0.80 | <1.85 | <7.5 | <0.67 | |
| 08/18/21 | 120 | <0.43 | <0.68 | <0.45 | <0.46 | <0.34 | <0.81 | <0.33 | <0.45 | Dup |
| MW-2 | | | | | | | | | | |
| 08/21/08 | 0.88 J | <0.40 | 1.62 | <0.20 | <0.20 | <0.20 | <0.60 | <0.10 | <0.40 | |
| 01/28/09 | 0.53 J | <0.40 | 1.15 | <0.20 | <0.20 | <0.20 | <0.60 | <0.10 | <0.40 | (1) |
| 04/08/09 | 0.87 J | <0.40 | 1.12 | <0.20 | <0.20 | <0.20 | <0.60 | <0.10 | <0.40 | |
| 06/14/13 | 1.5 | <0.43 | 0.48 J | <0.57 | <0.50 | <0.50 | <0.82 | <0.35 | <0.44 | |
| 09/12/13 | 1.0 | <0.43 | 0.45 J | <0.57 | <0.50 | <0.50 | <0.82 | <0.35 | <0.44 | |
| 03/13/14 | 0.88 J | <0.36 | 0.45 J | <0.50 | <0.50 | <0.50 | <0.82 | <0.35 | <0.44 | |

TABLE 2

SUMMARY OF GROUNDWATER MONITORING ANALYTICAL RESULTS

| Well ID | Concentration ($\mu\text{g/l}$) and Results Qualifier(s) for Detected Volatile Organic Compounds (VOCs) | | | | | | | | | Comments(Footnotes) |
|-------------|---|-------------------|-------------------------|------------------------|-------------|--------------|---------------|------------|--------------|---------------------|
| Sample Date | Tetrachloroethylene | Trichloroethylene | Dichlorodifluoromethane | 1,2,4-Trimethylbenzene | Benzene | Ethylbenzene | Xylenes | Styrene | Toluene | |
| NR 140 PAL | 0.5 | 0.5 | 200 | 96 | 0.5 | 140 | 1,000 | 10 | 200 | |
| NR 140 ES | 5.0 | 5.0 | 1,000 | 480 | 5.0 | 700 | 10,000 | 100 | 1,000 | |
| 07/10/14 | 2.0 | <0.33 | 0.43 J | <0.50 | <0.50 | <0.50 | <0.82 | <0.50 | <0.50 | |
| 12/04/14 | 1.1 | <0.33 | 0.50 J | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 04/12/16 | 0.95 J | <0.33 | 0.44 J | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 11/11/16 | 1.6 | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 07/06/17 | 1.9 | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 01/08/18 | 2.0 | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 08/01/18 | 2.5 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | (2) |
| 03/12/19 | 1.8 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |
| 09/18/19 | 2.6 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |
| 12/13/19 | 5.1 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |
| 07/08/20 | 5.6 | <0.26 | <0.50 | <0.84 | <0.25 | <0.32 | <0.73 | <3.0 | <0.27 | |
| 08/18/21 | 2.7 | <0.43 | <0.68 | <0.45 | <0.46 | <0.34 | <0.81 | <0.33 | <0.45 | |
| MW-3 | | | | | | | | | | |
| 08/21/08 | 4.80 | <0.40 | 0.36 J | 0.22 J | <u>0.88</u> | 1.09 | 4.39 | 0.14 J | 2.21 | |
| 01/28/09 | 3.80 | <0.40 | <0.30 | 0.21 J | 0.27 | 0.72 | 2.62 | 0.12 J | 0.65 J J | |
| 04/08/09 | 7.12 | <0.40 | <0.30 | <0.20 | <0.20 | 0.36 J | 1.34 J | <0.10 | <0.40 | |
| 06/14/13 | 8.9 | <0.43 | <0.40 | <0.57 | <0.50 | <0.50 | <0.82 | <0.35 | <0.44 | |
| 09/12/13 | 4.9 | <0.43 | <0.40 | <0.57 | <0.50 | <0.50 | <0.82 | <0.35 | <0.44 | |
| 03/13/14 | 10 | <0.36 | <0.40 | <0.50 | <0.50 | <0.50 | <0.82 | <0.35 | <0.44 | |
| 07/10/14 | 22.9 | <0.33 | <0.20 | <0.50 | <0.50 | <0.50 | <0.82 | <0.50 | <0.50 | |
| 12/04/14 | 37.4 | <0.33 | <0.20 | <0.50 | <0.50 | <0.50 | <0.82 | <0.50 | <0.50 | Dup |
| 04/12/16 | 18.6 | <0.33 | <0.20 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 11/11/16 | 10.1 | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 07/06/17 | 28.7 | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 01/08/18 | 35.8 | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | Dup |
| 08/01/18 | 29.1 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |
| 03/12/19 | 33.1 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |
| 09/18/19 | 81.4 | 1.1 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |
| 12/13/19 | 45.2 | 0.63 J | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | Dup |
| 07/08/20 | 63.0 | 0.30 J | <0.50 | <0.84 | <0.25 | <0.32 | <0.73 | <3.0 | <0.27 | |
| 08/18/21 | 4.2 | <0.43 | <0.68 | <0.45 | <0.46 | <0.34 | <0.81 | <0.33 | <0.45 | |

TABLE 2
SUMMARY OF GROUNDWATER MONITORING ANALYTICAL RESULTS

| Well ID | Concentration ($\mu\text{g/l}$) and Results Qualifier(s) for Detected Volatile Organic Compounds (VOCs) | | | | | | | | | Comments(Footnotes) |
|-------------|---|-------------------|-------------------------|------------------------|---------|--------------|---------|---------|---------|---------------------|
| Sample Date | Tetrachloroethylene | Trichloroethylene | Dichlorodifluoromethane | 1,2,4-Trimethylbenzene | Benzene | Ethylbenzene | Xylenes | Styrene | Toluene | |
| NR 140 PAL | 0.5 | 0.5 | 200 | 96 | 0.5 | 140 | 1,000 | 10 | 200 | |
| NR 140 ES | 5.0 | 5.0 | 1,000 | 480 | 5.0 | 700 | 10,000 | 100 | 1,000 | |
| MW-4 | | | | | | | | | | |
| 06/14/13 | <0.47 | <0.43 | <0.40 | <0.57 | <0.50 | <0.50 | <0.82 | <0.35 | <0.44 | Dup |
| 09/12/13 | 0.82 J | <0.43 | <0.40 | <0.57 | <0.50 | <0.50 | <0.82 | <0.35 | <0.44 | Dup |
| 03/13/14 | <0.47 | <0.36 | <0.40 | <0.50 | <0.50 | <0.50 | <0.82 | <0.35 | <0.44 | Dup |
| 07/10/14 | 1.3 | <0.33 | <0.20 | <0.50 | <0.50 | <0.50 | <0.82 | <0.50 | <0.50 | |
| 12/04/14 | 0.77 J | <0.33 | <0.20 | <0.50 | <0.50 | <0.50 | <0.82 | <0.50 | <0.50 | |
| 04/12/16 | 1.2 | <0.33 | <0.20 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 11/11/16 | 0.98 J | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 07/06/17 | 1.7 | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 01/08/18 | <0.50 | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 08/01/18 | 1.3 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | (2) |
| 03/12/19 | 0.96 J | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |
| 09/18/19 | 1.5 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |
| 12/13/19 | 1.5 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |
| 07/08/20 | 1.3 | <0.26 | <0.50 | <0.84 | <0.25 | <0.32 | <0.73 | <3.0 | <0.27 | |
| 08/18/21 | 1.7 | <0.43 | <0.68 | <0.45 | <0.46 | <0.34 | <0.81 | <0.33 | <0.45 | |
| MW-5 | | | | | | | | | | |
| 04/12/16 | <0.47 | <0.33 | <0.20 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 11/11/16 | <0.50 | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 09/18/19 | <0.33 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | 0.26 J | |
| 12/13/19 | <0.33 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |
| 07/08/20 | <0.33 | <0.26 | <0.50 | <0.84 | <0.25 | <0.32 | <0.73 | <3.0 | <0.27 | |
| 08/18/21 | <0.39 | <0.43 | <0.68 | <0.45 | <0.46 | <0.34 | <0.81 | <0.33 | <0.45 | |
| MW-6 | | | | | | | | | | |
| 04/12/16 | <0.47 | <0.33 | <0.20 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | Dup |
| 11/11/16 | 0.54 J | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 07/06/17 | 0.64 J | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 01/08/18 | 2.6 | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | |
| 08/01/18 | 2.15 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | Dup ⁽²⁾ |
| 03/12/19 | 3.0 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |
| 09/18/19 | 4.8 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |
| 12/13/19 | 4.7 | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | |

TABLE 2
SUMMARY OF GROUNDWATER MONITORING ANALYTICAL RESULTS

| Well ID | Concentration ($\mu\text{g/l}$) and Results Qualifier(s) for Detected Volatile Organic Compounds (VOCs) | | | | | | | | | Comments(Footnotes) | |
|-------------|---|-------------------|-------------------------|------------------------|------------|--------------|---------------|------------|--------------|-------------------------|--|
| Sample Date | Tetrachloroethylene | Trichloroethylene | Dichlorodifluoromethane | 1,2,4-Trimethylbenzene | Benzene | Ethylbenzene | Xylenes | Styrene | Toluene | | |
| NR 140 PAL | 0.5 | 0.5 | 200 | 96 | 0.5 | 140 | 1,000 | 10 | 200 | | |
| NR 140 ES | 5.0 | 5.0 | 1,000 | 480 | 5.0 | 700 | 10,000 | 100 | 1,000 | | |
| 07/08/20 | 5.3 | <0.26 | <0.50 | <0.84 | <0.25 | <0.32 | <0.73 | <3.0 | <0.27 | | |
| 07/08/20 | 5.2 | <0.26 | <0.50 | <0.84 | <0.25 | <0.32 | <0.73 | <3.0 | <0.27 | | |
| 08/18/21 | 6.0 | <0.43 | <0.68 | <0.45 | <0.46 | <0.34 | <0.81 | <0.33 | <0.45 | | |
| MW-7 | | | | | | | | | | | |
| 04/12/16 | <0.47 | <0.33 | <0.20 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | | |
| 11/11/16 | <0.50 | <0.33 | <0.22 | <0.50 | <0.50 | <0.50 | <1.50 | <0.50 | <0.50 | | |
| 09/18/19 | 0.59 J | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | | |
| 12/13/19 | 0.70 J | <0.26 | <0.50 | <0.84 | <0.25 | <0.22 | <0.73 | <0.47 | <0.17 | | |
| 07/08/20 | 0.80 J | <0.26 | <0.50 | <0.84 | <0.25 | <0.32 | <0.73 | <3.0 | <0.27 | | |
| 08/18/21 | 0.66 J | <0.43 | <0.68 | <0.45 | <0.46 | <0.34 | <0.81 | <0.33 | <0.45 | CM=0.86 $\mu\text{g/l}$ | |

TABLE 2
SUMMARY OF GROUNDWATER MONITORING ANALYTICAL RESULTS

NOTES:

Concentrations are in micrograms per liter ($\mu\text{g}/\ell$).

Detected concentrations at or above an NR 140 PAL are italicized and those at or above an NR 140 ES are bold.

Duplicate (Dup) results are averaged for statistical analysis/plotting, per December 2013 ITRC guidance.

All samples except those collected on 09/06/06 were analyzed for a full suite of VOCs using EPA Method 8021 or EPA 8260B. Only compounds detected in one or more samples are shown on this table.

J = Estimated concentration below laboratory quantitation level.

NA = Not analyzed.

NR 140 ES = Wisconsin Administrative Code NR 140 Enforcement Standard.

NR 140 PAL = Wisconsin Administrative Code NR 140 Preventive Action Limit.

PCE = Tetrachloroethylene.

TCE = Trichloroethylene.

U = Compound not detected at or above the average of the limits of detection measured in the sample and its duplicate.

FOOTNOTES:

(1) The 01/28/09 trip blank contained 0.84 J $\mu\text{g}/\ell$ of chloromethane as did MW-1 (0.47 J $\mu\text{g}/\ell$) and MW-2 (1.27 $\mu\text{g}/\ell$).

(2) Methylene chloride was detected at concentrations above its method detection limit but below its quantitation limit in the samples collected on 08/01/18 from MW-2 (at 1.0 $\mu\text{g}/\ell$), MW-4 (0.75 $\mu\text{g}/\ell$), and the two samples collected from MW-6 (1.4 and 0.72 $\mu\text{g}/\ell$).

(3) Chloromethane detected above its method detection limit but below its quantitation limit in the duplicate sample for MW-7 (2.5 $\mu\text{g}/\ell$).

KOELLER ONE, LLC
KOELLER SHOPPING CENTER
OSHKOSH, WISCONSIN

TABLE 3

SUMMARY OF HISTORICAL RNA DATA FIELD MEASUREMENTS

| Well ID | Temperature (°C) | Conductivity (mS/cm) | Dissolved Oxygen (mg/l) | pH | ORP (mV) | Comments/ Footnotes |
|-------------|---------------------|-------------------------|----------------------------|-----|-------------|------------------------|
| MW-1 | | | | | | |
| 06/15/13 | 14.5 | 0.375 | 2.70 | 7.3 | -36.2 | |
| 09/12/13 | 15.4 | 0.623 | 2.07 | 7.6 | -77.1 | |
| 07/06/17 | 14.8 | 0.513 | 3.26 | 7.7 | 57.2 | |
| 07/06/17 | 15.4 | 0.742 | 3.82 | 7.9 | 27.3 | (1) |
| 01/08/18 | 14.4 | 0.443 | 0.56 | 7.4 | 68.2 | |
| 01/08/18 | 14.0 | 0.721 | 0.73 | 7.8 | 54.7 | (1) |
| 08/01/18 | 14.6 | 0.730 | 1.08 | 7.3 | 171.8 | (2) |
| 03/12/19 | 12.0 | 0.698 | 0.80 | 7.6 | 151.1 | (2) |
| 03/12/19 | 11.5 | 0.729 | 1.21 | 7.6 | 87.6 | (3) |
| 09/18/19 | 16.5 | 0.787 | 2.70 | 7.7 | 165.1 | (2) |
| 12/13/19 | 15.3 | 0.950 | 1.95 | 7.7 | 218.7 | (2) |
| 08/18/21 | 16.7 | 0.746 | 1.49 | 7.5 | 138.8 | |
| MW-2 | | | | | | |
| 06/15/13 | 13.7 | 3.465 | 1.87 | 7.0 | -22.9 | |
| 09/12/13 | 15.6 | 2.964 | 3.49 | 6.9 | -56.5 | |
| 07/06/17 | 13.9 | 5.395 | 0.42 | 6.9 | 7.6 | |
| 07/06/17 | 13.5 | 5.141 | 0.58 | 6.8 | 31.3 | (1) |
| 01/08/18 | 14.4 | 4.837 | 0.40 | 6.3 | 137.5 | |
| 01/08/18 | 13.9 | 5.227 | 0.39 | 6.7 | 52.6 | (1) |
| 08/01/18 | 14.4 | 5.385 | 1.52 | 6.1 | 202.0 | (2) |
| 03/12/19 | 10.6 | 4.202 | 1.89 | 6.6 | 189.0 | (2) |
| 03/12/19 | 10.6 | 4.434 | 0.55 | 6.7 | 59.8 | (3) |
| 09/18/19 | 16.4 | 5.941 | 1.59 | 6.8 | 227.3 | (3) |
| 12/13/19 | 14.8 | 6.704 | 1.44 | 6.8 | 220.2 | (2) |
| 08/18/21 | 18.5 | 6.629 | 0.07 | 6.7 | 110.0 | |
| MW-3 | | | | | | |
| 06/15/13 | 14.0 | 2.013 | 1.35 | 6.9 | 24.0 | |
| 09/12/13 | 15.1 | 1.427 | 3.12 | 7.0 | -22.7 | |
| 07/06/17 | 14.4 | 3.871 | 1.54 | 6.9 | 39.3 | |
| 07/06/17 | 14.2 | 3.456 | 2.65 | 7.1 | 89.5 | (1) |
| 01/08/18 | 10.3 | 2.732 | 0.65 | 6.7 | 92.7 | |
| 01/08/18 | 13.4 | 2.323 | 1.87 | 7.2 | 65.8 | (1) |
| 08/01/18 | 14.9 | 0.754 | 0.55 | 7.0 | 144.1 | (2) |
| 03/12/19 | 10.6 | 2.623 | 0.17 | 7.2 | -82.6 | (2) |
| 09/18/19 | 16.8 | 0.915 | 2.34 | 7.3 | -174.3 | (2) |
| 12/13/19 | 14.6 | 4.229 | 1.85 | 7.2 | -28.2 | (2) |
| 08/18/21 | 16.8 | 1.173 | 0.05 | 7.3 | 109.0 | |
| MW-4 | | | | | | |
| 06/15/13 | 15.0 | 0.607 | 2.64 | 8.0 | 27.2 | |
| 09/12/13 | 16.0 | 0.547 | 2.43 | 7.8 | -6.3 | |
| 07/06/17 | 15.9 | 0.676 | 2.65 | 7.8 | 11.6 | |
| 07/06/17 | 14.5 | 0.551 | 0.68 | 7.8 | 55.0 | (1) |

TABLE 3

SUMMARY OF HISTORICAL RNA DATA FIELD MEASUREMENTS

| Well ID | Temperature (°C) | Conductivity (mS/cm) | Dissolved Oxygen (mg/l) | pH | ORP (mV) | Comments/ Footnotes | |
|------------------|---------------------|-------------------------|----------------------------|------|-------------|------------------------|-----|
| Measurement Date | | | | | | | |
| | 01/08/18 | 13.0 | 0.494 | 1.82 | 7.7 | 77.5 | (1) |
| | 08/01/18 | 15.7 | 0.456 | 1.18 | 7.4 | 171.9 | (2) |
| | 03/12/19 | 12.5 | 0.426 | 2.60 | 7.7 | 170.0 | (2) |
| | 03/12/19 | 12.0 | 0.412 | 4.30 | 7.7 | 136.3 | (3) |
| | 09/18/19 | 17.4 | 0.582 | 3.40 | 7.8 | 190.5 | (2) |
| | 12/13/19 | 15.6 | 0.558 | 2.59 | 7.9 | 212.9 | (2) |
| | 08/18/21 | 18.0 | 0.900 | 0.79 | 7.6 | 25.4 | |
| MW-5 | | | | | | | |
| | 07/06/17 | 16.1 | 1.660 | 4.9 | 7.3 | 34.3 | |
| | 09/18/19 | 16.6 | 2.455 | 1.62 | 7.1 | 175.3 | (2) |
| | 12/13/19 | 15.8 | 2.373 | 1.43 | 7.1 | 236.3 | (2) |
| | 08/18/21 | 18.1 | 1.804 | 1.84 | 6.6 | 31.2 | |
| MW-6 | | | | | | | |
| | 07/06/17 | 14.4 | 3.095 | 0.33 | 7.1 | 14.8 | |
| | 07/06/17 | 14.2 | 2.850 | 0.34 | 7.0 | 43.4 | (1) |
| | 01/08/18 | 13.6 | 1.343 | 0.23 | 6.8 | 135.0 | |
| | 01/08/18 | 13.8 | 1.385 | 0.74 | 7.3 | 61.1 | (1) |
| | 08/01/18 | 13.4 | 1.263 | 0.68 | 6.9 | 162.8 | (2) |
| | 03/12/19 | 11.3 | 1.137 | 1.20 | 7.2 | 177.5 | (2) |
| | 03/12/19 | 10.9 | 1.154 | 2.69 | 7.2 | 106.0 | (3) |
| | 09/18/19 | 15.3 | 1.203 | 1.72 | 7.4 | 209.2 | (3) |
| | 12/13/19 | 14.7 | 1.571 | 1.47 | 7.4 | 205.3 | (2) |
| | 08/18/21 | 15.4 | 2.922 | 0.24 | 6.7 | 74.5 | |
| MW-7 | | | | | | | |
| | 07/06/17 | 15.5 | 6.310 | 2.52 | 6.9 | 77.7 | |
| | 08/01/18 | 16.8 | 7.926 | 2.23 | 6.9 | 184.1 | (2) |
| | 09/18/19 | 17.7 | 7.278 | 8.41 | 7.0 | 242.0 | (2) |
| | 12/13/19 | 14.3 | 7.201 | 2.83 | 7.0 | 212.7 | (2) |
| | 08/18/21 | 18.6 | 6.100 | 1.65 | 6.9 | 23.7 | |

NOTES:

RNA = Remediation through natural attenuation.

Conductivity in millSiemens per centimeter (mS/cm).

Dissolved oxygen (DO) concentration in milligrams per liter (mg/l).

Oxidation reduction potential (ORP) in millivolts (mV).

Temperature in degrees Celsius (°C).

Water quality parameter data collected using a YSI 556 multi-parameter meter prior to 09/18/19, and a Hanna HI98194 on 09/18/19 and thereafter.

FOOTNOTES:

(1) RNA parameters collected in-situ after purging and sampling.

(2) RNA parameters collected after purging and prior to sampling.

(3) RNA parameters collected after purging and sampling.

KOELLER ONE LLC
KOELLER SHOPPING CENTER
OSHKOSH, WISCONSIN

TABLE 4

SUMMARY OF DETECTED PFAS COMPOUNDS IN MW-1
AUGUST 2021

| Compound | CAS Number | MW-1 | WI DHS Proposed PFAS Standard | |
|-------------------------------------|------------|------------|-------------------------------|------------------------|
| | | | ES (ng/l) | PAL (ng/l) |
| Perfluorooctanoic acid (PFOA) | 335-67-1 | <u>5.8</u> | <u>20⁽¹⁾</u> | <u>2⁽¹⁾</u> |
| Perfluorooctanesulfonic acid (PFOS) | 1763-23-1 | <u>9.7</u> | | |
| Perfluorononanoic acid (PFNA) | 375-95-1 | 0.95 J | 30 | <u>3</u> |
| Perfluorobutanoic acid (PFBA) | 375-22-4 | 5.1 | 10,000 | <u>2,000</u> |
| Perfluorohexanoic acid (PFHxA) | 307-24-4 | 2.3 J | 150,000 | <u>30,000</u> |
| Perfluorobutanesulfonic acid (PFBS) | 375-73-5 | 3.2 J | 450,000 | <u>90,000</u> |
| Perfluoropentanoic acid (PFPeA) | 2706-90-3 | 2.6 J | NRS | NRS |
| Perfluoroheptanoic acid (PFHpA) | 375-85-9 | 2.3 J | NRS | NRS |

NOTES:

The groundwater sample from MW-1 was collected on 8/18/21 and analyzed for 33 PFAS compounds using Modified 537 Method. Only compounds detected in the sample are shown in this table.

All concentrations are in nanograms per liter (ng/L).

NRS - No recommended standard

WI DHS Recommended preventive action limits (PALs) and enforcement standards (ES) taken from WDNR's March 1, 2021, PFAS update.

J - Estimated concentration between laboratory's quantitation and method detection levels.

FOOTNOTE:

(1) The proposed NR 140 PALs and ESs are for the combined concentrations of PFOA, PFOS, FOSA, NEtFOSA, NEtFOSAA, and NEtFOSE. FOSA, NEtFOSA, NEtFOSAA, and NEtFOSE were not detected in this sample and are therefore not included in this table.

ATTACHMENT A

ANALYTICAL RESULTS AND CHAIN OF CUSTODY RECORDS FOR GROUNDWATER
SAMPLES COLLECTED IN JULY 2020 AND AUGUST 2021

July 13, 2020

Tony Miller
Gannett Fleming
8040 Excelsior Drive, Ste 303
Madison, WI 53717

RE: Project: 47358.003 KOELLER ONE
Pace Project No.: 40210819

Dear Tony Miller:

Enclosed are the analytical results for sample(s) received by the laboratory on July 09, 2020. The results relate only to the samples included in this report. Results reported herein conform to the applicable TNI/NELAC Standards and the laboratory's Quality Manual, where applicable, unless otherwise noted in the body of the report.

The test results provided in this final report were generated by each of the following laboratories within the Pace Network:

- Pace Analytical Services - Green Bay

If you have any questions concerning this report, please feel free to contact me.

Sincerely,



Dan Milewsky
dan.milewsky@pacelabs.com
(920)469-2436
Project Manager

Enclosures

cc: Chelsea Payne, Gannett Fleming Inc.



REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

CERTIFICATIONS

Project: 47358.003 KOELLER ONE
Pace Project No.: 40210819

Pace Analytical Services Green Bay

1241 Bellevue Street, Green Bay, WI 54302
Florida/NELAP Certification #: E87948
Illinois Certification #: 200050
Kentucky UST Certification #: 82
Louisiana Certification #: 04168
Minnesota Certification #: 055-999-334
New York Certification #: 12064
North Dakota Certification #: R-150

Virginia VELAP ID: 460263
South Carolina Certification #: 83006001
Texas Certification #: T104704529-14-1
Wisconsin Certification #: 405132750
Wisconsin DATCP Certification #: 105-444
USDA Soil Permit #: P330-16-00157
Federal Fish & Wildlife Permit #: LE51774A-0

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

SAMPLE SUMMARY

Project: 47358.003 KOELLER ONE
Pace Project No.: 40210819

| Lab ID | Sample ID | Matrix | Date Collected | Date Received |
|-------------|------------|--------|----------------|----------------|
| 40210819001 | MW-1 | Water | 07/08/20 12:30 | 07/09/20 09:40 |
| 40210819002 | MW-2 | Water | 07/08/20 11:30 | 07/09/20 09:40 |
| 40210819003 | MW-3 | Water | 07/08/20 11:35 | 07/09/20 09:40 |
| 40210819004 | MW-4 | Water | 07/08/20 11:15 | 07/09/20 09:40 |
| 40210819005 | MW-5 | Water | 07/08/20 10:35 | 07/09/20 09:40 |
| 40210819006 | MW-6 | Water | 07/08/20 09:50 | 07/09/20 09:40 |
| 40210819007 | MW-6 DUP | Water | 07/08/20 09:50 | 07/09/20 09:40 |
| 40210819008 | MW-7 | Water | 07/08/20 09:10 | 07/09/20 09:40 |
| 40210819009 | TRIP BLANK | Water | 07/08/20 00:00 | 07/09/20 09:40 |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

SAMPLE ANALYTE COUNT

Project: 47358.003 KOELLER ONE
Pace Project No.: 40210819

| Lab ID | Sample ID | Method | Analysts | Analytes Reported |
|-------------|------------|----------|----------|-------------------|
| 40210819001 | MW-1 | EPA 8260 | LAP | 63 |
| 40210819002 | MW-2 | EPA 8260 | LAP | 63 |
| 40210819003 | MW-3 | EPA 8260 | LAP | 63 |
| 40210819004 | MW-4 | EPA 8260 | LAP | 63 |
| 40210819005 | MW-5 | EPA 8260 | LAP | 63 |
| 40210819006 | MW-6 | EPA 8260 | LAP | 63 |
| 40210819007 | MW-6 DUP | EPA 8260 | LAP | 63 |
| 40210819008 | MW-7 | EPA 8260 | LAP | 63 |
| 40210819009 | TRIP BLANK | EPA 8260 | LAP | 63 |

PASI-G = Pace Analytical Services - Green Bay

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

SUMMARY OF DETECTION

Project: 47358.003 KOELLER ONE
Pace Project No.: 40210819

| Lab Sample ID | Client Sample ID | | | | | |
|--------------------|-------------------|--------|-------|--------------|----------------|------------|
| Method | Parameters | Result | Units | Report Limit | Analyzed | Qualifiers |
| 40210819001 | MW-1 | | | | | |
| EPA 8260 | Tetrachloroethene | 181 | ug/L | 2.7 | 07/10/20 16:16 | |
| 40210819002 | MW-2 | | | | | |
| EPA 8260 | Tetrachloroethene | 5.6 | ug/L | 1.1 | 07/10/20 11:08 | |
| 40210819003 | MW-3 | | | | | |
| EPA 8260 | Tetrachloroethene | 63.0 | ug/L | 1.1 | 07/10/20 11:31 | |
| EPA 8260 | Trichloroethene | 0.30J | ug/L | 1.0 | 07/10/20 11:31 | |
| 40210819004 | MW-4 | | | | | |
| EPA 8260 | Tetrachloroethene | 1.3 | ug/L | 1.1 | 07/10/20 10:02 | |
| 40210819006 | MW-6 | | | | | |
| EPA 8260 | Tetrachloroethene | 5.3 | ug/L | 1.1 | 07/10/20 11:53 | |
| 40210819007 | MW-6 DUP | | | | | |
| EPA 8260 | Tetrachloroethene | 5.2 | ug/L | 1.1 | 07/10/20 12:14 | |
| 40210819008 | MW-7 | | | | | |
| EPA 8260 | Tetrachloroethene | 0.80J | ug/L | 1.1 | 07/10/20 10:46 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

| Sample: MW-1 | Lab ID: 40210819001 | Collected: 07/08/20 12:30 | Received: 07/09/20 09:40 | Matrix: Water | | | | | |
|-----------------------------|--------------------------------------|---------------------------|--------------------------|---------------|-----|----------|----------------|-----------|------|
| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | <0.67 | ug/L | 2.5 | 0.67 | 2.5 | | 07/10/20 16:16 | 630-20-6 | |
| 1,1,1-Trichloroethane | <0.61 | ug/L | 2.5 | 0.61 | 2.5 | | 07/10/20 16:16 | 71-55-6 | |
| 1,1,2,2-Tetrachloroethane | <0.69 | ug/L | 2.5 | 0.69 | 2.5 | | 07/10/20 16:16 | 79-34-5 | |
| 1,1,2-Trichloroethane | <1.4 | ug/L | 12.5 | 1.4 | 2.5 | | 07/10/20 16:16 | 79-00-5 | |
| 1,1-Dichloroethane | <0.68 | ug/L | 2.5 | 0.68 | 2.5 | | 07/10/20 16:16 | 75-34-3 | |
| 1,1-Dichloroethene | <0.61 | ug/L | 2.5 | 0.61 | 2.5 | | 07/10/20 16:16 | 75-35-4 | |
| 1,1-Dichloropropene | <1.4 | ug/L | 4.5 | 1.4 | 2.5 | | 07/10/20 16:16 | 563-58-6 | |
| 1,2,3-Trichlorobenzene | <5.5 | ug/L | 18.4 | 5.5 | 2.5 | | 07/10/20 16:16 | 87-61-6 | |
| 1,2,3-Trichloropropane | <1.5 | ug/L | 12.5 | 1.5 | 2.5 | | 07/10/20 16:16 | 96-18-4 | |
| 1,2,4-Trichlorobenzene | <2.4 | ug/L | 12.5 | 2.4 | 2.5 | | 07/10/20 16:16 | 120-82-1 | |
| 1,2,4-Trimethylbenzene | <2.1 | ug/L | 7.0 | 2.1 | 2.5 | | 07/10/20 16:16 | 95-63-6 | |
| 1,2-Dibromo-3-chloropropane | <4.4 | ug/L | 14.7 | 4.4 | 2.5 | | 07/10/20 16:16 | 96-12-8 | |
| 1,2-Dibromoethane (EDB) | <2.1 | ug/L | 6.9 | 2.1 | 2.5 | | 07/10/20 16:16 | 106-93-4 | |
| 1,2-Dichlorobenzene | <1.8 | ug/L | 5.9 | 1.8 | 2.5 | | 07/10/20 16:16 | 95-50-1 | |
| 1,2-Dichloroethane | <0.70 | ug/L | 2.5 | 0.70 | 2.5 | | 07/10/20 16:16 | 107-06-2 | |
| 1,2-Dichloropropane | <0.71 | ug/L | 2.5 | 0.71 | 2.5 | | 07/10/20 16:16 | 78-87-5 | |
| 1,3,5-Trimethylbenzene | <2.2 | ug/L | 7.3 | 2.2 | 2.5 | | 07/10/20 16:16 | 108-67-8 | |
| 1,3-Dichlorobenzene | <1.6 | ug/L | 5.2 | 1.6 | 2.5 | | 07/10/20 16:16 | 541-73-1 | |
| 1,3-Dichloropropene | <2.1 | ug/L | 6.9 | 2.1 | 2.5 | | 07/10/20 16:16 | 142-28-9 | |
| 1,4-Dichlorobenzene | <2.4 | ug/L | 7.9 | 2.4 | 2.5 | | 07/10/20 16:16 | 106-46-7 | |
| 2,2-Dichloropropane | <5.7 | ug/L | 18.9 | 5.7 | 2.5 | | 07/10/20 16:16 | 594-20-7 | |
| 2-Chlorotoluene | <2.3 | ug/L | 12.5 | 2.3 | 2.5 | | 07/10/20 16:16 | 95-49-8 | |
| 4-Chlorotoluene | <1.9 | ug/L | 6.3 | 1.9 | 2.5 | | 07/10/20 16:16 | 106-43-4 | |
| Benzene | <0.62 | ug/L | 2.5 | 0.62 | 2.5 | | 07/10/20 16:16 | 71-43-2 | |
| Bromobenzene | <0.60 | ug/L | 2.5 | 0.60 | 2.5 | | 07/10/20 16:16 | 108-86-1 | |
| Bromochloromethane | <0.91 | ug/L | 12.5 | 0.91 | 2.5 | | 07/10/20 16:16 | 74-97-5 | |
| Bromodichloromethane | <0.91 | ug/L | 3.0 | 0.91 | 2.5 | | 07/10/20 16:16 | 75-27-4 | |
| Bromoform | <9.9 | ug/L | 33.1 | 9.9 | 2.5 | | 07/10/20 16:16 | 75-25-2 | |
| Bromomethane | <2.4 | ug/L | 12.5 | 2.4 | 2.5 | | 07/10/20 16:16 | 74-83-9 | |
| Carbon tetrachloride | <2.7 | ug/L | 9.0 | 2.7 | 2.5 | | 07/10/20 16:16 | 56-23-5 | |
| Chlorobenzene | <1.8 | ug/L | 5.9 | 1.8 | 2.5 | | 07/10/20 16:16 | 108-90-7 | |
| Chloroethane | <3.4 | ug/L | 12.5 | 3.4 | 2.5 | | 07/10/20 16:16 | 75-00-3 | |
| Chloroform | <3.2 | ug/L | 12.5 | 3.2 | 2.5 | | 07/10/20 16:16 | 67-66-3 | |
| Chloromethane | <5.5 | ug/L | 18.2 | 5.5 | 2.5 | | 07/10/20 16:16 | 74-87-3 | |
| Dibromochloromethane | <6.5 | ug/L | 21.7 | 6.5 | 2.5 | | 07/10/20 16:16 | 124-48-1 | |
| Dibromomethane | <2.3 | ug/L | 7.8 | 2.3 | 2.5 | | 07/10/20 16:16 | 74-95-3 | |
| Dichlorodifluoromethane | <1.2 | ug/L | 12.5 | 1.2 | 2.5 | | 07/10/20 16:16 | 75-71-8 | |
| Ethylbenzene | <0.80 | ug/L | 2.7 | 0.80 | 2.5 | | 07/10/20 16:16 | 100-41-4 | |
| Hexachloro-1,3-butadiene | <3.7 | ug/L | 12.2 | 3.7 | 2.5 | | 07/10/20 16:16 | 87-68-3 | |
| Isopropylbenzene (Cumene) | <4.2 | ug/L | 14.0 | 4.2 | 2.5 | | 07/10/20 16:16 | 98-82-8 | |
| Methyl-tert-butyl ether | <3.1 | ug/L | 10.4 | 3.1 | 2.5 | | 07/10/20 16:16 | 1634-04-4 | |
| Methylene Chloride | <1.5 | ug/L | 12.5 | 1.5 | 2.5 | | 07/10/20 16:16 | 75-09-2 | |
| Naphthalene | <2.9 | ug/L | 12.5 | 2.9 | 2.5 | | 07/10/20 16:16 | 91-20-3 | |
| Styrene | <7.5 | ug/L | 25.1 | 7.5 | 2.5 | | 07/10/20 16:16 | 100-42-5 | |
| Tetrachloroethene | 181 | ug/L | 2.7 | 0.82 | 2.5 | | 07/10/20 16:16 | 127-18-4 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE
Pace Project No.: 40210819

Sample: MW-1 Lab ID: 40210819001 Collected: 07/08/20 12:30 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|---------------------------|--------------------------------------|-------|--------|------|-----|----------|----------------|-------------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| Toluene | <0.67 | ug/L | 2.2 | 0.67 | 2.5 | | 07/10/20 16:16 | 108-88-3 | |
| Trichloroethene | <0.64 | ug/L | 2.5 | 0.64 | 2.5 | | 07/10/20 16:16 | 79-01-6 | |
| Trichlorofluoromethane | <0.54 | ug/L | 2.5 | 0.54 | 2.5 | | 07/10/20 16:16 | 75-69-4 | |
| Vinyl chloride | <0.44 | ug/L | 2.5 | 0.44 | 2.5 | | 07/10/20 16:16 | 75-01-4 | |
| cis-1,2-Dichloroethene | <0.68 | ug/L | 2.5 | 0.68 | 2.5 | | 07/10/20 16:16 | 156-59-2 | |
| cis-1,3-Dichloropropene | <9.1 | ug/L | 30.2 | 9.1 | 2.5 | | 07/10/20 16:16 | 10061-01-5 | |
| m&p-Xylene | <1.2 | ug/L | 5.0 | 1.2 | 2.5 | | 07/10/20 16:16 | 179601-23-1 | |
| n-Butylbenzene | <1.8 | ug/L | 5.9 | 1.8 | 2.5 | | 07/10/20 16:16 | 104-51-8 | |
| n-Propylbenzene | <2.0 | ug/L | 12.5 | 2.0 | 2.5 | | 07/10/20 16:16 | 103-65-1 | |
| o-Xylene | <0.65 | ug/L | 2.5 | 0.65 | 2.5 | | 07/10/20 16:16 | 95-47-6 | |
| p-Isopropyltoluene | <2.0 | ug/L | 6.7 | 2.0 | 2.5 | | 07/10/20 16:16 | 99-87-6 | |
| sec-Butylbenzene | <2.1 | ug/L | 12.5 | 2.1 | 2.5 | | 07/10/20 16:16 | 135-98-8 | |
| tert-Butylbenzene | <0.76 | ug/L | 2.5 | 0.76 | 2.5 | | 07/10/20 16:16 | 98-06-6 | |
| trans-1,2-Dichloroethene | <1.2 | ug/L | 3.9 | 1.2 | 2.5 | | 07/10/20 16:16 | 156-60-5 | |
| trans-1,3-Dichloropropene | <10.9 | ug/L | 36.4 | 10.9 | 2.5 | | 07/10/20 16:16 | 10061-02-6 | |
| Surrogates | | | | | | | | | |
| 4-Bromofluorobenzene (S) | 90 | % | 70-130 | | 2.5 | | 07/10/20 16:16 | 460-00-4 | |
| Dibromofluoromethane (S) | 95 | % | 70-130 | | 2.5 | | 07/10/20 16:16 | 1868-53-7 | |
| Toluene-d8 (S) | 98 | % | 70-130 | | 2.5 | | 07/10/20 16:16 | 2037-26-5 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: MW-2 **Lab ID: 40210819002** Collected: 07/08/20 11:30 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|-----------------------------|--------------------------------------|-------|------|------|----|----------|----------------|-----------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 11:08 | 630-20-6 | |
| 1,1,1-Trichloroethane | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 11:08 | 71-55-6 | |
| 1,1,2,2-Tetrachloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 11:08 | 79-34-5 | |
| 1,1,2-Trichloroethane | <0.55 | ug/L | 5.0 | 0.55 | 1 | | 07/10/20 11:08 | 79-00-5 | |
| 1,1-Dichloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 11:08 | 75-34-3 | |
| 1,1-Dichloroethene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 11:08 | 75-35-4 | |
| 1,1-Dichloropropene | <0.54 | ug/L | 1.8 | 0.54 | 1 | | 07/10/20 11:08 | 563-58-6 | |
| 1,2,3-Trichlorobenzene | <2.2 | ug/L | 7.4 | 2.2 | 1 | | 07/10/20 11:08 | 87-61-6 | |
| 1,2,3-Trichloropropane | <0.59 | ug/L | 5.0 | 0.59 | 1 | | 07/10/20 11:08 | 96-18-4 | |
| 1,2,4-Trichlorobenzene | <0.95 | ug/L | 5.0 | 0.95 | 1 | | 07/10/20 11:08 | 120-82-1 | |
| 1,2,4-Trimethylbenzene | <0.84 | ug/L | 2.8 | 0.84 | 1 | | 07/10/20 11:08 | 95-63-6 | |
| 1,2-Dibromo-3-chloropropane | <1.8 | ug/L | 5.9 | 1.8 | 1 | | 07/10/20 11:08 | 96-12-8 | |
| 1,2-Dibromoethane (EDB) | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 11:08 | 106-93-4 | |
| 1,2-Dichlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 11:08 | 95-50-1 | |
| 1,2-Dichloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 11:08 | 107-06-2 | |
| 1,2-Dichloropropane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 11:08 | 78-87-5 | |
| 1,3,5-Trimethylbenzene | <0.87 | ug/L | 2.9 | 0.87 | 1 | | 07/10/20 11:08 | 108-67-8 | |
| 1,3-Dichlorobenzene | <0.63 | ug/L | 2.1 | 0.63 | 1 | | 07/10/20 11:08 | 541-73-1 | |
| 1,3-Dichloropropane | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 11:08 | 142-28-9 | |
| 1,4-Dichlorobenzene | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 11:08 | 106-46-7 | |
| 2,2-Dichloropropane | <2.3 | ug/L | 7.6 | 2.3 | 1 | | 07/10/20 11:08 | 594-20-7 | |
| 2-Chlorotoluene | <0.93 | ug/L | 5.0 | 0.93 | 1 | | 07/10/20 11:08 | 95-49-8 | |
| 4-Chlorotoluene | <0.76 | ug/L | 2.5 | 0.76 | 1 | | 07/10/20 11:08 | 106-43-4 | |
| Benzene | <0.25 | ug/L | 1.0 | 0.25 | 1 | | 07/10/20 11:08 | 71-43-2 | |
| Bromobenzene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 11:08 | 108-86-1 | |
| Bromochloromethane | <0.36 | ug/L | 5.0 | 0.36 | 1 | | 07/10/20 11:08 | 74-97-5 | |
| Bromodichloromethane | <0.36 | ug/L | 1.2 | 0.36 | 1 | | 07/10/20 11:08 | 75-27-4 | |
| Bromoform | <4.0 | ug/L | 13.2 | 4.0 | 1 | | 07/10/20 11:08 | 75-25-2 | |
| Bromomethane | <0.97 | ug/L | 5.0 | 0.97 | 1 | | 07/10/20 11:08 | 74-83-9 | |
| Carbon tetrachloride | <1.1 | ug/L | 3.6 | 1.1 | 1 | | 07/10/20 11:08 | 56-23-5 | |
| Chlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 11:08 | 108-90-7 | |
| Chloroethane | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 11:08 | 75-00-3 | |
| Chloroform | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 11:08 | 67-66-3 | |
| Chloromethane | <2.2 | ug/L | 7.3 | 2.2 | 1 | | 07/10/20 11:08 | 74-87-3 | |
| Dibromochloromethane | <2.6 | ug/L | 8.7 | 2.6 | 1 | | 07/10/20 11:08 | 124-48-1 | |
| Dibromomethane | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 11:08 | 74-95-3 | |
| Dichlorodifluoromethane | <0.50 | ug/L | 5.0 | 0.50 | 1 | | 07/10/20 11:08 | 75-71-8 | |
| Ethylbenzene | <0.32 | ug/L | 1.1 | 0.32 | 1 | | 07/10/20 11:08 | 100-41-4 | |
| Hexachloro-1,3-butadiene | <1.5 | ug/L | 4.9 | 1.5 | 1 | | 07/10/20 11:08 | 87-68-3 | |
| Isopropylbenzene (Cumene) | <1.7 | ug/L | 5.6 | 1.7 | 1 | | 07/10/20 11:08 | 98-82-8 | |
| Methyl-tert-butyl ether | <1.2 | ug/L | 4.2 | 1.2 | 1 | | 07/10/20 11:08 | 1634-04-4 | |
| Methylene Chloride | <0.58 | ug/L | 5.0 | 0.58 | 1 | | 07/10/20 11:08 | 75-09-2 | |
| Naphthalene | <1.2 | ug/L | 5.0 | 1.2 | 1 | | 07/10/20 11:08 | 91-20-3 | |
| Styrene | <3.0 | ug/L | 10.0 | 3.0 | 1 | | 07/10/20 11:08 | 100-42-5 | |
| Tetrachloroethene | 5.6 | ug/L | 1.1 | 0.33 | 1 | | 07/10/20 11:08 | 127-18-4 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: MW-2 **Lab ID: 40210819002** Collected: 07/08/20 11:30 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|---------------------------|--------------------------------------|-------|--------|------|----|----------|----------------|-------------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| Toluene | <0.27 | ug/L | 0.90 | 0.27 | 1 | | 07/10/20 11:08 | 108-88-3 | |
| Trichloroethene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 11:08 | 79-01-6 | |
| Trichlorofluoromethane | <0.21 | ug/L | 1.0 | 0.21 | 1 | | 07/10/20 11:08 | 75-69-4 | |
| Vinyl chloride | <0.17 | ug/L | 1.0 | 0.17 | 1 | | 07/10/20 11:08 | 75-01-4 | |
| cis-1,2-Dichloroethene | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 11:08 | 156-59-2 | |
| cis-1,3-Dichloropropene | <3.6 | ug/L | 12.1 | 3.6 | 1 | | 07/10/20 11:08 | 10061-01-5 | |
| m&p-Xylene | <0.47 | ug/L | 2.0 | 0.47 | 1 | | 07/10/20 11:08 | 179601-23-1 | |
| n-Butylbenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 11:08 | 104-51-8 | |
| n-Propylbenzene | <0.81 | ug/L | 5.0 | 0.81 | 1 | | 07/10/20 11:08 | 103-65-1 | |
| o-Xylene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 11:08 | 95-47-6 | |
| p-Isopropyltoluene | <0.80 | ug/L | 2.7 | 0.80 | 1 | | 07/10/20 11:08 | 99-87-6 | |
| sec-Butylbenzene | <0.85 | ug/L | 5.0 | 0.85 | 1 | | 07/10/20 11:08 | 135-98-8 | |
| tert-Butylbenzene | <0.30 | ug/L | 1.0 | 0.30 | 1 | | 07/10/20 11:08 | 98-06-6 | |
| trans-1,2-Dichloroethene | <0.46 | ug/L | 1.5 | 0.46 | 1 | | 07/10/20 11:08 | 156-60-5 | |
| trans-1,3-Dichloropropene | <4.4 | ug/L | 14.6 | 4.4 | 1 | | 07/10/20 11:08 | 10061-02-6 | |
| Surrogates | | | | | | | | | |
| 4-Bromofluorobenzene (S) | 90 | % | 70-130 | | 1 | | 07/10/20 11:08 | 460-00-4 | |
| Dibromofluoromethane (S) | 84 | % | 70-130 | | 1 | | 07/10/20 11:08 | 1868-53-7 | |
| Toluene-d8 (S) | 97 | % | 70-130 | | 1 | | 07/10/20 11:08 | 2037-26-5 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: MW-3 **Lab ID: 40210819003** Collected: 07/08/20 11:35 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|-----------------------------|--------------------------------------|-------|------|------|----|----------|----------------|-----------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 11:31 | 630-20-6 | |
| 1,1,1-Trichloroethane | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 11:31 | 71-55-6 | |
| 1,1,2,2-Tetrachloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 11:31 | 79-34-5 | |
| 1,1,2-Trichloroethane | <0.55 | ug/L | 5.0 | 0.55 | 1 | | 07/10/20 11:31 | 79-00-5 | |
| 1,1-Dichloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 11:31 | 75-34-3 | |
| 1,1-Dichloroethene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 11:31 | 75-35-4 | |
| 1,1-Dichloropropene | <0.54 | ug/L | 1.8 | 0.54 | 1 | | 07/10/20 11:31 | 563-58-6 | |
| 1,2,3-Trichlorobenzene | <2.2 | ug/L | 7.4 | 2.2 | 1 | | 07/10/20 11:31 | 87-61-6 | |
| 1,2,3-Trichloropropane | <0.59 | ug/L | 5.0 | 0.59 | 1 | | 07/10/20 11:31 | 96-18-4 | |
| 1,2,4-Trichlorobenzene | <0.95 | ug/L | 5.0 | 0.95 | 1 | | 07/10/20 11:31 | 120-82-1 | |
| 1,2,4-Trimethylbenzene | <0.84 | ug/L | 2.8 | 0.84 | 1 | | 07/10/20 11:31 | 95-63-6 | |
| 1,2-Dibromo-3-chloropropane | <1.8 | ug/L | 5.9 | 1.8 | 1 | | 07/10/20 11:31 | 96-12-8 | |
| 1,2-Dibromoethane (EDB) | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 11:31 | 106-93-4 | |
| 1,2-Dichlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 11:31 | 95-50-1 | |
| 1,2-Dichloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 11:31 | 107-06-2 | |
| 1,2-Dichloropropane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 11:31 | 78-87-5 | |
| 1,3,5-Trimethylbenzene | <0.87 | ug/L | 2.9 | 0.87 | 1 | | 07/10/20 11:31 | 108-67-8 | |
| 1,3-Dichlorobenzene | <0.63 | ug/L | 2.1 | 0.63 | 1 | | 07/10/20 11:31 | 541-73-1 | |
| 1,3-Dichloropropane | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 11:31 | 142-28-9 | |
| 1,4-Dichlorobenzene | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 11:31 | 106-46-7 | |
| 2,2-Dichloropropane | <2.3 | ug/L | 7.6 | 2.3 | 1 | | 07/10/20 11:31 | 594-20-7 | |
| 2-Chlorotoluene | <0.93 | ug/L | 5.0 | 0.93 | 1 | | 07/10/20 11:31 | 95-49-8 | |
| 4-Chlorotoluene | <0.76 | ug/L | 2.5 | 0.76 | 1 | | 07/10/20 11:31 | 106-43-4 | |
| Benzene | <0.25 | ug/L | 1.0 | 0.25 | 1 | | 07/10/20 11:31 | 71-43-2 | |
| Bromobenzene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 11:31 | 108-86-1 | |
| Bromochloromethane | <0.36 | ug/L | 5.0 | 0.36 | 1 | | 07/10/20 11:31 | 74-97-5 | |
| Bromodichloromethane | <0.36 | ug/L | 1.2 | 0.36 | 1 | | 07/10/20 11:31 | 75-27-4 | |
| Bromoform | <4.0 | ug/L | 13.2 | 4.0 | 1 | | 07/10/20 11:31 | 75-25-2 | |
| Bromomethane | <0.97 | ug/L | 5.0 | 0.97 | 1 | | 07/10/20 11:31 | 74-83-9 | |
| Carbon tetrachloride | <1.1 | ug/L | 3.6 | 1.1 | 1 | | 07/10/20 11:31 | 56-23-5 | |
| Chlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 11:31 | 108-90-7 | |
| Chloroethane | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 11:31 | 75-00-3 | |
| Chloroform | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 11:31 | 67-66-3 | |
| Chloromethane | <2.2 | ug/L | 7.3 | 2.2 | 1 | | 07/10/20 11:31 | 74-87-3 | |
| Dibromochloromethane | <2.6 | ug/L | 8.7 | 2.6 | 1 | | 07/10/20 11:31 | 124-48-1 | |
| Dibromomethane | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 11:31 | 74-95-3 | |
| Dichlorodifluoromethane | <0.50 | ug/L | 5.0 | 0.50 | 1 | | 07/10/20 11:31 | 75-71-8 | |
| Ethylbenzene | <0.32 | ug/L | 1.1 | 0.32 | 1 | | 07/10/20 11:31 | 100-41-4 | |
| Hexachloro-1,3-butadiene | <1.5 | ug/L | 4.9 | 1.5 | 1 | | 07/10/20 11:31 | 87-68-3 | |
| Isopropylbenzene (Cumene) | <1.7 | ug/L | 5.6 | 1.7 | 1 | | 07/10/20 11:31 | 98-82-8 | |
| Methyl-tert-butyl ether | <1.2 | ug/L | 4.2 | 1.2 | 1 | | 07/10/20 11:31 | 1634-04-4 | |
| Methylene Chloride | <0.58 | ug/L | 5.0 | 0.58 | 1 | | 07/10/20 11:31 | 75-09-2 | |
| Naphthalene | <1.2 | ug/L | 5.0 | 1.2 | 1 | | 07/10/20 11:31 | 91-20-3 | |
| Styrene | <3.0 | ug/L | 10.0 | 3.0 | 1 | | 07/10/20 11:31 | 100-42-5 | |
| Tetrachloroethene | 63.0 | ug/L | 1.1 | 0.33 | 1 | | 07/10/20 11:31 | 127-18-4 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: MW-3 Lab ID: 40210819003 Collected: 07/08/20 11:35 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|---------------------------|--------------------------------------|-------|--------|------|----|----------|----------------|-------------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| Toluene | <0.27 | ug/L | 0.90 | 0.27 | 1 | | 07/10/20 11:31 | 108-88-3 | |
| Trichloroethene | 0.30J | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 11:31 | 79-01-6 | |
| Trichlorofluoromethane | <0.21 | ug/L | 1.0 | 0.21 | 1 | | 07/10/20 11:31 | 75-69-4 | |
| Vinyl chloride | <0.17 | ug/L | 1.0 | 0.17 | 1 | | 07/10/20 11:31 | 75-01-4 | |
| cis-1,2-Dichloroethene | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 11:31 | 156-59-2 | |
| cis-1,3-Dichloropropene | <3.6 | ug/L | 12.1 | 3.6 | 1 | | 07/10/20 11:31 | 10061-01-5 | |
| m&p-Xylene | <0.47 | ug/L | 2.0 | 0.47 | 1 | | 07/10/20 11:31 | 179601-23-1 | |
| n-Butylbenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 11:31 | 104-51-8 | |
| n-Propylbenzene | <0.81 | ug/L | 5.0 | 0.81 | 1 | | 07/10/20 11:31 | 103-65-1 | |
| o-Xylene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 11:31 | 95-47-6 | |
| p-Isopropyltoluene | <0.80 | ug/L | 2.7 | 0.80 | 1 | | 07/10/20 11:31 | 99-87-6 | |
| sec-Butylbenzene | <0.85 | ug/L | 5.0 | 0.85 | 1 | | 07/10/20 11:31 | 135-98-8 | |
| tert-Butylbenzene | <0.30 | ug/L | 1.0 | 0.30 | 1 | | 07/10/20 11:31 | 98-06-6 | |
| trans-1,2-Dichloroethene | <0.46 | ug/L | 1.5 | 0.46 | 1 | | 07/10/20 11:31 | 156-60-5 | |
| trans-1,3-Dichloropropene | <4.4 | ug/L | 14.6 | 4.4 | 1 | | 07/10/20 11:31 | 10061-02-6 | |
| Surrogates | | | | | | | | | |
| 4-Bromofluorobenzene (S) | 92 | % | 70-130 | | 1 | | 07/10/20 11:31 | 460-00-4 | |
| Dibromofluoromethane (S) | 98 | % | 70-130 | | 1 | | 07/10/20 11:31 | 1868-53-7 | |
| Toluene-d8 (S) | 100 | % | 70-130 | | 1 | | 07/10/20 11:31 | 2037-26-5 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: MW-4 **Lab ID: 40210819004** Collected: 07/08/20 11:15 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|-----------------------------|--------------------------------------|-------|------|------|----|----------|----------------|-----------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 10:02 | 630-20-6 | |
| 1,1,1-Trichloroethane | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 10:02 | 71-55-6 | |
| 1,1,2,2-Tetrachloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 10:02 | 79-34-5 | |
| 1,1,2-Trichloroethane | <0.55 | ug/L | 5.0 | 0.55 | 1 | | 07/10/20 10:02 | 79-00-5 | |
| 1,1-Dichloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 10:02 | 75-34-3 | |
| 1,1-Dichloroethene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 10:02 | 75-35-4 | |
| 1,1-Dichloropropene | <0.54 | ug/L | 1.8 | 0.54 | 1 | | 07/10/20 10:02 | 563-58-6 | |
| 1,2,3-Trichlorobenzene | <2.2 | ug/L | 7.4 | 2.2 | 1 | | 07/10/20 10:02 | 87-61-6 | |
| 1,2,3-Trichloropropane | <0.59 | ug/L | 5.0 | 0.59 | 1 | | 07/10/20 10:02 | 96-18-4 | |
| 1,2,4-Trichlorobenzene | <0.95 | ug/L | 5.0 | 0.95 | 1 | | 07/10/20 10:02 | 120-82-1 | |
| 1,2,4-Trimethylbenzene | <0.84 | ug/L | 2.8 | 0.84 | 1 | | 07/10/20 10:02 | 95-63-6 | |
| 1,2-Dibromo-3-chloropropane | <1.8 | ug/L | 5.9 | 1.8 | 1 | | 07/10/20 10:02 | 96-12-8 | |
| 1,2-Dibromoethane (EDB) | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 10:02 | 106-93-4 | |
| 1,2-Dichlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 10:02 | 95-50-1 | |
| 1,2-Dichloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 10:02 | 107-06-2 | |
| 1,2-Dichloropropane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 10:02 | 78-87-5 | |
| 1,3,5-Trimethylbenzene | <0.87 | ug/L | 2.9 | 0.87 | 1 | | 07/10/20 10:02 | 108-67-8 | |
| 1,3-Dichlorobenzene | <0.63 | ug/L | 2.1 | 0.63 | 1 | | 07/10/20 10:02 | 541-73-1 | |
| 1,3-Dichloropropane | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 10:02 | 142-28-9 | |
| 1,4-Dichlorobenzene | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 10:02 | 106-46-7 | |
| 2,2-Dichloropropane | <2.3 | ug/L | 7.6 | 2.3 | 1 | | 07/10/20 10:02 | 594-20-7 | |
| 2-Chlorotoluene | <0.93 | ug/L | 5.0 | 0.93 | 1 | | 07/10/20 10:02 | 95-49-8 | |
| 4-Chlorotoluene | <0.76 | ug/L | 2.5 | 0.76 | 1 | | 07/10/20 10:02 | 106-43-4 | |
| Benzene | <0.25 | ug/L | 1.0 | 0.25 | 1 | | 07/10/20 10:02 | 71-43-2 | |
| Bromobenzene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 10:02 | 108-86-1 | |
| Bromochloromethane | <0.36 | ug/L | 5.0 | 0.36 | 1 | | 07/10/20 10:02 | 74-97-5 | |
| Bromodichloromethane | <0.36 | ug/L | 1.2 | 0.36 | 1 | | 07/10/20 10:02 | 75-27-4 | |
| Bromoform | <4.0 | ug/L | 13.2 | 4.0 | 1 | | 07/10/20 10:02 | 75-25-2 | |
| Bromomethane | <0.97 | ug/L | 5.0 | 0.97 | 1 | | 07/10/20 10:02 | 74-83-9 | |
| Carbon tetrachloride | <1.1 | ug/L | 3.6 | 1.1 | 1 | | 07/10/20 10:02 | 56-23-5 | |
| Chlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 10:02 | 108-90-7 | |
| Chloroethane | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 10:02 | 75-00-3 | |
| Chloroform | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 10:02 | 67-66-3 | |
| Chloromethane | <2.2 | ug/L | 7.3 | 2.2 | 1 | | 07/10/20 10:02 | 74-87-3 | |
| Dibromochloromethane | <2.6 | ug/L | 8.7 | 2.6 | 1 | | 07/10/20 10:02 | 124-48-1 | |
| Dibromomethane | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 10:02 | 74-95-3 | |
| Dichlorodifluoromethane | <0.50 | ug/L | 5.0 | 0.50 | 1 | | 07/10/20 10:02 | 75-71-8 | |
| Ethylbenzene | <0.32 | ug/L | 1.1 | 0.32 | 1 | | 07/10/20 10:02 | 100-41-4 | |
| Hexachloro-1,3-butadiene | <1.5 | ug/L | 4.9 | 1.5 | 1 | | 07/10/20 10:02 | 87-68-3 | |
| Isopropylbenzene (Cumene) | <1.7 | ug/L | 5.6 | 1.7 | 1 | | 07/10/20 10:02 | 98-82-8 | |
| Methyl-tert-butyl ether | <1.2 | ug/L | 4.2 | 1.2 | 1 | | 07/10/20 10:02 | 1634-04-4 | |
| Methylene Chloride | <0.58 | ug/L | 5.0 | 0.58 | 1 | | 07/10/20 10:02 | 75-09-2 | |
| Naphthalene | <1.2 | ug/L | 5.0 | 1.2 | 1 | | 07/10/20 10:02 | 91-20-3 | |
| Styrene | <3.0 | ug/L | 10.0 | 3.0 | 1 | | 07/10/20 10:02 | 100-42-5 | |
| Tetrachloroethene | 1.3 | ug/L | 1.1 | 0.33 | 1 | | 07/10/20 10:02 | 127-18-4 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE
Pace Project No.: 40210819

Sample: MW-4 Lab ID: 40210819004 Collected: 07/08/20 11:15 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|---------------------------|--------------------------------------|-------|--------|------|----|----------|----------------|-------------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| Toluene | <0.27 | ug/L | 0.90 | 0.27 | 1 | | 07/10/20 10:02 | 108-88-3 | |
| Trichloroethene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 10:02 | 79-01-6 | |
| Trichlorofluoromethane | <0.21 | ug/L | 1.0 | 0.21 | 1 | | 07/10/20 10:02 | 75-69-4 | |
| Vinyl chloride | <0.17 | ug/L | 1.0 | 0.17 | 1 | | 07/10/20 10:02 | 75-01-4 | |
| cis-1,2-Dichloroethene | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 10:02 | 156-59-2 | |
| cis-1,3-Dichloropropene | <3.6 | ug/L | 12.1 | 3.6 | 1 | | 07/10/20 10:02 | 10061-01-5 | |
| m&p-Xylene | <0.47 | ug/L | 2.0 | 0.47 | 1 | | 07/10/20 10:02 | 179601-23-1 | |
| n-Butylbenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 10:02 | 104-51-8 | |
| n-Propylbenzene | <0.81 | ug/L | 5.0 | 0.81 | 1 | | 07/10/20 10:02 | 103-65-1 | |
| o-Xylene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 10:02 | 95-47-6 | |
| p-Isopropyltoluene | <0.80 | ug/L | 2.7 | 0.80 | 1 | | 07/10/20 10:02 | 99-87-6 | |
| sec-Butylbenzene | <0.85 | ug/L | 5.0 | 0.85 | 1 | | 07/10/20 10:02 | 135-98-8 | |
| tert-Butylbenzene | <0.30 | ug/L | 1.0 | 0.30 | 1 | | 07/10/20 10:02 | 98-06-6 | |
| trans-1,2-Dichloroethene | <0.46 | ug/L | 1.5 | 0.46 | 1 | | 07/10/20 10:02 | 156-60-5 | |
| trans-1,3-Dichloropropene | <4.4 | ug/L | 14.6 | 4.4 | 1 | | 07/10/20 10:02 | 10061-02-6 | |
| Surrogates | | | | | | | | | |
| 4-Bromofluorobenzene (S) | 89 | % | 70-130 | | 1 | | 07/10/20 10:02 | 460-00-4 | |
| Dibromofluoromethane (S) | 94 | % | 70-130 | | 1 | | 07/10/20 10:02 | 1868-53-7 | |
| Toluene-d8 (S) | 97 | % | 70-130 | | 1 | | 07/10/20 10:02 | 2037-26-5 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

| Sample: MW-5 | Lab ID: 40210819005 | Collected: 07/08/20 10:35 | Received: 07/09/20 09:40 | Matrix: Water | | | | | |
|-----------------------------|--------------------------------------|---------------------------|--------------------------|---------------|----|----------|----------------|-----------|------|
| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 10:24 | 630-20-6 | |
| 1,1,1-Trichloroethane | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 10:24 | 71-55-6 | |
| 1,1,2,2-Tetrachloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 10:24 | 79-34-5 | |
| 1,1,2-Trichloroethane | <0.55 | ug/L | 5.0 | 0.55 | 1 | | 07/10/20 10:24 | 79-00-5 | |
| 1,1-Dichloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 10:24 | 75-34-3 | |
| 1,1-Dichloroethene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 10:24 | 75-35-4 | |
| 1,1-Dichloropropene | <0.54 | ug/L | 1.8 | 0.54 | 1 | | 07/10/20 10:24 | 563-58-6 | |
| 1,2,3-Trichlorobenzene | <2.2 | ug/L | 7.4 | 2.2 | 1 | | 07/10/20 10:24 | 87-61-6 | |
| 1,2,3-Trichloropropane | <0.59 | ug/L | 5.0 | 0.59 | 1 | | 07/10/20 10:24 | 96-18-4 | |
| 1,2,4-Trichlorobenzene | <0.95 | ug/L | 5.0 | 0.95 | 1 | | 07/10/20 10:24 | 120-82-1 | |
| 1,2,4-Trimethylbenzene | <0.84 | ug/L | 2.8 | 0.84 | 1 | | 07/10/20 10:24 | 95-63-6 | |
| 1,2-Dibromo-3-chloropropane | <1.8 | ug/L | 5.9 | 1.8 | 1 | | 07/10/20 10:24 | 96-12-8 | |
| 1,2-Dibromoethane (EDB) | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 10:24 | 106-93-4 | |
| 1,2-Dichlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 10:24 | 95-50-1 | |
| 1,2-Dichloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 10:24 | 107-06-2 | |
| 1,2-Dichloropropane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 10:24 | 78-87-5 | |
| 1,3,5-Trimethylbenzene | <0.87 | ug/L | 2.9 | 0.87 | 1 | | 07/10/20 10:24 | 108-67-8 | |
| 1,3-Dichlorobenzene | <0.63 | ug/L | 2.1 | 0.63 | 1 | | 07/10/20 10:24 | 541-73-1 | |
| 1,3-Dichloropropane | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 10:24 | 142-28-9 | |
| 1,4-Dichlorobenzene | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 10:24 | 106-46-7 | |
| 2,2-Dichloropropane | <2.3 | ug/L | 7.6 | 2.3 | 1 | | 07/10/20 10:24 | 594-20-7 | |
| 2-Chlorotoluene | <0.93 | ug/L | 5.0 | 0.93 | 1 | | 07/10/20 10:24 | 95-49-8 | |
| 4-Chlorotoluene | <0.76 | ug/L | 2.5 | 0.76 | 1 | | 07/10/20 10:24 | 106-43-4 | |
| Benzene | <0.25 | ug/L | 1.0 | 0.25 | 1 | | 07/10/20 10:24 | 71-43-2 | |
| Bromobenzene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 10:24 | 108-86-1 | |
| Bromochloromethane | <0.36 | ug/L | 5.0 | 0.36 | 1 | | 07/10/20 10:24 | 74-97-5 | |
| Bromodichloromethane | <0.36 | ug/L | 1.2 | 0.36 | 1 | | 07/10/20 10:24 | 75-27-4 | |
| Bromoform | <4.0 | ug/L | 13.2 | 4.0 | 1 | | 07/10/20 10:24 | 75-25-2 | |
| Bromomethane | <0.97 | ug/L | 5.0 | 0.97 | 1 | | 07/10/20 10:24 | 74-83-9 | |
| Carbon tetrachloride | <1.1 | ug/L | 3.6 | 1.1 | 1 | | 07/10/20 10:24 | 56-23-5 | |
| Chlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 10:24 | 108-90-7 | |
| Chloroethane | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 10:24 | 75-00-3 | |
| Chloroform | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 10:24 | 67-66-3 | |
| Chloromethane | <2.2 | ug/L | 7.3 | 2.2 | 1 | | 07/10/20 10:24 | 74-87-3 | |
| Dibromochloromethane | <2.6 | ug/L | 8.7 | 2.6 | 1 | | 07/10/20 10:24 | 124-48-1 | |
| Dibromomethane | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 10:24 | 74-95-3 | |
| Dichlorodifluoromethane | <0.50 | ug/L | 5.0 | 0.50 | 1 | | 07/10/20 10:24 | 75-71-8 | |
| Ethylbenzene | <0.32 | ug/L | 1.1 | 0.32 | 1 | | 07/10/20 10:24 | 100-41-4 | |
| Hexachloro-1,3-butadiene | <1.5 | ug/L | 4.9 | 1.5 | 1 | | 07/10/20 10:24 | 87-68-3 | |
| Isopropylbenzene (Cumene) | <1.7 | ug/L | 5.6 | 1.7 | 1 | | 07/10/20 10:24 | 98-82-8 | |
| Methyl-tert-butyl ether | <1.2 | ug/L | 4.2 | 1.2 | 1 | | 07/10/20 10:24 | 1634-04-4 | |
| Methylene Chloride | <0.58 | ug/L | 5.0 | 0.58 | 1 | | 07/10/20 10:24 | 75-09-2 | |
| Naphthalene | <1.2 | ug/L | 5.0 | 1.2 | 1 | | 07/10/20 10:24 | 91-20-3 | |
| Styrene | <3.0 | ug/L | 10.0 | 3.0 | 1 | | 07/10/20 10:24 | 100-42-5 | |
| Tetrachloroethene | <0.33 | ug/L | 1.1 | 0.33 | 1 | | 07/10/20 10:24 | 127-18-4 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: MW-5 Lab ID: 40210819005 Collected: 07/08/20 10:35 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|---------------------------|--------------------------------------|-------|--------|------|----|----------|----------------|-------------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| Toluene | <0.27 | ug/L | 0.90 | 0.27 | 1 | | 07/10/20 10:24 | 108-88-3 | |
| Trichloroethene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 10:24 | 79-01-6 | |
| Trichlorofluoromethane | <0.21 | ug/L | 1.0 | 0.21 | 1 | | 07/10/20 10:24 | 75-69-4 | |
| Vinyl chloride | <0.17 | ug/L | 1.0 | 0.17 | 1 | | 07/10/20 10:24 | 75-01-4 | |
| cis-1,2-Dichloroethene | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 10:24 | 156-59-2 | |
| cis-1,3-Dichloropropene | <3.6 | ug/L | 12.1 | 3.6 | 1 | | 07/10/20 10:24 | 10061-01-5 | |
| m&p-Xylene | <0.47 | ug/L | 2.0 | 0.47 | 1 | | 07/10/20 10:24 | 179601-23-1 | |
| n-Butylbenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 10:24 | 104-51-8 | |
| n-Propylbenzene | <0.81 | ug/L | 5.0 | 0.81 | 1 | | 07/10/20 10:24 | 103-65-1 | |
| o-Xylene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 10:24 | 95-47-6 | |
| p-Isopropyltoluene | <0.80 | ug/L | 2.7 | 0.80 | 1 | | 07/10/20 10:24 | 99-87-6 | |
| sec-Butylbenzene | <0.85 | ug/L | 5.0 | 0.85 | 1 | | 07/10/20 10:24 | 135-98-8 | |
| tert-Butylbenzene | <0.30 | ug/L | 1.0 | 0.30 | 1 | | 07/10/20 10:24 | 98-06-6 | |
| trans-1,2-Dichloroethene | <0.46 | ug/L | 1.5 | 0.46 | 1 | | 07/10/20 10:24 | 156-60-5 | |
| trans-1,3-Dichloropropene | <4.4 | ug/L | 14.6 | 4.4 | 1 | | 07/10/20 10:24 | 10061-02-6 | |
| Surrogates | | | | | | | | | |
| 4-Bromofluorobenzene (S) | 90 | % | 70-130 | | 1 | | 07/10/20 10:24 | 460-00-4 | |
| Dibromofluoromethane (S) | 94 | % | 70-130 | | 1 | | 07/10/20 10:24 | 1868-53-7 | |
| Toluene-d8 (S) | 98 | % | 70-130 | | 1 | | 07/10/20 10:24 | 2037-26-5 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: MW-6 **Lab ID: 40210819006** Collected: 07/08/20 09:50 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|-----------------------------|--------------------------------------|-------|------|------|----|----------|----------------|-----------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 11:53 | 630-20-6 | |
| 1,1,1-Trichloroethane | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 11:53 | 71-55-6 | |
| 1,1,2,2-Tetrachloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 11:53 | 79-34-5 | |
| 1,1,2-Trichloroethane | <0.55 | ug/L | 5.0 | 0.55 | 1 | | 07/10/20 11:53 | 79-00-5 | |
| 1,1-Dichloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 11:53 | 75-34-3 | |
| 1,1-Dichloroethene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 11:53 | 75-35-4 | |
| 1,1-Dichloropropene | <0.54 | ug/L | 1.8 | 0.54 | 1 | | 07/10/20 11:53 | 563-58-6 | |
| 1,2,3-Trichlorobenzene | <2.2 | ug/L | 7.4 | 2.2 | 1 | | 07/10/20 11:53 | 87-61-6 | |
| 1,2,3-Trichloropropane | <0.59 | ug/L | 5.0 | 0.59 | 1 | | 07/10/20 11:53 | 96-18-4 | |
| 1,2,4-Trichlorobenzene | <0.95 | ug/L | 5.0 | 0.95 | 1 | | 07/10/20 11:53 | 120-82-1 | |
| 1,2,4-Trimethylbenzene | <0.84 | ug/L | 2.8 | 0.84 | 1 | | 07/10/20 11:53 | 95-63-6 | |
| 1,2-Dibromo-3-chloropropane | <1.8 | ug/L | 5.9 | 1.8 | 1 | | 07/10/20 11:53 | 96-12-8 | |
| 1,2-Dibromoethane (EDB) | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 11:53 | 106-93-4 | |
| 1,2-Dichlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 11:53 | 95-50-1 | |
| 1,2-Dichloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 11:53 | 107-06-2 | |
| 1,2-Dichloropropane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 11:53 | 78-87-5 | |
| 1,3,5-Trimethylbenzene | <0.87 | ug/L | 2.9 | 0.87 | 1 | | 07/10/20 11:53 | 108-67-8 | |
| 1,3-Dichlorobenzene | <0.63 | ug/L | 2.1 | 0.63 | 1 | | 07/10/20 11:53 | 541-73-1 | |
| 1,3-Dichloropropane | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 11:53 | 142-28-9 | |
| 1,4-Dichlorobenzene | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 11:53 | 106-46-7 | |
| 2,2-Dichloropropane | <2.3 | ug/L | 7.6 | 2.3 | 1 | | 07/10/20 11:53 | 594-20-7 | |
| 2-Chlorotoluene | <0.93 | ug/L | 5.0 | 0.93 | 1 | | 07/10/20 11:53 | 95-49-8 | |
| 4-Chlorotoluene | <0.76 | ug/L | 2.5 | 0.76 | 1 | | 07/10/20 11:53 | 106-43-4 | |
| Benzene | <0.25 | ug/L | 1.0 | 0.25 | 1 | | 07/10/20 11:53 | 71-43-2 | |
| Bromobenzene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 11:53 | 108-86-1 | |
| Bromochloromethane | <0.36 | ug/L | 5.0 | 0.36 | 1 | | 07/10/20 11:53 | 74-97-5 | |
| Bromodichloromethane | <0.36 | ug/L | 1.2 | 0.36 | 1 | | 07/10/20 11:53 | 75-27-4 | |
| Bromoform | <4.0 | ug/L | 13.2 | 4.0 | 1 | | 07/10/20 11:53 | 75-25-2 | |
| Bromomethane | <0.97 | ug/L | 5.0 | 0.97 | 1 | | 07/10/20 11:53 | 74-83-9 | |
| Carbon tetrachloride | <1.1 | ug/L | 3.6 | 1.1 | 1 | | 07/10/20 11:53 | 56-23-5 | |
| Chlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 11:53 | 108-90-7 | |
| Chloroethane | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 11:53 | 75-00-3 | |
| Chloroform | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 11:53 | 67-66-3 | |
| Chloromethane | <2.2 | ug/L | 7.3 | 2.2 | 1 | | 07/10/20 11:53 | 74-87-3 | |
| Dibromochloromethane | <2.6 | ug/L | 8.7 | 2.6 | 1 | | 07/10/20 11:53 | 124-48-1 | |
| Dibromomethane | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 11:53 | 74-95-3 | |
| Dichlorodifluoromethane | <0.50 | ug/L | 5.0 | 0.50 | 1 | | 07/10/20 11:53 | 75-71-8 | |
| Ethylbenzene | <0.32 | ug/L | 1.1 | 0.32 | 1 | | 07/10/20 11:53 | 100-41-4 | |
| Hexachloro-1,3-butadiene | <1.5 | ug/L | 4.9 | 1.5 | 1 | | 07/10/20 11:53 | 87-68-3 | |
| Isopropylbenzene (Cumene) | <1.7 | ug/L | 5.6 | 1.7 | 1 | | 07/10/20 11:53 | 98-82-8 | |
| Methyl-tert-butyl ether | <1.2 | ug/L | 4.2 | 1.2 | 1 | | 07/10/20 11:53 | 1634-04-4 | |
| Methylene Chloride | <0.58 | ug/L | 5.0 | 0.58 | 1 | | 07/10/20 11:53 | 75-09-2 | |
| Naphthalene | <1.2 | ug/L | 5.0 | 1.2 | 1 | | 07/10/20 11:53 | 91-20-3 | |
| Styrene | <3.0 | ug/L | 10.0 | 3.0 | 1 | | 07/10/20 11:53 | 100-42-5 | |
| Tetrachloroethene | 5.3 | ug/L | 1.1 | 0.33 | 1 | | 07/10/20 11:53 | 127-18-4 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: MW-6 **Lab ID: 40210819006** Collected: 07/08/20 09:50 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|---------------------------|--------------------------------------|-------|--------|------|----|----------|----------------|-------------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| Toluene | <0.27 | ug/L | 0.90 | 0.27 | 1 | | 07/10/20 11:53 | 108-88-3 | |
| Trichloroethene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 11:53 | 79-01-6 | |
| Trichlorofluoromethane | <0.21 | ug/L | 1.0 | 0.21 | 1 | | 07/10/20 11:53 | 75-69-4 | |
| Vinyl chloride | <0.17 | ug/L | 1.0 | 0.17 | 1 | | 07/10/20 11:53 | 75-01-4 | |
| cis-1,2-Dichloroethene | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 11:53 | 156-59-2 | |
| cis-1,3-Dichloropropene | <3.6 | ug/L | 12.1 | 3.6 | 1 | | 07/10/20 11:53 | 10061-01-5 | |
| m&p-Xylene | <0.47 | ug/L | 2.0 | 0.47 | 1 | | 07/10/20 11:53 | 179601-23-1 | |
| n-Butylbenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 11:53 | 104-51-8 | |
| n-Propylbenzene | <0.81 | ug/L | 5.0 | 0.81 | 1 | | 07/10/20 11:53 | 103-65-1 | |
| o-Xylene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 11:53 | 95-47-6 | |
| p-Isopropyltoluene | <0.80 | ug/L | 2.7 | 0.80 | 1 | | 07/10/20 11:53 | 99-87-6 | |
| sec-Butylbenzene | <0.85 | ug/L | 5.0 | 0.85 | 1 | | 07/10/20 11:53 | 135-98-8 | |
| tert-Butylbenzene | <0.30 | ug/L | 1.0 | 0.30 | 1 | | 07/10/20 11:53 | 98-06-6 | |
| trans-1,2-Dichloroethene | <0.46 | ug/L | 1.5 | 0.46 | 1 | | 07/10/20 11:53 | 156-60-5 | |
| trans-1,3-Dichloropropene | <4.4 | ug/L | 14.6 | 4.4 | 1 | | 07/10/20 11:53 | 10061-02-6 | |
| Surrogates | | | | | | | | | |
| 4-Bromofluorobenzene (S) | 89 | % | 70-130 | | 1 | | 07/10/20 11:53 | 460-00-4 | |
| Dibromofluoromethane (S) | 95 | % | 70-130 | | 1 | | 07/10/20 11:53 | 1868-53-7 | |
| Toluene-d8 (S) | 97 | % | 70-130 | | 1 | | 07/10/20 11:53 | 2037-26-5 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: MW-6 DUP Lab ID: 40210819007 Collected: 07/08/20 09:50 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|-----------------------------|--------------------------------------|-------|------|------|----|----------|----------------|-----------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 12:14 | 630-20-6 | |
| 1,1,1-Trichloroethane | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 12:14 | 71-55-6 | |
| 1,1,2,2-Tetrachloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 12:14 | 79-34-5 | |
| 1,1,2-Trichloroethane | <0.55 | ug/L | 5.0 | 0.55 | 1 | | 07/10/20 12:14 | 79-00-5 | |
| 1,1-Dichloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 12:14 | 75-34-3 | |
| 1,1-Dichloroethene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 12:14 | 75-35-4 | |
| 1,1-Dichloropropene | <0.54 | ug/L | 1.8 | 0.54 | 1 | | 07/10/20 12:14 | 563-58-6 | |
| 1,2,3-Trichlorobenzene | <2.2 | ug/L | 7.4 | 2.2 | 1 | | 07/10/20 12:14 | 87-61-6 | |
| 1,2,3-Trichloropropane | <0.59 | ug/L | 5.0 | 0.59 | 1 | | 07/10/20 12:14 | 96-18-4 | |
| 1,2,4-Trichlorobenzene | <0.95 | ug/L | 5.0 | 0.95 | 1 | | 07/10/20 12:14 | 120-82-1 | |
| 1,2,4-Trimethylbenzene | <0.84 | ug/L | 2.8 | 0.84 | 1 | | 07/10/20 12:14 | 95-63-6 | |
| 1,2-Dibromo-3-chloropropane | <1.8 | ug/L | 5.9 | 1.8 | 1 | | 07/10/20 12:14 | 96-12-8 | |
| 1,2-Dibromoethane (EDB) | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 12:14 | 106-93-4 | |
| 1,2-Dichlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 12:14 | 95-50-1 | |
| 1,2-Dichloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 12:14 | 107-06-2 | |
| 1,2-Dichloropropane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 12:14 | 78-87-5 | |
| 1,3,5-Trimethylbenzene | <0.87 | ug/L | 2.9 | 0.87 | 1 | | 07/10/20 12:14 | 108-67-8 | |
| 1,3-Dichlorobenzene | <0.63 | ug/L | 2.1 | 0.63 | 1 | | 07/10/20 12:14 | 541-73-1 | |
| 1,3-Dichloropropane | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 12:14 | 142-28-9 | |
| 1,4-Dichlorobenzene | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 12:14 | 106-46-7 | |
| 2,2-Dichloropropane | <2.3 | ug/L | 7.6 | 2.3 | 1 | | 07/10/20 12:14 | 594-20-7 | |
| 2-Chlorotoluene | <0.93 | ug/L | 5.0 | 0.93 | 1 | | 07/10/20 12:14 | 95-49-8 | |
| 4-Chlorotoluene | <0.76 | ug/L | 2.5 | 0.76 | 1 | | 07/10/20 12:14 | 106-43-4 | |
| Benzene | <0.25 | ug/L | 1.0 | 0.25 | 1 | | 07/10/20 12:14 | 71-43-2 | |
| Bromobenzene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 12:14 | 108-86-1 | |
| Bromochloromethane | <0.36 | ug/L | 5.0 | 0.36 | 1 | | 07/10/20 12:14 | 74-97-5 | |
| Bromodichloromethane | <0.36 | ug/L | 1.2 | 0.36 | 1 | | 07/10/20 12:14 | 75-27-4 | |
| Bromoform | <4.0 | ug/L | 13.2 | 4.0 | 1 | | 07/10/20 12:14 | 75-25-2 | |
| Bromomethane | <0.97 | ug/L | 5.0 | 0.97 | 1 | | 07/10/20 12:14 | 74-83-9 | |
| Carbon tetrachloride | <1.1 | ug/L | 3.6 | 1.1 | 1 | | 07/10/20 12:14 | 56-23-5 | |
| Chlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 12:14 | 108-90-7 | |
| Chloroethane | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 12:14 | 75-00-3 | |
| Chloroform | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 12:14 | 67-66-3 | |
| Chloromethane | <2.2 | ug/L | 7.3 | 2.2 | 1 | | 07/10/20 12:14 | 74-87-3 | |
| Dibromochloromethane | <2.6 | ug/L | 8.7 | 2.6 | 1 | | 07/10/20 12:14 | 124-48-1 | |
| Dibromomethane | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 12:14 | 74-95-3 | |
| Dichlorodifluoromethane | <0.50 | ug/L | 5.0 | 0.50 | 1 | | 07/10/20 12:14 | 75-71-8 | |
| Ethylbenzene | <0.32 | ug/L | 1.1 | 0.32 | 1 | | 07/10/20 12:14 | 100-41-4 | |
| Hexachloro-1,3-butadiene | <1.5 | ug/L | 4.9 | 1.5 | 1 | | 07/10/20 12:14 | 87-68-3 | |
| Isopropylbenzene (Cumene) | <1.7 | ug/L | 5.6 | 1.7 | 1 | | 07/10/20 12:14 | 98-82-8 | |
| Methyl-tert-butyl ether | <1.2 | ug/L | 4.2 | 1.2 | 1 | | 07/10/20 12:14 | 1634-04-4 | |
| Methylene Chloride | <0.58 | ug/L | 5.0 | 0.58 | 1 | | 07/10/20 12:14 | 75-09-2 | |
| Naphthalene | <1.2 | ug/L | 5.0 | 1.2 | 1 | | 07/10/20 12:14 | 91-20-3 | |
| Styrene | <3.0 | ug/L | 10.0 | 3.0 | 1 | | 07/10/20 12:14 | 100-42-5 | |
| Tetrachloroethene | 5.2 | ug/L | 1.1 | 0.33 | 1 | | 07/10/20 12:14 | 127-18-4 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: MW-6 DUP Lab ID: 40210819007 Collected: 07/08/20 09:50 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|---------------------------|--------------------------------------|-------|--------|------|----|----------|----------------|-------------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| Toluene | <0.27 | ug/L | 0.90 | 0.27 | 1 | | 07/10/20 12:14 | 108-88-3 | |
| Trichloroethene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 12:14 | 79-01-6 | |
| Trichlorofluoromethane | <0.21 | ug/L | 1.0 | 0.21 | 1 | | 07/10/20 12:14 | 75-69-4 | |
| Vinyl chloride | <0.17 | ug/L | 1.0 | 0.17 | 1 | | 07/10/20 12:14 | 75-01-4 | |
| cis-1,2-Dichloroethene | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 12:14 | 156-59-2 | |
| cis-1,3-Dichloropropene | <3.6 | ug/L | 12.1 | 3.6 | 1 | | 07/10/20 12:14 | 10061-01-5 | |
| m&p-Xylene | <0.47 | ug/L | 2.0 | 0.47 | 1 | | 07/10/20 12:14 | 179601-23-1 | |
| n-Butylbenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 12:14 | 104-51-8 | |
| n-Propylbenzene | <0.81 | ug/L | 5.0 | 0.81 | 1 | | 07/10/20 12:14 | 103-65-1 | |
| o-Xylene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 12:14 | 95-47-6 | |
| p-Isopropyltoluene | <0.80 | ug/L | 2.7 | 0.80 | 1 | | 07/10/20 12:14 | 99-87-6 | |
| sec-Butylbenzene | <0.85 | ug/L | 5.0 | 0.85 | 1 | | 07/10/20 12:14 | 135-98-8 | |
| tert-Butylbenzene | <0.30 | ug/L | 1.0 | 0.30 | 1 | | 07/10/20 12:14 | 98-06-6 | |
| trans-1,2-Dichloroethene | <0.46 | ug/L | 1.5 | 0.46 | 1 | | 07/10/20 12:14 | 156-60-5 | |
| trans-1,3-Dichloropropene | <4.4 | ug/L | 14.6 | 4.4 | 1 | | 07/10/20 12:14 | 10061-02-6 | |
| Surrogates | | | | | | | | | |
| 4-Bromofluorobenzene (S) | 89 | % | 70-130 | | 1 | | 07/10/20 12:14 | 460-00-4 | |
| Dibromofluoromethane (S) | 94 | % | 70-130 | | 1 | | 07/10/20 12:14 | 1868-53-7 | |
| Toluene-d8 (S) | 98 | % | 70-130 | | 1 | | 07/10/20 12:14 | 2037-26-5 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: MW-7 **Lab ID: 40210819008** Collected: 07/08/20 09:10 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|-----------------------------|--------------------------------------|-------|------|------|----|----------|----------------|-----------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 10:46 | 630-20-6 | |
| 1,1,1-Trichloroethane | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 10:46 | 71-55-6 | |
| 1,1,2,2-Tetrachloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 10:46 | 79-34-5 | |
| 1,1,2-Trichloroethane | <0.55 | ug/L | 5.0 | 0.55 | 1 | | 07/10/20 10:46 | 79-00-5 | |
| 1,1-Dichloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 10:46 | 75-34-3 | |
| 1,1-Dichloroethene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 10:46 | 75-35-4 | |
| 1,1-Dichloropropene | <0.54 | ug/L | 1.8 | 0.54 | 1 | | 07/10/20 10:46 | 563-58-6 | |
| 1,2,3-Trichlorobenzene | <2.2 | ug/L | 7.4 | 2.2 | 1 | | 07/10/20 10:46 | 87-61-6 | |
| 1,2,3-Trichloropropane | <0.59 | ug/L | 5.0 | 0.59 | 1 | | 07/10/20 10:46 | 96-18-4 | |
| 1,2,4-Trichlorobenzene | <0.95 | ug/L | 5.0 | 0.95 | 1 | | 07/10/20 10:46 | 120-82-1 | |
| 1,2,4-Trimethylbenzene | <0.84 | ug/L | 2.8 | 0.84 | 1 | | 07/10/20 10:46 | 95-63-6 | |
| 1,2-Dibromo-3-chloropropane | <1.8 | ug/L | 5.9 | 1.8 | 1 | | 07/10/20 10:46 | 96-12-8 | |
| 1,2-Dibromoethane (EDB) | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 10:46 | 106-93-4 | |
| 1,2-Dichlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 10:46 | 95-50-1 | |
| 1,2-Dichloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 10:46 | 107-06-2 | |
| 1,2-Dichloropropane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 10:46 | 78-87-5 | |
| 1,3,5-Trimethylbenzene | <0.87 | ug/L | 2.9 | 0.87 | 1 | | 07/10/20 10:46 | 108-67-8 | |
| 1,3-Dichlorobenzene | <0.63 | ug/L | 2.1 | 0.63 | 1 | | 07/10/20 10:46 | 541-73-1 | |
| 1,3-Dichloropropane | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 10:46 | 142-28-9 | |
| 1,4-Dichlorobenzene | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 10:46 | 106-46-7 | |
| 2,2-Dichloropropane | <2.3 | ug/L | 7.6 | 2.3 | 1 | | 07/10/20 10:46 | 594-20-7 | |
| 2-Chlorotoluene | <0.93 | ug/L | 5.0 | 0.93 | 1 | | 07/10/20 10:46 | 95-49-8 | |
| 4-Chlorotoluene | <0.76 | ug/L | 2.5 | 0.76 | 1 | | 07/10/20 10:46 | 106-43-4 | |
| Benzene | <0.25 | ug/L | 1.0 | 0.25 | 1 | | 07/10/20 10:46 | 71-43-2 | |
| Bromobenzene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 10:46 | 108-86-1 | |
| Bromochloromethane | <0.36 | ug/L | 5.0 | 0.36 | 1 | | 07/10/20 10:46 | 74-97-5 | |
| Bromodichloromethane | <0.36 | ug/L | 1.2 | 0.36 | 1 | | 07/10/20 10:46 | 75-27-4 | |
| Bromoform | <4.0 | ug/L | 13.2 | 4.0 | 1 | | 07/10/20 10:46 | 75-25-2 | |
| Bromomethane | <0.97 | ug/L | 5.0 | 0.97 | 1 | | 07/10/20 10:46 | 74-83-9 | |
| Carbon tetrachloride | <1.1 | ug/L | 3.6 | 1.1 | 1 | | 07/10/20 10:46 | 56-23-5 | |
| Chlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 10:46 | 108-90-7 | |
| Chloroethane | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 10:46 | 75-00-3 | |
| Chloroform | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 10:46 | 67-66-3 | |
| Chloromethane | <2.2 | ug/L | 7.3 | 2.2 | 1 | | 07/10/20 10:46 | 74-87-3 | |
| Dibromochloromethane | <2.6 | ug/L | 8.7 | 2.6 | 1 | | 07/10/20 10:46 | 124-48-1 | |
| Dibromomethane | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 10:46 | 74-95-3 | |
| Dichlorodifluoromethane | <0.50 | ug/L | 5.0 | 0.50 | 1 | | 07/10/20 10:46 | 75-71-8 | |
| Ethylbenzene | <0.32 | ug/L | 1.1 | 0.32 | 1 | | 07/10/20 10:46 | 100-41-4 | |
| Hexachloro-1,3-butadiene | <1.5 | ug/L | 4.9 | 1.5 | 1 | | 07/10/20 10:46 | 87-68-3 | |
| Isopropylbenzene (Cumene) | <1.7 | ug/L | 5.6 | 1.7 | 1 | | 07/10/20 10:46 | 98-82-8 | |
| Methyl-tert-butyl ether | <1.2 | ug/L | 4.2 | 1.2 | 1 | | 07/10/20 10:46 | 1634-04-4 | |
| Methylene Chloride | <0.58 | ug/L | 5.0 | 0.58 | 1 | | 07/10/20 10:46 | 75-09-2 | |
| Naphthalene | <1.2 | ug/L | 5.0 | 1.2 | 1 | | 07/10/20 10:46 | 91-20-3 | |
| Styrene | <3.0 | ug/L | 10.0 | 3.0 | 1 | | 07/10/20 10:46 | 100-42-5 | |
| Tetrachloroethene | 0.80J | ug/L | 1.1 | 0.33 | 1 | | 07/10/20 10:46 | 127-18-4 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: MW-7 **Lab ID: 40210819008** Collected: 07/08/20 09:10 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|---------------------------|--------------------------------------|-------|--------|------|----|----------|----------------|-------------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| Toluene | <0.27 | ug/L | 0.90 | 0.27 | 1 | | 07/10/20 10:46 | 108-88-3 | |
| Trichloroethene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 10:46 | 79-01-6 | |
| Trichlorofluoromethane | <0.21 | ug/L | 1.0 | 0.21 | 1 | | 07/10/20 10:46 | 75-69-4 | |
| Vinyl chloride | <0.17 | ug/L | 1.0 | 0.17 | 1 | | 07/10/20 10:46 | 75-01-4 | |
| cis-1,2-Dichloroethene | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 10:46 | 156-59-2 | |
| cis-1,3-Dichloropropene | <3.6 | ug/L | 12.1 | 3.6 | 1 | | 07/10/20 10:46 | 10061-01-5 | |
| m&p-Xylene | <0.47 | ug/L | 2.0 | 0.47 | 1 | | 07/10/20 10:46 | 179601-23-1 | |
| n-Butylbenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 10:46 | 104-51-8 | |
| n-Propylbenzene | <0.81 | ug/L | 5.0 | 0.81 | 1 | | 07/10/20 10:46 | 103-65-1 | |
| o-Xylene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 10:46 | 95-47-6 | |
| p-Isopropyltoluene | <0.80 | ug/L | 2.7 | 0.80 | 1 | | 07/10/20 10:46 | 99-87-6 | |
| sec-Butylbenzene | <0.85 | ug/L | 5.0 | 0.85 | 1 | | 07/10/20 10:46 | 135-98-8 | |
| tert-Butylbenzene | <0.30 | ug/L | 1.0 | 0.30 | 1 | | 07/10/20 10:46 | 98-06-6 | |
| trans-1,2-Dichloroethene | <0.46 | ug/L | 1.5 | 0.46 | 1 | | 07/10/20 10:46 | 156-60-5 | |
| trans-1,3-Dichloropropene | <4.4 | ug/L | 14.6 | 4.4 | 1 | | 07/10/20 10:46 | 10061-02-6 | |
| Surrogates | | | | | | | | | |
| 4-Bromofluorobenzene (S) | 89 | % | 70-130 | | 1 | | 07/10/20 10:46 | 460-00-4 | |
| Dibromofluoromethane (S) | 96 | % | 70-130 | | 1 | | 07/10/20 10:46 | 1868-53-7 | |
| Toluene-d8 (S) | 96 | % | 70-130 | | 1 | | 07/10/20 10:46 | 2037-26-5 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: TRIP BLANK Lab ID: 40210819009 Collected: 07/08/20 00:00 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|-----------------------------|--------------------------------------|-------|------|------|----|----------|----------------|-----------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 09:40 | 630-20-6 | |
| 1,1,1-Trichloroethane | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 09:40 | 71-55-6 | |
| 1,1,2,2-Tetrachloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 09:40 | 79-34-5 | |
| 1,1,2-Trichloroethane | <0.55 | ug/L | 5.0 | 0.55 | 1 | | 07/10/20 09:40 | 79-00-5 | |
| 1,1-Dichloroethane | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 09:40 | 75-34-3 | |
| 1,1-Dichloroethene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 09:40 | 75-35-4 | |
| 1,1-Dichloropropene | <0.54 | ug/L | 1.8 | 0.54 | 1 | | 07/10/20 09:40 | 563-58-6 | |
| 1,2,3-Trichlorobenzene | <2.2 | ug/L | 7.4 | 2.2 | 1 | | 07/10/20 09:40 | 87-61-6 | |
| 1,2,3-Trichloropropane | <0.59 | ug/L | 5.0 | 0.59 | 1 | | 07/10/20 09:40 | 96-18-4 | |
| 1,2,4-Trichlorobenzene | <0.95 | ug/L | 5.0 | 0.95 | 1 | | 07/10/20 09:40 | 120-82-1 | |
| 1,2,4-Trimethylbenzene | <0.84 | ug/L | 2.8 | 0.84 | 1 | | 07/10/20 09:40 | 95-63-6 | |
| 1,2-Dibromo-3-chloropropane | <1.8 | ug/L | 5.9 | 1.8 | 1 | | 07/10/20 09:40 | 96-12-8 | |
| 1,2-Dibromoethane (EDB) | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 09:40 | 106-93-4 | |
| 1,2-Dichlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 09:40 | 95-50-1 | |
| 1,2-Dichloroethane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 09:40 | 107-06-2 | |
| 1,2-Dichloropropane | <0.28 | ug/L | 1.0 | 0.28 | 1 | | 07/10/20 09:40 | 78-87-5 | |
| 1,3,5-Trimethylbenzene | <0.87 | ug/L | 2.9 | 0.87 | 1 | | 07/10/20 09:40 | 108-67-8 | |
| 1,3-Dichlorobenzene | <0.63 | ug/L | 2.1 | 0.63 | 1 | | 07/10/20 09:40 | 541-73-1 | |
| 1,3-Dichloropropane | <0.83 | ug/L | 2.8 | 0.83 | 1 | | 07/10/20 09:40 | 142-28-9 | |
| 1,4-Dichlorobenzene | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 09:40 | 106-46-7 | |
| 2,2-Dichloropropane | <2.3 | ug/L | 7.6 | 2.3 | 1 | | 07/10/20 09:40 | 594-20-7 | |
| 2-Chlorotoluene | <0.93 | ug/L | 5.0 | 0.93 | 1 | | 07/10/20 09:40 | 95-49-8 | |
| 4-Chlorotoluene | <0.76 | ug/L | 2.5 | 0.76 | 1 | | 07/10/20 09:40 | 106-43-4 | |
| Benzene | <0.25 | ug/L | 1.0 | 0.25 | 1 | | 07/10/20 09:40 | 71-43-2 | |
| Bromobenzene | <0.24 | ug/L | 1.0 | 0.24 | 1 | | 07/10/20 09:40 | 108-86-1 | |
| Bromochloromethane | <0.36 | ug/L | 5.0 | 0.36 | 1 | | 07/10/20 09:40 | 74-97-5 | |
| Bromodichloromethane | <0.36 | ug/L | 1.2 | 0.36 | 1 | | 07/10/20 09:40 | 75-27-4 | |
| Bromoform | <4.0 | ug/L | 13.2 | 4.0 | 1 | | 07/10/20 09:40 | 75-25-2 | |
| Bromomethane | <0.97 | ug/L | 5.0 | 0.97 | 1 | | 07/10/20 09:40 | 74-83-9 | |
| Carbon tetrachloride | <1.1 | ug/L | 3.6 | 1.1 | 1 | | 07/10/20 09:40 | 56-23-5 | |
| Chlorobenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 09:40 | 108-90-7 | |
| Chloroethane | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 09:40 | 75-00-3 | |
| Chloroform | <1.3 | ug/L | 5.0 | 1.3 | 1 | | 07/10/20 09:40 | 67-66-3 | |
| Chloromethane | <2.2 | ug/L | 7.3 | 2.2 | 1 | | 07/10/20 09:40 | 74-87-3 | |
| Dibromochloromethane | <2.6 | ug/L | 8.7 | 2.6 | 1 | | 07/10/20 09:40 | 124-48-1 | |
| Dibromomethane | <0.94 | ug/L | 3.1 | 0.94 | 1 | | 07/10/20 09:40 | 74-95-3 | |
| Dichlorodifluoromethane | <0.50 | ug/L | 5.0 | 0.50 | 1 | | 07/10/20 09:40 | 75-71-8 | |
| Ethylbenzene | <0.32 | ug/L | 1.1 | 0.32 | 1 | | 07/10/20 09:40 | 100-41-4 | |
| Hexachloro-1,3-butadiene | <1.5 | ug/L | 4.9 | 1.5 | 1 | | 07/10/20 09:40 | 87-68-3 | |
| Isopropylbenzene (Cumene) | <1.7 | ug/L | 5.6 | 1.7 | 1 | | 07/10/20 09:40 | 98-82-8 | |
| Methyl-tert-butyl ether | <1.2 | ug/L | 4.2 | 1.2 | 1 | | 07/10/20 09:40 | 1634-04-4 | |
| Methylene Chloride | <0.58 | ug/L | 5.0 | 0.58 | 1 | | 07/10/20 09:40 | 75-09-2 | |
| Naphthalene | <1.2 | ug/L | 5.0 | 1.2 | 1 | | 07/10/20 09:40 | 91-20-3 | |
| Styrene | <3.0 | ug/L | 10.0 | 3.0 | 1 | | 07/10/20 09:40 | 100-42-5 | |
| Tetrachloroethene | <0.33 | ug/L | 1.1 | 0.33 | 1 | | 07/10/20 09:40 | 127-18-4 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

ANALYTICAL RESULTS

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

Sample: TRIP BLANK Lab ID: 40210819009 Collected: 07/08/20 00:00 Received: 07/09/20 09:40 Matrix: Water

| Parameters | Results | Units | LOQ | LOD | DF | Prepared | Analyzed | CAS No. | Qual |
|---------------------------|--------------------------------------|-------|--------|------|----|----------|----------------|-------------|------|
| 8260 MSV | Analytical Method: EPA 8260 | | | | | | | | |
| | Pace Analytical Services - Green Bay | | | | | | | | |
| Toluene | <0.27 | ug/L | 0.90 | 0.27 | 1 | | 07/10/20 09:40 | 108-88-3 | |
| Trichloroethene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 09:40 | 79-01-6 | |
| Trichlorofluoromethane | <0.21 | ug/L | 1.0 | 0.21 | 1 | | 07/10/20 09:40 | 75-69-4 | |
| Vinyl chloride | <0.17 | ug/L | 1.0 | 0.17 | 1 | | 07/10/20 09:40 | 75-01-4 | |
| cis-1,2-Dichloroethene | <0.27 | ug/L | 1.0 | 0.27 | 1 | | 07/10/20 09:40 | 156-59-2 | |
| cis-1,3-Dichloropropene | <3.6 | ug/L | 12.1 | 3.6 | 1 | | 07/10/20 09:40 | 10061-01-5 | |
| m&p-Xylene | <0.47 | ug/L | 2.0 | 0.47 | 1 | | 07/10/20 09:40 | 179601-23-1 | |
| n-Butylbenzene | <0.71 | ug/L | 2.4 | 0.71 | 1 | | 07/10/20 09:40 | 104-51-8 | |
| n-Propylbenzene | <0.81 | ug/L | 5.0 | 0.81 | 1 | | 07/10/20 09:40 | 103-65-1 | |
| o-Xylene | <0.26 | ug/L | 1.0 | 0.26 | 1 | | 07/10/20 09:40 | 95-47-6 | |
| p-Isopropyltoluene | <0.80 | ug/L | 2.7 | 0.80 | 1 | | 07/10/20 09:40 | 99-87-6 | |
| sec-Butylbenzene | <0.85 | ug/L | 5.0 | 0.85 | 1 | | 07/10/20 09:40 | 135-98-8 | |
| tert-Butylbenzene | <0.30 | ug/L | 1.0 | 0.30 | 1 | | 07/10/20 09:40 | 98-06-6 | |
| trans-1,2-Dichloroethene | <0.46 | ug/L | 1.5 | 0.46 | 1 | | 07/10/20 09:40 | 156-60-5 | |
| trans-1,3-Dichloropropene | <4.4 | ug/L | 14.6 | 4.4 | 1 | | 07/10/20 09:40 | 10061-02-6 | |
| Surrogates | | | | | | | | | |
| 4-Bromofluorobenzene (S) | 91 | % | 70-130 | | 1 | | 07/10/20 09:40 | 460-00-4 | HS |
| Dibromofluoromethane (S) | 88 | % | 70-130 | | 1 | | 07/10/20 09:40 | 1868-53-7 | |
| Toluene-d8 (S) | 98 | % | 70-130 | | 1 | | 07/10/20 09:40 | 2037-26-5 | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

QUALITY CONTROL DATA

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

QC Batch: 359774 Analysis Method: EPA 8260

QC Batch Method: EPA 8260 Analysis Description: 8260 MSV

Laboratory: Pace Analytical Services - Green Bay

Associated Lab Samples: 40210819001, 40210819002, 40210819003, 40210819004, 40210819005, 40210819006, 40210819007,
40210819008, 40210819009

METHOD BLANK: 2080653

Matrix: Water

Associated Lab Samples: 40210819001, 40210819002, 40210819003, 40210819004, 40210819005, 40210819006, 40210819007,
40210819008, 40210819009

| Parameter | Units | Blank | Reporting | | Qualifiers |
|-----------------------------|-------|--------|-----------|----------------|------------|
| | | Result | Limit | Analyzed | |
| 1,1,1,2-Tetrachloroethane | ug/L | <0.27 | 1.0 | 07/10/20 07:06 | |
| 1,1,1-Trichloroethane | ug/L | <0.24 | 1.0 | 07/10/20 07:06 | |
| 1,1,2,2-Tetrachloroethane | ug/L | <0.28 | 1.0 | 07/10/20 07:06 | |
| 1,1,2-Trichloroethane | ug/L | <0.55 | 5.0 | 07/10/20 07:06 | |
| 1,1-Dichloroethane | ug/L | <0.27 | 1.0 | 07/10/20 07:06 | |
| 1,1-Dichloroethene | ug/L | <0.24 | 1.0 | 07/10/20 07:06 | |
| 1,1-Dichloropropene | ug/L | <0.54 | 1.8 | 07/10/20 07:06 | |
| 1,2,3-Trichlorobenzene | ug/L | <2.2 | 7.4 | 07/10/20 07:06 | |
| 1,2,3-Trichloropropane | ug/L | <0.59 | 5.0 | 07/10/20 07:06 | |
| 1,2,4-Trichlorobenzene | ug/L | <0.95 | 5.0 | 07/10/20 07:06 | |
| 1,2,4-Trimethylbenzene | ug/L | <0.84 | 2.8 | 07/10/20 07:06 | |
| 1,2-Dibromo-3-chloropropane | ug/L | <1.8 | 5.9 | 07/10/20 07:06 | |
| 1,2-Dibromoethane (EDB) | ug/L | <0.83 | 2.8 | 07/10/20 07:06 | |
| 1,2-Dichlorobenzene | ug/L | <0.71 | 2.4 | 07/10/20 07:06 | |
| 1,2-Dichloroethane | ug/L | <0.28 | 1.0 | 07/10/20 07:06 | |
| 1,2-Dichloropropane | ug/L | <0.28 | 1.0 | 07/10/20 07:06 | |
| 1,3,5-Trimethylbenzene | ug/L | <0.87 | 2.9 | 07/10/20 07:06 | |
| 1,3-Dichlorobenzene | ug/L | <0.63 | 2.1 | 07/10/20 07:06 | |
| 1,3-Dichloropropane | ug/L | <0.83 | 2.8 | 07/10/20 07:06 | |
| 1,4-Dichlorobenzene | ug/L | <0.94 | 3.1 | 07/10/20 07:06 | |
| 2,2-Dichloropropane | ug/L | <2.3 | 7.6 | 07/10/20 07:06 | |
| 2-Chlorotoluene | ug/L | <0.93 | 5.0 | 07/10/20 07:06 | |
| 4-Chlorotoluene | ug/L | <0.76 | 2.5 | 07/10/20 07:06 | |
| Benzene | ug/L | <0.25 | 1.0 | 07/10/20 07:06 | |
| Bromobenzene | ug/L | <0.24 | 1.0 | 07/10/20 07:06 | |
| Bromochloromethane | ug/L | <0.36 | 5.0 | 07/10/20 07:06 | |
| Bromodichloromethane | ug/L | <0.36 | 1.2 | 07/10/20 07:06 | |
| Bromoform | ug/L | <4.0 | 13.2 | 07/10/20 07:06 | |
| Bromomethane | ug/L | <0.97 | 5.0 | 07/10/20 07:06 | |
| Carbon tetrachloride | ug/L | <1.1 | 3.6 | 07/10/20 07:06 | |
| Chlorobenzene | ug/L | <0.71 | 2.4 | 07/10/20 07:06 | |
| Chloroethane | ug/L | <1.3 | 5.0 | 07/10/20 07:06 | |
| Chloroform | ug/L | <1.3 | 5.0 | 07/10/20 07:06 | |
| Chloromethane | ug/L | <2.2 | 7.3 | 07/10/20 07:06 | |
| cis-1,2-Dichloroethene | ug/L | <0.27 | 1.0 | 07/10/20 07:06 | |
| cis-1,3-Dichloropropene | ug/L | <3.6 | 12.1 | 07/10/20 07:06 | |
| Dibromochloromethane | ug/L | <2.6 | 8.7 | 07/10/20 07:06 | |
| Dibromomethane | ug/L | <0.94 | 3.1 | 07/10/20 07:06 | |
| Dichlorodifluoromethane | ug/L | <0.50 | 5.0 | 07/10/20 07:06 | |

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,

without the written consent of Pace Analytical Services, LLC.

QUALITY CONTROL DATA

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

METHOD BLANK: 2080653

Matrix: Water

Associated Lab Samples: 40210819001, 40210819002, 40210819003, 40210819004, 40210819005, 40210819006, 40210819007,
40210819008, 40210819009

| Parameter | Units | Blank Result | Reporting Limit | Analyzed | Qualifiers |
|---------------------------|-------|--------------|-----------------|----------------|------------|
| Ethylbenzene | ug/L | <0.32 | 1.1 | 07/10/20 07:06 | |
| Hexachloro-1,3-butadiene | ug/L | <1.5 | 4.9 | 07/10/20 07:06 | |
| Isopropylbenzene (Cumene) | ug/L | <1.7 | 5.6 | 07/10/20 07:06 | |
| m&p-Xylene | ug/L | <0.47 | 2.0 | 07/10/20 07:06 | |
| Methyl-tert-butyl ether | ug/L | <1.2 | 4.2 | 07/10/20 07:06 | |
| Methylene Chloride | ug/L | <0.58 | 5.0 | 07/10/20 07:06 | |
| n-Butylbenzene | ug/L | <0.71 | 2.4 | 07/10/20 07:06 | |
| n-Propylbenzene | ug/L | <0.81 | 5.0 | 07/10/20 07:06 | |
| Naphthalene | ug/L | <1.2 | 5.0 | 07/10/20 07:06 | |
| o-Xylene | ug/L | <0.26 | 1.0 | 07/10/20 07:06 | |
| p-Isopropyltoluene | ug/L | <0.80 | 2.7 | 07/10/20 07:06 | |
| sec-Butylbenzene | ug/L | <0.85 | 5.0 | 07/10/20 07:06 | |
| Styrene | ug/L | <3.0 | 10.0 | 07/10/20 07:06 | |
| tert-Butylbenzene | ug/L | <0.30 | 1.0 | 07/10/20 07:06 | |
| Tetrachloroethene | ug/L | <0.33 | 1.1 | 07/10/20 07:06 | |
| Toluene | ug/L | <0.27 | 0.90 | 07/10/20 07:06 | |
| trans-1,2-Dichloroethene | ug/L | <0.46 | 1.5 | 07/10/20 07:06 | |
| trans-1,3-Dichloropropene | ug/L | <4.4 | 14.6 | 07/10/20 07:06 | |
| Trichloroethene | ug/L | <0.26 | 1.0 | 07/10/20 07:06 | |
| Trichlorofluoromethane | ug/L | <0.21 | 1.0 | 07/10/20 07:06 | |
| Vinyl chloride | ug/L | <0.17 | 1.0 | 07/10/20 07:06 | |
| 4-Bromofluorobenzene (S) | % | 90 | 70-130 | 07/10/20 07:06 | |
| Dibromofluoromethane (S) | % | 92 | 70-130 | 07/10/20 07:06 | |
| Toluene-d8 (S) | % | 97 | 70-130 | 07/10/20 07:06 | |

LABORATORY CONTROL SAMPLE: 2080654

| Parameter | Units | Spike Conc. | LCS Result | LCS % Rec | % Rec Limits | Qualifiers |
|-----------------------------|-------|-------------|------------|-----------|--------------|------------|
| 1,1,1-Trichloroethane | ug/L | 50 | 50.9 | 102 | 70-130 | |
| 1,1,2,2-Tetrachloroethane | ug/L | 50 | 44.5 | 89 | 64-131 | |
| 1,1,2-Trichloroethane | ug/L | 50 | 48.7 | 97 | 70-130 | |
| 1,1-Dichloroethane | ug/L | 50 | 44.4 | 89 | 69-163 | |
| 1,1-Dichloroethene | ug/L | 50 | 49.7 | 99 | 77-123 | |
| 1,2,4-Trichlorobenzene | ug/L | 50 | 48.6 | 97 | 68-130 | |
| 1,2-Dibromo-3-chloropropane | ug/L | 50 | 42.7 | 85 | 63-130 | |
| 1,2-Dibromoethane (EDB) | ug/L | 50 | 50.1 | 100 | 70-130 | |
| 1,2-Dichlorobenzene | ug/L | 50 | 50.7 | 101 | 70-130 | |
| 1,2-Dichloroethane | ug/L | 50 | 40.5 | 81 | 78-142 | |
| 1,2-Dichloropropane | ug/L | 50 | 45.2 | 90 | 86-134 | |
| 1,3-Dichlorobenzene | ug/L | 50 | 51.1 | 102 | 70-130 | |
| 1,4-Dichlorobenzene | ug/L | 50 | 51.9 | 104 | 70-130 | |
| Benzene | ug/L | 50 | 45.7 | 91 | 70-130 | |
| Bromodichloromethane | ug/L | 50 | 46.5 | 93 | 70-130 | |

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,

without the written consent of Pace Analytical Services, LLC.

QUALITY CONTROL DATA

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

LABORATORY CONTROL SAMPLE: 2080654

| Parameter | Units | Spike Conc. | LCS Result | LCS % Rec | % Rec Limits | Qualifiers |
|---------------------------|-------|-------------|------------|-----------|--------------|------------|
| Bromoform | ug/L | 50 | 46.7 | 93 | 70-130 | |
| Bromomethane | ug/L | 50 | 35.6 | 71 | 39-129 | |
| Carbon tetrachloride | ug/L | 50 | 46.2 | 92 | 70-132 | |
| Chlorobenzene | ug/L | 50 | 53.2 | 106 | 70-130 | |
| Chloroethane | ug/L | 50 | 43.9 | 88 | 66-140 | |
| Chloroform | ug/L | 50 | 48.8 | 98 | 75-132 | |
| Chloromethane | ug/L | 50 | 36.9 | 74 | 32-143 | |
| cis-1,2-Dichloroethene | ug/L | 50 | 47.5 | 95 | 70-130 | |
| cis-1,3-Dichloropropene | ug/L | 50 | 46.1 | 92 | 70-130 | |
| Dibromochloromethane | ug/L | 50 | 45.9 | 92 | 70-130 | |
| Dichlorodifluoromethane | ug/L | 50 | 46.1 | 92 | 10-141 | |
| Ethylbenzene | ug/L | 50 | 51.7 | 103 | 80-120 | |
| Isopropylbenzene (Cumene) | ug/L | 50 | 53.3 | 107 | 70-130 | |
| m&p-Xylene | ug/L | 100 | 106 | 106 | 70-130 | |
| Methyl-tert-butyl ether | ug/L | 50 | 40.2 | 80 | 61-129 | |
| Methylene Chloride | ug/L | 50 | 45.4 | 91 | 70-130 | |
| o-Xylene | ug/L | 50 | 51.7 | 103 | 70-130 | |
| Styrene | ug/L | 50 | 51.6 | 103 | 70-130 | |
| Tetrachloroethene | ug/L | 50 | 49.6 | 99 | 70-130 | |
| Toluene | ug/L | 50 | 51.6 | 103 | 80-120 | |
| trans-1,2-Dichloroethene | ug/L | 50 | 49.9 | 100 | 70-130 | |
| trans-1,3-Dichloropropene | ug/L | 50 | 43.9 | 88 | 69-130 | |
| Trichloroethene | ug/L | 50 | 52.0 | 104 | 70-130 | |
| Trichlorofluoromethane | ug/L | 50 | 51.9 | 104 | 75-145 | |
| Vinyl chloride | ug/L | 50 | 48.9 | 98 | 51-140 | |
| 4-Bromofluorobenzene (S) | % | | | 94 | 70-130 | |
| Dibromofluoromethane (S) | % | | | 93 | 70-130 | |
| Toluene-d8 (S) | % | | | 98 | 70-130 | |

MATRIX SPIKE & MATRIX SPIKE DUPLICATE: 2081562 2081563

| Parameter | Units | MS | | MSD | | MS | | MSD | | % Rec Limits | RPD | Max RPD | Qual |
|-----------------------------|-------|-------------|--------|-------------|-------------|-----------|------------|-------|--------|--------------|-----|---------|------|
| | | 40210819002 | Result | Spike Conc. | Spike Conc. | MS Result | MSD Result | % Rec | % Rec | | | | |
| 1,1,1-Trichloroethane | ug/L | <0.24 | 50 | 50 | 51.4 | 52.0 | 103 | 104 | 70-130 | 1 | 20 | | |
| 1,1,2,2-Tetrachloroethane | ug/L | <0.28 | 50 | 50 | 54.2 | 52.5 | 108 | 105 | 64-137 | 3 | 20 | | |
| 1,1,2-Trichloroethane | ug/L | <0.55 | 50 | 50 | 51.1 | 51.7 | 102 | 103 | 70-137 | 1 | 20 | | |
| 1,1-Dichloroethane | ug/L | <0.27 | 50 | 50 | 45.0 | 45.0 | 90 | 90 | 69-163 | 0 | 20 | | |
| 1,1-Dichloroethene | ug/L | <0.24 | 50 | 50 | 49.2 | 50.1 | 98 | 100 | 77-129 | 2 | 20 | | |
| 1,2,4-Trichlorobenzene | ug/L | <0.95 | 50 | 50 | 48.5 | 49.9 | 97 | 100 | 68-130 | 3 | 20 | | |
| 1,2-Dibromo-3-chloropropane | ug/L | <1.8 | 50 | 50 | 44.3 | 43.9 | 89 | 88 | 60-130 | 1 | 20 | | |
| 1,2-Dibromoethane (EDB) | ug/L | <0.83 | 50 | 50 | 52.1 | 52.0 | 104 | 104 | 70-130 | 0 | 20 | | |
| 1,2-Dichlorobenzene | ug/L | <0.71 | 50 | 50 | 52.5 | 52.3 | 105 | 105 | 70-130 | 0 | 20 | | |
| 1,2-Dichloroethane | ug/L | <0.28 | 50 | 50 | 42.7 | 43.0 | 85 | 86 | 78-145 | 1 | 20 | | |
| 1,2-Dichloropropane | ug/L | <0.28 | 50 | 50 | 46.3 | 46.8 | 93 | 94 | 86-135 | 1 | 20 | | |

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

QUALITY CONTROL DATA

Project: 47358.003 KOELLER ONE

Pace Project No.: 40210819

MATRIX SPIKE & MATRIX SPIKE DUPLICATE: 2081562 2081563

| Parameter | Units | MS | | MSD | | MS Result | MSD % Rec | MSD % Rec | % Rec Limits | Max | |
|------------------------------|-------|-------------|----------------|----------------|--------------|--------------|--------------|--------------|-----------------|-----|-----|
| | | 40210819002 | Spike Conc. | Spike Conc. | MS Result | | | | | RPD | RPD |
| 1,3-Dichlorobenzene | ug/L | <0.63 | 50 | 50 | 52.5 | 52.3 | 105 | 105 | 70-130 | 0 | 20 |
| 1,4-Dichlorobenzene | ug/L | <0.94 | 50 | 50 | 53.8 | 53.0 | 108 | 106 | 70-130 | 1 | 20 |
| Benzene | ug/L | <0.25 | 50 | 50 | 46.2 | 47.1 | 92 | 94 | 70-136 | 2 | 20 |
| Bromodichloromethane | ug/L | <0.36 | 50 | 50 | 47.2 | 47.4 | 94 | 95 | 70-130 | 1 | 20 |
| Bromoform | ug/L | <4.0 | 50 | 50 | 50.4 | 50.6 | 101 | 101 | 69-130 | 0 | 20 |
| Bromomethane | ug/L | <0.97 | 50 | 50 | 38.0 | 38.1 | 76 | 76 | 39-138 | 0 | 20 |
| Carbon tetrachloride | ug/L | <1.1 | 50 | 50 | 47.0 | 46.4 | 94 | 93 | 70-142 | 1 | 20 |
| Chlorobenzene | ug/L | <0.71 | 50 | 50 | 54.0 | 55.8 | 108 | 112 | 70-130 | 3 | 20 |
| Chloroethane | ug/L | <1.3 | 50 | 50 | 46.5 | 46.6 | 93 | 93 | 61-149 | 0 | 20 |
| Chloroform | ug/L | <1.3 | 50 | 50 | 54.9 | 48.6 | 110 | 97 | 75-133 | 12 | 20 |
| Chloromethane | ug/L | <2.2 | 50 | 50 | 37.5 | 37.1 | 75 | 74 | 32-143 | 1 | 20 |
| cis-1,2-Dichloroethene | ug/L | <0.27 | 50 | 50 | 49.3 | 50.2 | 99 | 100 | 70-130 | 2 | 20 |
| cis-1,3-Dichloropropene | ug/L | <3.6 | 50 | 50 | 47.1 | 47.3 | 94 | 95 | 70-130 | 0 | 20 |
| Dibromochloromethane | ug/L | <2.6 | 50 | 50 | 48.3 | 48.4 | 97 | 97 | 70-130 | 0 | 20 |
| Dichlorodifluoromethane | ug/L | <0.50 | 50 | 50 | 45.4 | 44.8 | 91 | 90 | 10-141 | 1 | 20 |
| Ethylbenzene | ug/L | <0.32 | 50 | 50 | 52.6 | 53.5 | 105 | 107 | 80-120 | 2 | 20 |
| Isopropylbenzene (Cumene) | ug/L | <1.7 | 50 | 50 | 54.7 | 55.1 | 109 | 110 | 70-130 | 1 | 20 |
| m&p-Xylene | ug/L | <0.47 | 100 | 100 | 110 | 110 | 110 | 110 | 70-130 | 0 | 20 |
| Methyl-tert-butyl ether | ug/L | <1.2 | 50 | 50 | 41.1 | 41.6 | 82 | 83 | 61-136 | 1 | 20 |
| Methylene Chloride | ug/L | <0.58 | 50 | 50 | 46.7 | 46.5 | 93 | 93 | 68-137 | 0 | 20 |
| o-Xylene | ug/L | <0.26 | 50 | 50 | 53.9 | 54.2 | 108 | 108 | 70-130 | 0 | 20 |
| Styrene | ug/L | <3.0 | 50 | 50 | 53.2 | 53.7 | 106 | 107 | 70-130 | 1 | 20 |
| Tetrachloroethene | ug/L | 5.6 | 50 | 50 | 56.9 | 57.6 | 103 | 104 | 70-130 | 1 | 20 |
| Toluene | ug/L | <0.27 | 50 | 50 | 52.9 | 52.8 | 106 | 106 | 80-120 | 0 | 20 |
| trans-1,2-Dichloroethene | ug/L | <0.46 | 50 | 50 | 51.8 | 51.9 | 104 | 104 | 70-130 | 0 | 20 |
| trans-1,3-Dichloropropene | ug/L | <4.4 | 50 | 50 | 45.5 | 45.4 | 91 | 91 | 69-130 | 0 | 20 |
| Trichloroethene | ug/L | <0.26 | 50 | 50 | 51.1 | 51.3 | 102 | 103 | 70-130 | 0 | 20 |
| Trichlorofluoromethane | ug/L | <0.21 | 50 | 50 | 52.9 | 53.5 | 106 | 107 | 74-157 | 1 | 20 |
| Vinyl chloride | ug/L | <0.17 | 50 | 50 | 48.1 | 48.4 | 96 | 97 | 51-140 | 0 | 20 |
| 4-Bromofluorobenzene (S) | % | | | | | | 96 | 96 | 70-130 | | |
| Dibromofluoromethane (S) | % | | | | | | 93 | 94 | 70-130 | | |
| Toluene-d8 (S) | % | | | | | | 99 | 100 | 70-130 | | |

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

QUALIFIERS

Project: 47358.003 KOELLER ONE
Pace Project No.: 40210819

DEFINITIONS

DF - Dilution Factor, if reported, represents the factor applied to the reported data due to dilution of the sample aliquot.

ND - Not Detected at or above LOD.

J - Estimated concentration at or above the LOD and below the LOQ.

LOD - Limit of Detection adjusted for dilution factor, percent moisture, initial weight and final volume.

LOQ - Limit of Quantitation adjusted for dilution factor, percent moisture, initial weight and final volume.

S - Surrogate

1,2-Diphenylhydrazine decomposes to and cannot be separated from Azobenzene using Method 8270. The result for each analyte is a combined concentration.

Consistent with EPA guidelines, unrounded data are displayed and have been used to calculate % recovery and RPD values.

LCS(D) - Laboratory Control Sample (Duplicate)

MS(D) - Matrix Spike (Duplicate)

DUP - Sample Duplicate

RPD - Relative Percent Difference

NC - Not Calculable.

SG - Silica Gel - Clean-Up

U - Indicates the compound was analyzed for, but not detected at or above the adjusted LOD.

N-Nitrosodiphenylamine decomposes and cannot be separated from Diphenylamine using Method 8270. The result reported for each analyte is a combined concentration.

Pace Analytical is TNI accredited. Contact your Pace PM for the current list of accredited analytes.

TNI - The NELAC Institute.

ANALYTE QUALIFIERS

HS Results are from sample aliquot taken from VOA vial with headspace (air bubble greater than 6 mm diameter).

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

QUALITY CONTROL DATA CROSS REFERENCE TABLE

Project: 47358.003 KOELLER ONE

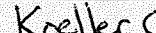
Pace Project No.: 40210819

| Lab ID | Sample ID | QC Batch Method | QC Batch | Analytical Method | Analytical Batch |
|-------------|------------|-----------------|----------|-------------------|------------------|
| 40210819001 | MW-1 | EPA 8260 | 359774 | | |
| 40210819002 | MW-2 | EPA 8260 | 359774 | | |
| 40210819003 | MW-3 | EPA 8260 | 359774 | | |
| 40210819004 | MW-4 | EPA 8260 | 359774 | | |
| 40210819005 | MW-5 | EPA 8260 | 359774 | | |
| 40210819006 | MW-6 | EPA 8260 | 359774 | | |
| 40210819007 | MW-6 DUP | EPA 8260 | 359774 | | |
| 40210819008 | MW-7 | EPA 8260 | 359774 | | |
| 40210819009 | TRIP BLANK | EPA 8260 | 359774 | | |

REPORT OF LABORATORY ANALYSIS

This report shall not be reproduced, except in full,
without the written consent of Pace Analytical Services, LLC.

(Please Print Clearly)

| | |
|---------------------|---|
| Company Name: | Gannett Fleming |
| Branch/Location: | Madison, WI |
| Project Contact: | Anthony Miller |
| Phone: | 608-327-5050 |
| Project Number: | 47358.003 |
| Project Name: | Koeller One |
| Project State: | WI |
| Sampled By (Print): | Chelsea Payne |
| Sampled By (Sign): |  |
| PO #: | |
| | Regulatory Program: |



UPPER MIDWEST REGION

MN: 612-607-1700 WI: 920-469-2436

Page 1 of

Page 30 of 32

CHAIN OF CUSTODY

***Preservation Codes**

| | | | | | | |
|-----------------------------|----------------------|----------------------------------|--------------------|------------|------------|--------|
| A=None | B=HCl | C=H ₂ SO ₄ | D=HNO ₃ | E=DI Water | F=Methanol | G=NaOH |
| H=Sodium Bisulfate Solution | I=Sodium Thiosulfate | J=Other | | | | |

| | | | | | |
|--|----------------------------------|----------------------------|---|----------------------------|-----------------------------|
| Rush Turnaround Time Requested - Prelims (Rush TAT subject to approval/surcharge) | Relinquished By: | Date/Time: | Received By: | Date/Time: | PACE Project No. |
| Date Needed: | <i>Chamberlain</i> | 7/18/20 13:30 | | | 40210219 |
| Transmit Prelim Rush Results by (complete what you want): | Relinquished By: <i>FedEx</i> | Date/Time: 7/19/20 0940 | Received By: <i>Eric Miller/Pace</i> | Date/Time: 7/19/20 0940 | Receipt Temp = <i>RT</i> °C |
| Email #1: | Relinquished By: | Date/Time: | Received By: | Date/Time: | Sample Receipt pH |
| Email #2: | | | | | OK / Adjusted |
| Telephone: | Relinquished By: | Date/Time: | Received By: | Date/Time: | Cooler Custody Seal |
| Fax: | | | | | Present / Not Present |
| Samples on HOLD are subject to special pricing and release of liability | Relinquished By: | Date/Time: | Received By: | Date/Time: | Intact / Not Intact |

Sample Preservation Receipt Form

Client Name: Gannett Fleming

All containers needing preservation have been checked and noted below: Yes No N/A

Project # 40210879

Pace Analytical Services, LLC
1241 Bellevue Street, Suite 9
Green Bay, WI 54302

Page 31 of 32

| Pace Lab # | Glass | | | | | Plastic | | | | | Vials | | | | | Jars | | | General | | | Initial when completed: | Date/ Time: | | | | |
|------------|-------|------|------|------|------|---------|------|------|------|------|-------|------|------|------|------|------|------|------|---------|------|------|-------------------------|----------------|------|------|--------------|--|
| | AG1U | BG1U | AG1H | AG4S | AG4U | AG5U | AG2S | BG3U | BP1U | BP3U | BP3B | BP3N | BP3S | VG9A | DG9T | VG9U | VG9H | VG9M | VG9D | JGFU | JG9U | WGFU | WPFU | SP5T | ZPLC | GN | |
| 001 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 002 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 003 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 004 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 005 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 006 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 007 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 008 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 009 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 010 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 011 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 012 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 013 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 014 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 015 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 016 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 017 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 018 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 019 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |
| 020 | | | | | | | | | | | | | | | | | | | | | | | | | | 2.5 / 5 / 10 | |

Exceptions to preservation check: VOA, Coliform, TOC, TOX, TOH, O&G, WI DRO, Phenolics, Other:

Headspace in VOA Vials (>6mm) : Yes No N/A *If yes look in headspace column

| | | | | | | | |
|------|---------------------------|------|------------------------|------|-------------------------|------|-------------------------------|
| AG1U | 1 liter amber glass | BP1U | 1 liter plastic unpres | VG9A | 40 mL clear ascorbic | JGFU | 4 oz amber jar unpres |
| BG1U | 1 liter clear glass | BP3U | 250 mL plastic unpres | DG9T | 40 mL amber Na Thio | JG9U | 9 oz amber jar unpres |
| AG1H | 1 liter amber glass HCL | BP3B | 250 mL plastic NaOH | VG9U | 40 mL clear vial unpres | WGFU | 4 oz clear jar unpres |
| AG4S | 125 mL amber glass H2SO4 | BP3N | 250 mL plastic HNO3 | VG9H | 40 mL clear vial HCL | WPFU | 4 oz plastic jar unpres |
| AG4U | 120 mL amber glass unpres | BP3S | 250 mL plastic H2SO4 | VG9M | 40 mL clear vial MeOH | SP5T | 120 mL plastic Na Thiosulfate |
| AG5U | 100 mL amber glass unpres | | | VG9D | 40 mL clear vial DI | ZPLC | ziploc bag |
| AG2S | 500 mL amber glass H2SO4 | | | | | GN | |
| BG3U | 250 mL clear glass unpres | | | | | | |



Document Name:
Sample Condition Upon Receipt (SCUR)
Document No.:
ENV-FRM-GBAY-0014-Rev.00

Document Revised: 26Mar2020
Author:
Pace Green Bay Quality Office

Sample Condition Upon Receipt Form (SCUR)

Project #:

WO# : 40210819

Client Name: Gannett Fleming

Courier: CS Logistics Fed Ex Speedee UPS Waltco
 Client Pace Other:

Tracking #: 803262940323

Custody Seal on Cooler/Box Present: yes no Seals intact: yes no

Custody Seal on Samples Present: yes no Seals intact: yes no

Packing Material: Bubble Wrap Bubble Bags None Other

Thermometer Used SR - M4 Type of Ice: Wet Blue Dry None

Cooler Temperature Uncorr: 20.1 /Corr:



40210819

Temp Blank Present: yes no

Biological Tissue is Frozen: yes no

Temp should be above freezing to 6°C.

Biota Samples may be received at ≤ 0°C if shipped on Dry Ice.

Samples on ice, cooling process has begun

Person examining contents:

Date: 7/9/2019 Initials: EMW

Labeled By Initials: MJ

| | | |
|--|--|------------------|
| Chain of Custody Present: | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 1. |
| Chain of Custody Filled Out: | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 2. |
| Chain of Custody Relinquished: | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 3. |
| Sampler Name & Signature on COC: | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 4. |
| Samples Arrived within Hold Time: - VOA Samples frozen upon receipt | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No | 5. Date/Time: |
| Short Hold Time Analysis (<72hr): | <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No | 6. |
| Rush Turn Around Time Requested: | <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No | 7. |
| Sufficient Volume: | | 8. |
| For Analysis: <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No MS/MSD: <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A | | |
| Correct Containers Used: -Pace Containers Used: -Pace IR Containers Used: | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No | 9. |
| Containers Intact: | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No | 10. |
| Filtered volume received for Dissolved tests | <input type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 11. |
| Sample Labels match COC: -Includes date/time/ID/Analysis Matrix: <u>WS</u> | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 12. |
| Trip Blank Present: | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 13. |
| Trip Blank Custody Seals Present | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | |
| Pace Trip Blank Lot # (if purchased): <u>411</u> | | |

Client Notification/ Resolution:

If checked, see attached form for additional comments

Person Contacted: _____ Date/Time: _____

Comments/ Resolution:

PM Review is documented electronically in LIMs. By releasing the project, the PM acknowledges they have reviewed the sample log.



The laboratory report and
QA/QC data were reviewed
and approved by AWM on
08/31/21

31-Aug-2021

Anthony Miller
Gannett Fleming, Inc.
8040 Excelsior Drive
Suite 303
Madison, WI 53717-1338

Re: **Koeller (47358.003)**

Work Order: **21081889**

Dear Anthony,

ALS Environmental received 10 samples on 20-Aug-2021 08:30 AM for the analyses presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental - Holland and for only the analyses requested.

Sample results are compliant with industry accepted practices and Quality Control results achieved laboratory specifications. Any exceptions are noted in the Case Narrative, or noted with qualifiers in the report or QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained from ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

The total number of pages in this report is 53.

If you have any questions regarding this report, please feel free to contact me:

ADDRESS: 3352 128th Avenue, Holland, MI, USA
PHONE: +1 (616) 399-6070 FAX: +1 (616) 399-6185

Sincerely,

A handwritten signature in black ink that reads "Jodi Blouw".

Electronically approved by: Jodi Blouw

Jodi Blouw

Report of Laboratory Analysis

Certificate No: MN 026-999-449

ALS GROUP USA, CORP Part of the ALS Laboratory Group A Campbell Brothers Limited Company

Environmental

www.alsglobal.com

RIGHT SOLUTIONS RIGHT PARTNER

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Work Order: **21081889**

Work Order Sample Summary

| Lab Samp ID | Client Sample ID | Matrix | Tag Number | Collection Date | Date Received | Hold |
|--------------------|-------------------------|---------------|-------------------|------------------------|----------------------|--------------------------|
| 21081889-01 | MW-1 | Water | | 8/18/2021 14:00 | 8/20/2021 08:30 | <input type="checkbox"/> |
| 21081889-02 | MW-2 | Water | | 8/18/2021 12:55 | 8/20/2021 08:30 | <input type="checkbox"/> |
| 21081889-03 | MW-3 | Water | | 8/18/2021 13:25 | 8/20/2021 08:30 | <input type="checkbox"/> |
| 21081889-04 | MW-4 | Water | | 8/18/2021 11:40 | 8/20/2021 08:30 | <input type="checkbox"/> |
| 21081889-05 | MW-5 | Water | | 8/18/2021 10:25 | 8/20/2021 08:30 | <input type="checkbox"/> |
| 21081889-06 | MW-6 | Water | | 8/18/2021 12:10 | 8/20/2021 08:30 | <input type="checkbox"/> |
| 21081889-07 | MW-7 | Water | | 8/18/2021 11:05 | 8/20/2021 08:30 | <input type="checkbox"/> |
| 21081889-08 | MW-1 DUP | Water | | 8/18/2021 14:05 | 8/20/2021 08:30 | <input type="checkbox"/> |
| 21081889-09 | Field Blank | Water | | 8/18/2021 | 8/20/2021 08:30 | <input type="checkbox"/> |
| 21081889-10 | Trip Blank | Water | | 8/18/2021 | 8/20/2021 08:30 | <input type="checkbox"/> |

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
WorkOrder: 21081889

**QUALIFIERS,
ACRONYMS, UNITS**

| <u>Qualifier</u> | <u>Description</u> |
|-------------------------|---|
| * | Value exceeds Regulatory Limit |
| ** | Estimated Value |
| a | Analyte is non-accredited |
| B | Analyte detected in the associated Method Blank above the Reporting Limit |
| E | Value above quantitation range |
| H | Analyzed outside of Holding Time |
| Hr | BOD/CBOD - Sample was reset outside Hold Time, value should be considered estimated. |
| J | Analyte is present at an estimated concentration between the MDL and Report Limit |
| ND | Not Detected at the Reporting Limit |
| O | Sample amount is > 4 times amount spiked |
| P | Dual Column results percent difference > 40% |
| R | RPD above laboratory control limit |
| S | Spike Recovery outside laboratory control limits |
| U | Analyzed but not detected above the MDL |
| X | Analyte was detected in the Method Blank between the MDL and Reporting Limit, sample results may exhibit background or reagent contamination at the observed level. |

| <u>Acronym</u> | <u>Description</u> |
|-----------------------|-------------------------------------|
| DUP | Method Duplicate |
| LCS | Laboratory Control Sample |
| LCSD | Laboratory Control Sample Duplicate |
| LOD | Limit of Detection (see MDL) |
| LOQ | Limit of Quantitation (see PQL) |
| MBLK | Method Blank |
| MDL | Method Detection Limit |
| MS | Matrix Spike |
| MSD | Matrix Spike Duplicate |
| PQL | Practical Quantitation Limit |
| RPD | Relative Percent Difference |
| TDL | Target Detection Limit |
| TNTC | Too Numerous To Count |
| A | APHA Standard Methods |
| D | ASTM |
| E | EPA |
| SW | SW-846 Update III |

| <u>Units Reported</u> | <u>Description</u> |
|------------------------------|---------------------------|
| µg/L | Micrograms per Liter |
| ng/L | Nanograms per Liter |

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Work Order: 21081889

Case Narrative

Samples for the above noted Work Order were received on 08/20/2021. The attached "Sample Receipt Checklist" documents the status of custody seals, container integrity, preservation, and temperature compliance.

Samples were analyzed according to the analytical methodology previously transmitted in the "Work Order Acknowledgement". Methodologies are also documented in the "Analytical Result" section for each sample. Quality control results are listed in the "QC Report" section. Sample association for the reported quality control is located at the end of each batch summary. If applicable, results are appropriately qualified in the Analytical Result and QC Report sections. The "Qualifiers" section documents the various qualifiers, units, and acronyms utilized in reporting. A copy of the laboratory's scope of accreditation is available upon request.

With the following exceptions, all sample analyses achieved analytical criteria.

Volatile Organics:

No deviations or anomalies were noted.

Extractable Organics:

Batch 182387, Method E537 Mod, Sample LCS-182387: The LCS recovery was above the upper control limit. All the sample results in the batch were non-detect. No qualification is necessary for this analyte: FtS 10:2

No other deviations or anomalies were noted.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-1
Collection Date: 8/18/2021 02:00 PM

Work Order: 21081889
Lab ID: 21081889-01
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|--|-------------|------|-------------|--------------|-------|-----------------|-----------------|
| PFAS BY EPA 537 MODIFIED | | | | | | | |
| Fluorotelomer Sulphonic Acid 4:2 (FtS 4:2) | U | | 0.93 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Fluorotelomer Sulphonic Acid 6:2 (FtS 6:2) | U | | 0.66 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Fluorotelomer Sulphonic Acid 8:2 (FtS 8:2) | U | | 1.1 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Fluorotelomer Sulphonic Acid 10:2 (FtS 10:2) | U | | 0.89 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorobutanesulfonic Acid (PFBS) | 3.2 | J | 0.35 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorobutanoic Acid (PFBA) | 5.1 | | 2.6 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorodecanesulfonic Acid (PFDS) | U | | 1.4 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorodecanoic Acid (PFDA) | U | | 1.2 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorododecanesulfonic Acid (PFDoS) | U | | 1.4 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorododecanoic Acid (PFDoA) | U | | 1.4 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluoroheptanesulfonic Acid (PFHpS) | U | | 0.56 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluoroheptanoic Acid (PFHpA) | 2.3 | J | 0.43 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorohexadecanoic Acid (PFHxDA) | U | | 0.38 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorohexanesulfonic Acid (PFHxS) | U | | 0.36 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorohexanoic Acid (PFHxA) | 2.3 | J | 1.2 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorononanesulfonic Acid (PFNS) | U | | 0.49 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorononanoic Acid (PFNA) | 0.95 | J | 0.86 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorooctadecanoic Acid (PFODA) | U | | 0.64 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorooctanesulfonamide (PFOSA) | U | | 0.70 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorooctanesulfonic Acid (PFOS) | 9.7 | | 0.88 | 2.0 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorooctanoic Acid (PFOA) | 5.8 | | 0.62 | 2.0 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluoropentanesulfonic Acid (PFPeS) | U | | 0.55 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluoropentanoic Acid (PFPeA) | 2.6 | J | 1.3 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorotetradecanoic Acid (PFTeA) | U | | 2.6 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluorotridecanoic Acid (PFTriA) | U | | 0.76 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Perfluoroundecanoic Acid (PFUnA) | U | | 0.96 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| N-ethylperfluoro-1-octanesulfonamide | U | | 1.1 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| N-Ethylperfluorooctanesulfonamidoacetic Acid | U | | 0.62 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| N-Ethylperfluorooctanesulfonamidoethanol | U | | 0.51 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| N-methylperfluoro-1-octanesulfonamide | U | | 0.78 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-1
Collection Date: 8/18/2021 02:00 PM

Work Order: 21081889
Lab ID: 21081889-01
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|--|--------|------------------------|------|--------------|-------|--------------------|-----------------|
| N-Methylperfluorooctanesulfonamidoacetic Acid | U | | 0.64 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| N-Methylperfluorooctanesulfonamidoethanol | U | | 0.48 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Hexafluoropropylene oxide dimer acid (HFPO-DA) | U | | 1.2 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| 4,8-Dioxa-3H-perfluorononanoic Acid (DONA) | U | | 0.56 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| 11Cl-Pf3OUDS | U | | 0.46 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| 9Cl-PF3ONS | U | | 0.44 | 4.9 | ng/L | 1 | 8/23/2021 20:54 |
| Surr: 13C2-FtS 4:2 | 120 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C2-FtS 6:2 | 104 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C2-FtS 8:2 | 99.0 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C2-PFDA | 86.3 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C2-PFDoA | 66.9 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C2-PFHxA | 82.3 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C2-PFHxDA | 75.3 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C2-PFTeA | 72.4 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C2-PFUra | 85.0 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C3-HFPO-DA | 72.6 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C3-PFBS | 78.7 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C4-PFBA | 80.2 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C4-PFHpa | 70.6 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C4-PFOA | 79.8 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C4-PFOS | 83.6 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C5-PFNA | 79.1 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C5-PFPeA | 80.1 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 13C8-FOSA | 71.3 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: 18O2-PFHxS | 87.1 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: d5-N-EtFOSA | 64.1 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: d5-N-EtFOSAA | 94.8 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: d9-N-EtFOSE | 66.7 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: d3-N-MeFOSA | 67.4 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: d3-N-MeFOSAA | 87.2 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| Surr: d7-N-MeFOSE | 76.6 | | | 50-150 | %REC | 1 | 8/23/2021 20:54 |
| 1,4-DIOXANE BY SELECT ION MONITORING | | Method: SW8260B | | | | Analyst: BG | |
| 1,4-Dioxane | U | | 0.44 | 1.0 | µg/L | 1 | 8/30/2021 17:31 |
| Surr: Toluene-d8 | 99.4 | | | 74-124 | %REC | 1 | 8/30/2021 17:31 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-1
Collection Date: 8/18/2021 02:00 PM

Work Order: 21081889
Lab ID: 21081889-01
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------------|--------|------|------|-----------------|-------|-----------------|-----------------|
| VOLATILE ORGANIC COMPOUNDS | | | | | | | |
| | | | | Method: SW8260C | | | Analyst: MF |
| 1,1,1,2-Tetrachloroethane | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 02:08 |
| 1,1,1-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 02:08 |
| 1,1,2,2-Tetrachloroethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 02:08 |
| 1,1,2-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 02:08 |
| 1,1-Dichloroethane | U | | 0.44 | 1.5 | µg/L | 1 | 8/27/2021 02:08 |
| 1,1-Dichloroethene | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 02:08 |
| 1,1-Dichloropropene | U | | 0.37 | 1.2 | µg/L | 1 | 8/27/2021 02:08 |
| 1,2,3-Trichlorobenzene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 02:08 |
| 1,2,3-Trichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 02:08 |
| 1,2,4-Trichlorobenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 02:08 |
| 1,2,4-Trimethylbenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 02:08 |
| 1,2-Dibromo-3-chloropropane | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 02:08 |
| 1,2-Dibromoethane | U | | 0.41 | 1.4 | µg/L | 1 | 8/27/2021 02:08 |
| 1,2-Dichlorobenzene | U | | 0.32 | 1.1 | µg/L | 1 | 8/27/2021 02:08 |
| 1,2-Dichloroethane | U | | 0.44 | 1.4 | µg/L | 1 | 8/27/2021 02:08 |
| 1,2-Dichloropropane | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 02:08 |
| 1,3,5-Trimethylbenzene | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 02:08 |
| 1,3-Dichlorobenzene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 02:08 |
| 1,3-Dichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 02:08 |
| 1,4-Dichlorobenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 02:08 |
| 2,2-Dichloropropane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 02:08 |
| 2-Butanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 02:08 |
| 2-Chlorotoluene | U | | 0.36 | 1.2 | µg/L | 1 | 8/27/2021 02:08 |
| 4-Chlorotoluene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 02:08 |
| 4-Methyl-2-pentanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 02:08 |
| Acetone | U | | 6.2 | 21 | µg/L | 1 | 8/27/2021 02:08 |
| Benzene | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 02:08 |
| Bromobenzene | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 02:08 |
| Bromochloromethane | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 02:08 |
| Bromodichloromethane | U | | 0.49 | 1.6 | µg/L | 1 | 8/27/2021 02:08 |
| Bromoform | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 02:08 |
| Bromomethane | U | | 0.90 | 3.0 | µg/L | 1 | 8/27/2021 02:08 |
| Carbon tetrachloride | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 02:08 |
| Chlorobenzene | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 02:08 |
| Chloroethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 02:08 |
| Chloroform | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 02:08 |
| Chloromethane | U | | 0.83 | 2.8 | µg/L | 1 | 8/27/2021 02:08 |
| cis-1,2-Dichloroethene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 02:08 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-1
Collection Date: 8/18/2021 02:00 PM

Work Order: 21081889
Lab ID: 21081889-01
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------|------------|------|------------|--------------|-------------|-----------------|-----------------|
| cis-1,3-Dichloropropene | U | | 0.57 | 1.9 | µg/L | 1 | 8/27/2021 02:08 |
| Dibromochloromethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 02:08 |
| Dibromomethane | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 02:08 |
| Dichlorodifluoromethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 02:08 |
| Ethylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 02:08 |
| Hexachlorobutadiene | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 02:08 |
| Isopropylbenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 02:08 |
| m,p-Xylene | U | | 0.81 | 2.7 | µg/L | 1 | 8/27/2021 02:08 |
| Methyl tert-butyl ether | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 02:08 |
| Methylene chloride | U | | 0.86 | 2.9 | µg/L | 1 | 8/27/2021 02:08 |
| Naphthalene | U | | 0.77 | 2.6 | µg/L | 1 | 8/27/2021 02:08 |
| n-Butylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 02:08 |
| n-Propylbenzene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 02:08 |
| o-Xylene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 02:08 |
| p-Isopropyltoluene | U | | 0.26 | 0.88 | µg/L | 1 | 8/27/2021 02:08 |
| sec-Butylbenzene | U | | 0.30 | 1.0 | µg/L | 1 | 8/27/2021 02:08 |
| Styrene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 02:08 |
| tert-Butylbenzene | U | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 02:08 |
| Tetrachloroethene | 120 | | 2.0 | 6.6 | µg/L | 5 | 8/27/2021 15:31 |
| Toluene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 02:08 |
| trans-1,2-Dichloroethene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 02:08 |
| trans-1,3-Dichloropropene | U | | 0.38 | 2.7 | µg/L | 1 | 8/27/2021 02:08 |
| Trichloroethene | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 02:08 |
| Trichlorofluoromethane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 02:08 |
| Vinyl chloride | U | | 0.53 | 1.8 | µg/L | 1 | 8/27/2021 02:08 |
| Xylenes, Total | U | | 0.81 | 4.4 | µg/L | 1 | 8/27/2021 02:08 |
| Surr: 1,2-Dichloroethane-d4 | 103 | | | 75-120 | %REC | 1 | 8/27/2021 02:08 |
| Surr: 1,2-Dichloroethane-d4 | 99.0 | | | 75-120 | %REC | 5 | 8/27/2021 15:31 |
| Surr: 4-Bromofluorobenzene | 95.5 | | | 80-110 | %REC | 1 | 8/27/2021 02:08 |
| Surr: 4-Bromofluorobenzene | 101 | | | 80-110 | %REC | 5 | 8/27/2021 15:31 |
| Surr: Dibromofluoromethane | 102 | | | 85-115 | %REC | 1 | 8/27/2021 02:08 |
| Surr: Dibromofluoromethane | 99.1 | | | 85-115 | %REC | 5 | 8/27/2021 15:31 |
| Surr: Toluene-d8 | 104 | | | 85-110 | %REC | 1 | 8/27/2021 02:08 |
| Surr: Toluene-d8 | 100 | | | 85-110 | %REC | 5 | 8/27/2021 15:31 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-2
Collection Date: 8/18/2021 12:55 PM

Work Order: 21081889
Lab ID: 21081889-02
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------------|--------|------|------|-----------------|-------|-----------------|-----------------|
| VOLATILE ORGANIC COMPOUNDS | | | | | | | |
| | | | | Method: SW8260C | | | Analyst: MF |
| 1,1,1,2-Tetrachloroethane | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 15:51 |
| 1,1,1-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 15:51 |
| 1,1,2,2-Tetrachloroethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 15:51 |
| 1,1,2-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 15:51 |
| 1,1-Dichloroethane | U | | 0.44 | 1.5 | µg/L | 1 | 8/27/2021 15:51 |
| 1,1-Dichloroethene | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 15:51 |
| 1,1-Dichloropropene | U | | 0.37 | 1.2 | µg/L | 1 | 8/27/2021 15:51 |
| 1,2,3-Trichlorobenzene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 15:51 |
| 1,2,3-Trichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 15:51 |
| 1,2,4-Trichlorobenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 15:51 |
| 1,2,4-Trimethylbenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 15:51 |
| 1,2-Dibromo-3-chloropropane | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 15:51 |
| 1,2-Dibromoethane | U | | 0.41 | 1.4 | µg/L | 1 | 8/27/2021 15:51 |
| 1,2-Dichlorobenzene | U | | 0.32 | 1.1 | µg/L | 1 | 8/27/2021 15:51 |
| 1,2-Dichloroethane | U | | 0.44 | 1.4 | µg/L | 1 | 8/27/2021 15:51 |
| 1,2-Dichloropropane | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 15:51 |
| 1,3,5-Trimethylbenzene | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 15:51 |
| 1,3-Dichlorobenzene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 15:51 |
| 1,3-Dichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 15:51 |
| 1,4-Dichlorobenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 15:51 |
| 2,2-Dichloropropane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 15:51 |
| 2-Butanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 15:51 |
| 2-Chlorotoluene | U | | 0.36 | 1.2 | µg/L | 1 | 8/27/2021 15:51 |
| 4-Chlorotoluene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 15:51 |
| 4-Methyl-2-pentanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 15:51 |
| Acetone | U | | 6.2 | 21 | µg/L | 1 | 8/27/2021 15:51 |
| Benzene | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 15:51 |
| Bromobenzene | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 15:51 |
| Bromochloromethane | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 15:51 |
| Bromodichloromethane | U | | 0.49 | 1.6 | µg/L | 1 | 8/27/2021 15:51 |
| Bromoform | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 15:51 |
| Bromomethane | U | | 0.90 | 3.0 | µg/L | 1 | 8/27/2021 15:51 |
| Carbon tetrachloride | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 15:51 |
| Chlorobenzene | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 15:51 |
| Chloroethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 15:51 |
| Chloroform | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 15:51 |
| Chloromethane | U | | 0.83 | 2.8 | µg/L | 1 | 8/27/2021 15:51 |
| cis-1,2-Dichloroethene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 15:51 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-2
Collection Date: 8/18/2021 12:55 PM

Work Order: 21081889
Lab ID: 21081889-02
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------|------------|------|-------------|--------------|-------------|-----------------|-----------------|
| cis-1,3-Dichloropropene | U | | 0.57 | 1.9 | µg/L | 1 | 8/27/2021 15:51 |
| Dibromochloromethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 15:51 |
| Dibromomethane | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 15:51 |
| Dichlorodifluoromethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 15:51 |
| Ethylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 15:51 |
| Hexachlorobutadiene | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 15:51 |
| Isopropylbenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 15:51 |
| m,p-Xylene | U | | 0.81 | 2.7 | µg/L | 1 | 8/27/2021 15:51 |
| Methyl tert-butyl ether | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 15:51 |
| Methylene chloride | U | | 0.86 | 2.9 | µg/L | 1 | 8/27/2021 15:51 |
| Naphthalene | U | | 0.77 | 2.6 | µg/L | 1 | 8/27/2021 15:51 |
| n-Butylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 15:51 |
| n-Propylbenzene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 15:51 |
| o-Xylene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 15:51 |
| p-Isopropyltoluene | U | | 0.26 | 0.88 | µg/L | 1 | 8/27/2021 15:51 |
| sec-Butylbenzene | U | | 0.30 | 1.0 | µg/L | 1 | 8/27/2021 15:51 |
| Styrene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 15:51 |
| tert-Butylbenzene | U | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 15:51 |
| Tetrachloroethene | 2.7 | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 15:51 |
| Toluene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 15:51 |
| trans-1,2-Dichloroethene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 15:51 |
| trans-1,3-Dichloropropene | U | | 0.38 | 2.7 | µg/L | 1 | 8/27/2021 15:51 |
| Trichloroethene | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 15:51 |
| Trichlorofluoromethane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 15:51 |
| Vinyl chloride | U | | 0.53 | 1.8 | µg/L | 1 | 8/27/2021 15:51 |
| Xylenes, Total | U | | 0.81 | 4.4 | µg/L | 1 | 8/27/2021 15:51 |
| Surr: 1,2-Dichloroethane-d4 | 101 | | | 75-120 | %REC | 1 | 8/27/2021 15:51 |
| Surr: 4-Bromofluorobenzene | 100 | | | 80-110 | %REC | 1 | 8/27/2021 15:51 |
| Surr: Dibromofluoromethane | 104 | | | 85-115 | %REC | 1 | 8/27/2021 15:51 |
| Surr: Toluene-d8 | 101 | | | 85-110 | %REC | 1 | 8/27/2021 15:51 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-3
Collection Date: 8/18/2021 01:25 PM

Work Order: 21081889
Lab ID: 21081889-03
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------------|--------|------|------|-----------------|-------|-----------------|-----------------|
| VOLATILE ORGANIC COMPOUNDS | | | | | | | |
| | | | | Method: SW8260C | | | Analyst: MF |
| 1,1,1,2-Tetrachloroethane | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 02:48 |
| 1,1,1-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 02:48 |
| 1,1,2,2-Tetrachloroethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 02:48 |
| 1,1,2-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 02:48 |
| 1,1-Dichloroethane | U | | 0.44 | 1.5 | µg/L | 1 | 8/27/2021 02:48 |
| 1,1-Dichloroethene | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 02:48 |
| 1,1-Dichloropropene | U | | 0.37 | 1.2 | µg/L | 1 | 8/27/2021 02:48 |
| 1,2,3-Trichlorobenzene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 02:48 |
| 1,2,3-Trichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 02:48 |
| 1,2,4-Trichlorobenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 02:48 |
| 1,2,4-Trimethylbenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 02:48 |
| 1,2-Dibromo-3-chloropropane | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 02:48 |
| 1,2-Dibromoethane | U | | 0.41 | 1.4 | µg/L | 1 | 8/27/2021 02:48 |
| 1,2-Dichlorobenzene | U | | 0.32 | 1.1 | µg/L | 1 | 8/27/2021 02:48 |
| 1,2-Dichloroethane | U | | 0.44 | 1.4 | µg/L | 1 | 8/27/2021 02:48 |
| 1,2-Dichloropropane | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 02:48 |
| 1,3,5-Trimethylbenzene | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 02:48 |
| 1,3-Dichlorobenzene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 02:48 |
| 1,3-Dichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 02:48 |
| 1,4-Dichlorobenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 02:48 |
| 2,2-Dichloropropane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 02:48 |
| 2-Butanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 02:48 |
| 2-Chlorotoluene | U | | 0.36 | 1.2 | µg/L | 1 | 8/27/2021 02:48 |
| 4-Chlorotoluene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 02:48 |
| 4-Methyl-2-pentanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 02:48 |
| Acetone | U | | 6.2 | 21 | µg/L | 1 | 8/27/2021 02:48 |
| Benzene | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 02:48 |
| Bromobenzene | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 02:48 |
| Bromochloromethane | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 02:48 |
| Bromodichloromethane | U | | 0.49 | 1.6 | µg/L | 1 | 8/27/2021 02:48 |
| Bromoform | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 02:48 |
| Bromomethane | U | | 0.90 | 3.0 | µg/L | 1 | 8/27/2021 02:48 |
| Carbon tetrachloride | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 02:48 |
| Chlorobenzene | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 02:48 |
| Chloroethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 02:48 |
| Chloroform | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 02:48 |
| Chloromethane | U | | 0.83 | 2.8 | µg/L | 1 | 8/27/2021 02:48 |
| cis-1,2-Dichloroethene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 02:48 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-3
Collection Date: 8/18/2021 01:25 PM

Work Order: 21081889
Lab ID: 21081889-03
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------|------------|------|-------------|--------------|-------------|-----------------|------------------------|
| cis-1,3-Dichloropropene | U | | 0.57 | 1.9 | µg/L | 1 | 8/27/2021 02:48 |
| Dibromochloromethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 02:48 |
| Dibromomethane | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 02:48 |
| Dichlorodifluoromethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 02:48 |
| Ethylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 02:48 |
| Hexachlorobutadiene | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 02:48 |
| Isopropylbenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 02:48 |
| m,p-Xylene | U | | 0.81 | 2.7 | µg/L | 1 | 8/27/2021 02:48 |
| Methyl tert-butyl ether | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 02:48 |
| Methylene chloride | U | | 0.86 | 2.9 | µg/L | 1 | 8/27/2021 02:48 |
| Naphthalene | U | | 0.77 | 2.6 | µg/L | 1 | 8/27/2021 02:48 |
| n-Butylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 02:48 |
| n-Propylbenzene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 02:48 |
| o-Xylene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 02:48 |
| p-Isopropyltoluene | U | | 0.26 | 0.88 | µg/L | 1 | 8/27/2021 02:48 |
| sec-Butylbenzene | U | | 0.30 | 1.0 | µg/L | 1 | 8/27/2021 02:48 |
| Styrene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 02:48 |
| tert-Butylbenzene | U | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 02:48 |
| Tetrachloroethene | 4.2 | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 02:48 |
| Toluene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 02:48 |
| trans-1,2-Dichloroethene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 02:48 |
| trans-1,3-Dichloropropene | U | | 0.38 | 2.7 | µg/L | 1 | 8/27/2021 02:48 |
| Trichloroethene | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 02:48 |
| Trichlorofluoromethane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 02:48 |
| Vinyl chloride | U | | 0.53 | 1.8 | µg/L | 1 | 8/27/2021 02:48 |
| Xylenes, Total | U | | 0.81 | 4.4 | µg/L | 1 | 8/27/2021 02:48 |
| Surr: 1,2-Dichloroethane-d4 | 102 | | | 75-120 | %REC | 1 | 8/27/2021 02:48 |
| Surr: 4-Bromofluorobenzene | 95.0 | | | 80-110 | %REC | 1 | 8/27/2021 02:48 |
| Surr: Dibromofluoromethane | 103 | | | 85-115 | %REC | 1 | 8/27/2021 02:48 |
| Surr: Toluene-d8 | 100 | | | 85-110 | %REC | 1 | 8/27/2021 02:48 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-4
Collection Date: 8/18/2021 11:40 AM

Work Order: 21081889
Lab ID: 21081889-04
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------------|--------|------|------|-----------------|-------|-----------------|-----------------|
| VOLATILE ORGANIC COMPOUNDS | | | | | | | |
| | | | | Method: SW8260C | | | Analyst: MF |
| 1,1,1,2-Tetrachloroethane | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 03:07 |
| 1,1,1-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 03:07 |
| 1,1,2,2-Tetrachloroethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:07 |
| 1,1,2-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 03:07 |
| 1,1-Dichloroethane | U | | 0.44 | 1.5 | µg/L | 1 | 8/27/2021 03:07 |
| 1,1-Dichloroethene | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 03:07 |
| 1,1-Dichloropropene | U | | 0.37 | 1.2 | µg/L | 1 | 8/27/2021 03:07 |
| 1,2,3-Trichlorobenzene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 03:07 |
| 1,2,3-Trichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:07 |
| 1,2,4-Trichlorobenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:07 |
| 1,2,4-Trimethylbenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:07 |
| 1,2-Dibromo-3-chloropropane | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 03:07 |
| 1,2-Dibromoethane | U | | 0.41 | 1.4 | µg/L | 1 | 8/27/2021 03:07 |
| 1,2-Dichlorobenzene | U | | 0.32 | 1.1 | µg/L | 1 | 8/27/2021 03:07 |
| 1,2-Dichloroethane | U | | 0.44 | 1.4 | µg/L | 1 | 8/27/2021 03:07 |
| 1,2-Dichloropropane | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 03:07 |
| 1,3,5-Trimethylbenzene | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 03:07 |
| 1,3-Dichlorobenzene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 03:07 |
| 1,3-Dichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:07 |
| 1,4-Dichlorobenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 03:07 |
| 2,2-Dichloropropane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 03:07 |
| 2-Butanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 03:07 |
| 2-Chlorotoluene | U | | 0.36 | 1.2 | µg/L | 1 | 8/27/2021 03:07 |
| 4-Chlorotoluene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 03:07 |
| 4-Methyl-2-pentanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 03:07 |
| Acetone | U | | 6.2 | 21 | µg/L | 1 | 8/27/2021 03:07 |
| Benzene | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 03:07 |
| Bromobenzene | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 03:07 |
| Bromochloromethane | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:07 |
| Bromodichloromethane | U | | 0.49 | 1.6 | µg/L | 1 | 8/27/2021 03:07 |
| Bromoform | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 03:07 |
| Bromomethane | U | | 0.90 | 3.0 | µg/L | 1 | 8/27/2021 03:07 |
| Carbon tetrachloride | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 03:07 |
| Chlorobenzene | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:07 |
| Chloroethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 03:07 |
| Chloroform | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 03:07 |
| Chloromethane | U | | 0.83 | 2.8 | µg/L | 1 | 8/27/2021 03:07 |
| cis-1,2-Dichloroethene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 03:07 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-4
Collection Date: 8/18/2021 11:40 AM

Work Order: 21081889
Lab ID: 21081889-04
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------|------------|------|-------------|--------------|-------------|-----------------|------------------------|
| cis-1,3-Dichloropropene | U | | 0.57 | 1.9 | µg/L | 1 | 8/27/2021 03:07 |
| Dibromochloromethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:07 |
| Dibromomethane | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 03:07 |
| Dichlorodifluoromethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 03:07 |
| Ethylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 03:07 |
| Hexachlorobutadiene | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 03:07 |
| Isopropylbenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 03:07 |
| m,p-Xylene | U | | 0.81 | 2.7 | µg/L | 1 | 8/27/2021 03:07 |
| Methyl tert-butyl ether | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:07 |
| Methylene chloride | U | | 0.86 | 2.9 | µg/L | 1 | 8/27/2021 03:07 |
| Naphthalene | U | | 0.77 | 2.6 | µg/L | 1 | 8/27/2021 03:07 |
| n-Butylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 03:07 |
| n-Propylbenzene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 03:07 |
| o-Xylene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 03:07 |
| p-Isopropyltoluene | U | | 0.26 | 0.88 | µg/L | 1 | 8/27/2021 03:07 |
| sec-Butylbenzene | U | | 0.30 | 1.0 | µg/L | 1 | 8/27/2021 03:07 |
| Styrene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 03:07 |
| tert-Butylbenzene | U | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 03:07 |
| Tetrachloroethene | 1.7 | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 03:07 |
| Toluene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:07 |
| trans-1,2-Dichloroethene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 03:07 |
| trans-1,3-Dichloropropene | U | | 0.38 | 2.7 | µg/L | 1 | 8/27/2021 03:07 |
| Trichloroethene | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 03:07 |
| Trichlorofluoromethane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 03:07 |
| Vinyl chloride | U | | 0.53 | 1.8 | µg/L | 1 | 8/27/2021 03:07 |
| Xylenes, Total | U | | 0.81 | 4.4 | µg/L | 1 | 8/27/2021 03:07 |
| Surr: 1,2-Dichloroethane-d4 | 98.6 | | | 75-120 | %REC | 1 | 8/27/2021 03:07 |
| Surr: 4-Bromofluorobenzene | 99.0 | | | 80-110 | %REC | 1 | 8/27/2021 03:07 |
| Surr: Dibromofluoromethane | 103 | | | 85-115 | %REC | 1 | 8/27/2021 03:07 |
| Surr: Toluene-d8 | 103 | | | 85-110 | %REC | 1 | 8/27/2021 03:07 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-5
Collection Date: 8/18/2021 10:25 AM

Work Order: 21081889
Lab ID: 21081889-05
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------------|--------|------|------|-----------------|-------|-----------------|-----------------|
| VOLATILE ORGANIC COMPOUNDS | | | | | | | |
| | | | | Method: SW8260C | | | Analyst: MF |
| 1,1,1,2-Tetrachloroethane | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 03:27 |
| 1,1,1-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 03:27 |
| 1,1,2,2-Tetrachloroethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:27 |
| 1,1,2-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 03:27 |
| 1,1-Dichloroethane | U | | 0.44 | 1.5 | µg/L | 1 | 8/27/2021 03:27 |
| 1,1-Dichloroethene | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 03:27 |
| 1,1-Dichloropropene | U | | 0.37 | 1.2 | µg/L | 1 | 8/27/2021 03:27 |
| 1,2,3-Trichlorobenzene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 03:27 |
| 1,2,3-Trichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:27 |
| 1,2,4-Trichlorobenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:27 |
| 1,2,4-Trimethylbenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:27 |
| 1,2-Dibromo-3-chloropropane | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 03:27 |
| 1,2-Dibromoethane | U | | 0.41 | 1.4 | µg/L | 1 | 8/27/2021 03:27 |
| 1,2-Dichlorobenzene | U | | 0.32 | 1.1 | µg/L | 1 | 8/27/2021 03:27 |
| 1,2-Dichloroethane | U | | 0.44 | 1.4 | µg/L | 1 | 8/27/2021 03:27 |
| 1,2-Dichloropropane | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 03:27 |
| 1,3,5-Trimethylbenzene | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 03:27 |
| 1,3-Dichlorobenzene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 03:27 |
| 1,3-Dichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:27 |
| 1,4-Dichlorobenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 03:27 |
| 2,2-Dichloropropane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 03:27 |
| 2-Butanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 03:27 |
| 2-Chlorotoluene | U | | 0.36 | 1.2 | µg/L | 1 | 8/27/2021 03:27 |
| 4-Chlorotoluene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 03:27 |
| 4-Methyl-2-pentanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 03:27 |
| Acetone | U | | 6.2 | 21 | µg/L | 1 | 8/27/2021 03:27 |
| Benzene | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 03:27 |
| Bromobenzene | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 03:27 |
| Bromochloromethane | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:27 |
| Bromodichloromethane | U | | 0.49 | 1.6 | µg/L | 1 | 8/27/2021 03:27 |
| Bromoform | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 03:27 |
| Bromomethane | U | | 0.90 | 3.0 | µg/L | 1 | 8/27/2021 03:27 |
| Carbon tetrachloride | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 03:27 |
| Chlorobenzene | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:27 |
| Chloroethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 03:27 |
| Chloroform | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 03:27 |
| Chloromethane | U | | 0.83 | 2.8 | µg/L | 1 | 8/27/2021 03:27 |
| cis-1,2-Dichloroethene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 03:27 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-5
Collection Date: 8/18/2021 10:25 AM

Work Order: 21081889
Lab ID: 21081889-05
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------|--------|------|------|--------------|-------|-----------------|-----------------|
| cis-1,3-Dichloropropene | U | | 0.57 | 1.9 | µg/L | 1 | 8/27/2021 03:27 |
| Dibromochloromethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:27 |
| Dibromomethane | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 03:27 |
| Dichlorodifluoromethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 03:27 |
| Ethylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 03:27 |
| Hexachlorobutadiene | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 03:27 |
| Isopropylbenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 03:27 |
| m,p-Xylene | U | | 0.81 | 2.7 | µg/L | 1 | 8/27/2021 03:27 |
| Methyl tert-butyl ether | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:27 |
| Methylene chloride | U | | 0.86 | 2.9 | µg/L | 1 | 8/27/2021 03:27 |
| Naphthalene | U | | 0.77 | 2.6 | µg/L | 1 | 8/27/2021 03:27 |
| n-Butylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 03:27 |
| n-Propylbenzene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 03:27 |
| o-Xylene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 03:27 |
| p-Isopropyltoluene | U | | 0.26 | 0.88 | µg/L | 1 | 8/27/2021 03:27 |
| sec-Butylbenzene | U | | 0.30 | 1.0 | µg/L | 1 | 8/27/2021 03:27 |
| Styrene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 03:27 |
| tert-Butylbenzene | U | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 03:27 |
| Tetrachloroethene | U | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 03:27 |
| Toluene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:27 |
| trans-1,2-Dichloroethene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 03:27 |
| trans-1,3-Dichloropropene | U | | 0.38 | 2.7 | µg/L | 1 | 8/27/2021 03:27 |
| Trichloroethene | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 03:27 |
| Trichlorofluoromethane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 03:27 |
| Vinyl chloride | U | | 0.53 | 1.8 | µg/L | 1 | 8/27/2021 03:27 |
| Xylenes, Total | U | | 0.81 | 4.4 | µg/L | 1 | 8/27/2021 03:27 |
| Surr: 1,2-Dichloroethane-d4 | 98.7 | | | 75-120 | %REC | 1 | 8/27/2021 03:27 |
| Surr: 4-Bromofluorobenzene | 93.2 | | | 80-110 | %REC | 1 | 8/27/2021 03:27 |
| Surr: Dibromofluoromethane | 98.2 | | | 85-115 | %REC | 1 | 8/27/2021 03:27 |
| Surr: Toluene-d8 | 101 | | | 85-110 | %REC | 1 | 8/27/2021 03:27 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-6
Collection Date: 8/18/2021 12:10 PM

Work Order: 21081889
Lab ID: 21081889-06
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------------|--------|------|------|-----------------|-------|-----------------|-----------------|
| VOLATILE ORGANIC COMPOUNDS | | | | | | | |
| | | | | Method: SW8260C | | | Analyst: MF |
| 1,1,1,2-Tetrachloroethane | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 03:47 |
| 1,1,1-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 03:47 |
| 1,1,2,2-Tetrachloroethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:47 |
| 1,1,2-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 03:47 |
| 1,1-Dichloroethane | U | | 0.44 | 1.5 | µg/L | 1 | 8/27/2021 03:47 |
| 1,1-Dichloroethene | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 03:47 |
| 1,1-Dichloropropene | U | | 0.37 | 1.2 | µg/L | 1 | 8/27/2021 03:47 |
| 1,2,3-Trichlorobenzene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 03:47 |
| 1,2,3-Trichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:47 |
| 1,2,4-Trichlorobenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:47 |
| 1,2,4-Trimethylbenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:47 |
| 1,2-Dibromo-3-chloropropane | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 03:47 |
| 1,2-Dibromoethane | U | | 0.41 | 1.4 | µg/L | 1 | 8/27/2021 03:47 |
| 1,2-Dichlorobenzene | U | | 0.32 | 1.1 | µg/L | 1 | 8/27/2021 03:47 |
| 1,2-Dichloroethane | U | | 0.44 | 1.4 | µg/L | 1 | 8/27/2021 03:47 |
| 1,2-Dichloropropane | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 03:47 |
| 1,3,5-Trimethylbenzene | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 03:47 |
| 1,3-Dichlorobenzene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 03:47 |
| 1,3-Dichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:47 |
| 1,4-Dichlorobenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 03:47 |
| 2,2-Dichloropropane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 03:47 |
| 2-Butanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 03:47 |
| 2-Chlorotoluene | U | | 0.36 | 1.2 | µg/L | 1 | 8/27/2021 03:47 |
| 4-Chlorotoluene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 03:47 |
| 4-Methyl-2-pentanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 03:47 |
| Acetone | U | | 6.2 | 21 | µg/L | 1 | 8/27/2021 03:47 |
| Benzene | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 03:47 |
| Bromobenzene | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 03:47 |
| Bromochloromethane | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:47 |
| Bromodichloromethane | U | | 0.49 | 1.6 | µg/L | 1 | 8/27/2021 03:47 |
| Bromoform | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 03:47 |
| Bromomethane | U | | 0.90 | 3.0 | µg/L | 1 | 8/27/2021 03:47 |
| Carbon tetrachloride | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 03:47 |
| Chlorobenzene | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:47 |
| Chloroethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 03:47 |
| Chloroform | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 03:47 |
| Chloromethane | U | | 0.83 | 2.8 | µg/L | 1 | 8/27/2021 03:47 |
| cis-1,2-Dichloroethene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 03:47 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-6
Collection Date: 8/18/2021 12:10 PM

Work Order: 21081889
Lab ID: 21081889-06
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------|------------|------|-------------|--------------|-------------|-----------------|-----------------|
| cis-1,3-Dichloropropene | U | | 0.57 | 1.9 | µg/L | 1 | 8/27/2021 03:47 |
| Dibromochloromethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 03:47 |
| Dibromomethane | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 03:47 |
| Dichlorodifluoromethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 03:47 |
| Ethylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 03:47 |
| Hexachlorobutadiene | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 03:47 |
| Isopropylbenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 03:47 |
| m,p-Xylene | U | | 0.81 | 2.7 | µg/L | 1 | 8/27/2021 03:47 |
| Methyl tert-butyl ether | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:47 |
| Methylene chloride | U | | 0.86 | 2.9 | µg/L | 1 | 8/27/2021 03:47 |
| Naphthalene | U | | 0.77 | 2.6 | µg/L | 1 | 8/27/2021 03:47 |
| n-Butylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 03:47 |
| n-Propylbenzene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 03:47 |
| o-Xylene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 03:47 |
| p-Isopropyltoluene | U | | 0.26 | 0.88 | µg/L | 1 | 8/27/2021 03:47 |
| sec-Butylbenzene | U | | 0.30 | 1.0 | µg/L | 1 | 8/27/2021 03:47 |
| Styrene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 03:47 |
| tert-Butylbenzene | U | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 03:47 |
| Tetrachloroethene | 6.0 | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 03:47 |
| Toluene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 03:47 |
| trans-1,2-Dichloroethene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 03:47 |
| trans-1,3-Dichloropropene | U | | 0.38 | 2.7 | µg/L | 1 | 8/27/2021 03:47 |
| Trichloroethene | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 03:47 |
| Trichlorofluoromethane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 03:47 |
| Vinyl chloride | U | | 0.53 | 1.8 | µg/L | 1 | 8/27/2021 03:47 |
| Xylenes, Total | U | | 0.81 | 4.4 | µg/L | 1 | 8/27/2021 03:47 |
| Surr: 1,2-Dichloroethane-d4 | 97.3 | | | 75-120 | %REC | 1 | 8/27/2021 03:47 |
| Surr: 4-Bromofluorobenzene | 95.7 | | | 80-110 | %REC | 1 | 8/27/2021 03:47 |
| Surr: Dibromofluoromethane | 101 | | | 85-115 | %REC | 1 | 8/27/2021 03:47 |
| Surr: Toluene-d8 | 99.6 | | | 85-110 | %REC | 1 | 8/27/2021 03:47 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-7
Collection Date: 8/18/2021 11:05 AM

Work Order: 21081889
Lab ID: 21081889-07
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------------|-------------|------|-------------|------------------------|-------------|-----------------|--------------------|
| VOLATILE ORGANIC COMPOUNDS | | | | | | | |
| | | | | Method: SW8260C | | | Analyst: MF |
| 1,1,1,2-Tetrachloroethane | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 04:07 |
| 1,1,1-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 04:07 |
| 1,1,2,2-Tetrachloroethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 04:07 |
| 1,1,2-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 04:07 |
| 1,1-Dichloroethane | U | | 0.44 | 1.5 | µg/L | 1 | 8/27/2021 04:07 |
| 1,1-Dichloroethene | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 04:07 |
| 1,1-Dichloropropene | U | | 0.37 | 1.2 | µg/L | 1 | 8/27/2021 04:07 |
| 1,2,3-Trichlorobenzene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 04:07 |
| 1,2,3-Trichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 04:07 |
| 1,2,4-Trichlorobenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 04:07 |
| 1,2,4-Trimethylbenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 04:07 |
| 1,2-Dibromo-3-chloropropane | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 04:07 |
| 1,2-Dibromoethane | U | | 0.41 | 1.4 | µg/L | 1 | 8/27/2021 04:07 |
| 1,2-Dichlorobenzene | U | | 0.32 | 1.1 | µg/L | 1 | 8/27/2021 04:07 |
| 1,2-Dichloroethane | U | | 0.44 | 1.4 | µg/L | 1 | 8/27/2021 04:07 |
| 1,2-Dichloropropane | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 04:07 |
| 1,3,5-Trimethylbenzene | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 04:07 |
| 1,3-Dichlorobenzene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 04:07 |
| 1,3-Dichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 04:07 |
| 1,4-Dichlorobenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 04:07 |
| 2,2-Dichloropropane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 04:07 |
| 2-Butanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 04:07 |
| 2-Chlorotoluene | U | | 0.36 | 1.2 | µg/L | 1 | 8/27/2021 04:07 |
| 4-Chlorotoluene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 04:07 |
| 4-Methyl-2-pentanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 04:07 |
| Acetone | U | | 6.2 | 21 | µg/L | 1 | 8/27/2021 04:07 |
| Benzene | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 04:07 |
| Bromobenzene | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 04:07 |
| Bromochloromethane | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 04:07 |
| Bromodichloromethane | U | | 0.49 | 1.6 | µg/L | 1 | 8/27/2021 04:07 |
| Bromoform | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 04:07 |
| Bromomethane | U | | 0.90 | 3.0 | µg/L | 1 | 8/27/2021 04:07 |
| Carbon tetrachloride | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 04:07 |
| Chlorobenzene | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 04:07 |
| Chloroethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 04:07 |
| Chloroform | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 04:07 |
| Chloromethane | 0.86 | J | 0.83 | 2.8 | µg/L | 1 | 8/27/2021 04:07 |
| cis-1,2-Dichloroethene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 04:07 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-7
Collection Date: 8/18/2021 11:05 AM

Work Order: 21081889
Lab ID: 21081889-07
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------|-------------|------|-------------|--------------|-------------|-----------------|-----------------|
| cis-1,3-Dichloropropene | U | | 0.57 | 1.9 | µg/L | 1 | 8/27/2021 04:07 |
| Dibromochloromethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 04:07 |
| Dibromomethane | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 04:07 |
| Dichlorodifluoromethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 04:07 |
| Ethylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 04:07 |
| Hexachlorobutadiene | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 04:07 |
| Isopropylbenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 04:07 |
| m,p-Xylene | U | | 0.81 | 2.7 | µg/L | 1 | 8/27/2021 04:07 |
| Methyl tert-butyl ether | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 04:07 |
| Methylene chloride | U | | 0.86 | 2.9 | µg/L | 1 | 8/27/2021 04:07 |
| Naphthalene | U | | 0.77 | 2.6 | µg/L | 1 | 8/27/2021 04:07 |
| n-Butylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 04:07 |
| n-Propylbenzene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 04:07 |
| o-Xylene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 04:07 |
| p-Isopropyltoluene | U | | 0.26 | 0.88 | µg/L | 1 | 8/27/2021 04:07 |
| sec-Butylbenzene | U | | 0.30 | 1.0 | µg/L | 1 | 8/27/2021 04:07 |
| Styrene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 04:07 |
| tert-Butylbenzene | U | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 04:07 |
| Tetrachloroethene | 0.66 | J | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 04:07 |
| Toluene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 04:07 |
| trans-1,2-Dichloroethene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 04:07 |
| trans-1,3-Dichloropropene | U | | 0.38 | 2.7 | µg/L | 1 | 8/27/2021 04:07 |
| Trichloroethene | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 04:07 |
| Trichlorofluoromethane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 04:07 |
| Vinyl chloride | U | | 0.53 | 1.8 | µg/L | 1 | 8/27/2021 04:07 |
| Xylenes, Total | U | | 0.81 | 4.4 | µg/L | 1 | 8/27/2021 04:07 |
| Surr: 1,2-Dichloroethane-d4 | 100 | | | 75-120 | %REC | 1 | 8/27/2021 04:07 |
| Surr: 4-Bromofluorobenzene | 100 | | | 80-110 | %REC | 1 | 8/27/2021 04:07 |
| Surr: Dibromofluoromethane | 102 | | | 85-115 | %REC | 1 | 8/27/2021 04:07 |
| Surr: Toluene-d8 | 104 | | | 85-110 | %REC | 1 | 8/27/2021 04:07 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-1 DUP
Collection Date: 8/18/2021 02:05 PM

Work Order: 21081889
Lab ID: 21081889-08
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------------|--------|------|------|-----------------|-------|-----------------|-----------------|
| VOLATILE ORGANIC COMPOUNDS | | | | | | | |
| | | | | Method: SW8260C | | | Analyst: MF |
| 1,1,1,2-Tetrachloroethane | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 04:27 |
| 1,1,1-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 04:27 |
| 1,1,2,2-Tetrachloroethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 04:27 |
| 1,1,2-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 04:27 |
| 1,1-Dichloroethane | U | | 0.44 | 1.5 | µg/L | 1 | 8/27/2021 04:27 |
| 1,1-Dichloroethene | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 04:27 |
| 1,1-Dichloropropene | U | | 0.37 | 1.2 | µg/L | 1 | 8/27/2021 04:27 |
| 1,2,3-Trichlorobenzene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 04:27 |
| 1,2,3-Trichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 04:27 |
| 1,2,4-Trichlorobenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 04:27 |
| 1,2,4-Trimethylbenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 04:27 |
| 1,2-Dibromo-3-chloropropane | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 04:27 |
| 1,2-Dibromoethane | U | | 0.41 | 1.4 | µg/L | 1 | 8/27/2021 04:27 |
| 1,2-Dichlorobenzene | U | | 0.32 | 1.1 | µg/L | 1 | 8/27/2021 04:27 |
| 1,2-Dichloroethane | U | | 0.44 | 1.4 | µg/L | 1 | 8/27/2021 04:27 |
| 1,2-Dichloropropane | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 04:27 |
| 1,3,5-Trimethylbenzene | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 04:27 |
| 1,3-Dichlorobenzene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 04:27 |
| 1,3-Dichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 04:27 |
| 1,4-Dichlorobenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 04:27 |
| 2,2-Dichloropropane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 04:27 |
| 2-Butanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 04:27 |
| 2-Chlorotoluene | U | | 0.36 | 1.2 | µg/L | 1 | 8/27/2021 04:27 |
| 4-Chlorotoluene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 04:27 |
| 4-Methyl-2-pentanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 04:27 |
| Acetone | U | | 6.2 | 21 | µg/L | 1 | 8/27/2021 04:27 |
| Benzene | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 04:27 |
| Bromobenzene | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 04:27 |
| Bromochloromethane | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 04:27 |
| Bromodichloromethane | U | | 0.49 | 1.6 | µg/L | 1 | 8/27/2021 04:27 |
| Bromoform | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 04:27 |
| Bromomethane | U | | 0.90 | 3.0 | µg/L | 1 | 8/27/2021 04:27 |
| Carbon tetrachloride | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 04:27 |
| Chlorobenzene | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 04:27 |
| Chloroethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 04:27 |
| Chloroform | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 04:27 |
| Chloromethane | U | | 0.83 | 2.8 | µg/L | 1 | 8/27/2021 04:27 |
| cis-1,2-Dichloroethene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 04:27 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: MW-1 DUP
Collection Date: 8/18/2021 02:05 PM

Work Order: 21081889
Lab ID: 21081889-08
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------|------------|------|------------|--------------|-------------|-----------------|-----------------|
| cis-1,3-Dichloropropene | U | | 0.57 | 1.9 | µg/L | 1 | 8/27/2021 04:27 |
| Dibromochloromethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 04:27 |
| Dibromomethane | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 04:27 |
| Dichlorodifluoromethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 04:27 |
| Ethylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 04:27 |
| Hexachlorobutadiene | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 04:27 |
| Isopropylbenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 04:27 |
| m,p-Xylene | U | | 0.81 | 2.7 | µg/L | 1 | 8/27/2021 04:27 |
| Methyl tert-butyl ether | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 04:27 |
| Methylene chloride | U | | 0.86 | 2.9 | µg/L | 1 | 8/27/2021 04:27 |
| Naphthalene | U | | 0.77 | 2.6 | µg/L | 1 | 8/27/2021 04:27 |
| n-Butylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 04:27 |
| n-Propylbenzene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 04:27 |
| o-Xylene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 04:27 |
| p-Isopropyltoluene | U | | 0.26 | 0.88 | µg/L | 1 | 8/27/2021 04:27 |
| sec-Butylbenzene | U | | 0.30 | 1.0 | µg/L | 1 | 8/27/2021 04:27 |
| Styrene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 04:27 |
| tert-Butylbenzene | U | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 04:27 |
| Tetrachloroethene | 120 | | 2.0 | 6.6 | µg/L | 5 | 8/27/2021 16:11 |
| Toluene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 04:27 |
| trans-1,2-Dichloroethene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 04:27 |
| trans-1,3-Dichloropropene | U | | 0.38 | 2.7 | µg/L | 1 | 8/27/2021 04:27 |
| Trichloroethene | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 04:27 |
| Trichlorofluoromethane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 04:27 |
| Vinyl chloride | U | | 0.53 | 1.8 | µg/L | 1 | 8/27/2021 04:27 |
| Xylenes, Total | U | | 0.81 | 4.4 | µg/L | 1 | 8/27/2021 04:27 |
| Surr: 1,2-Dichloroethane-d4 | 103 | | | 75-120 | %REC | 1 | 8/27/2021 04:27 |
| Surr: 1,2-Dichloroethane-d4 | 102 | | | 75-120 | %REC | 5 | 8/27/2021 16:11 |
| Surr: 4-Bromofluorobenzene | 97.7 | | | 80-110 | %REC | 1 | 8/27/2021 04:27 |
| Surr: 4-Bromofluorobenzene | 98.3 | | | 80-110 | %REC | 5 | 8/27/2021 16:11 |
| Surr: Dibromofluoromethane | 106 | | | 85-115 | %REC | 1 | 8/27/2021 04:27 |
| Surr: Dibromofluoromethane | 99.0 | | | 85-115 | %REC | 5 | 8/27/2021 16:11 |
| Surr: Toluene-d8 | 101 | | | 85-110 | %REC | 1 | 8/27/2021 04:27 |
| Surr: Toluene-d8 | 102 | | | 85-110 | %REC | 5 | 8/27/2021 16:11 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: Field Blank
Collection Date: 8/18/2021

Work Order: 21081889
Lab ID: 21081889-09
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|--|--------|------|------|--------------|-------|-----------------|-----------------|
| PFAS BY EPA 537 MODIFIED | | | | | | | |
| Fluorotelomer Sulphonic Acid 4:2 (FtS 4:2) | U | | 0.98 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Fluorotelomer Sulphonic Acid 6:2 (FtS 6:2) | U | | 0.69 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Fluorotelomer Sulphonic Acid 8:2 (FtS 8:2) | U | | 1.2 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Fluorotelomer Sulphonic Acid 10:2 (FtS 10:2) | U | | 0.94 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorobutanesulfonic Acid (PFBS) | U | | 0.37 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorobutanoic Acid (PFBA) | U | | 2.7 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorodecanesulfonic Acid (PFDS) | U | | 1.4 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorodecanoic Acid (PFDA) | U | | 1.3 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorododecanesulfonic Acid (PFDoS) | U | | 1.5 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorododecanoic Acid (PFDoA) | U | | 1.5 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluoroheptanesulfonic Acid (PFHpS) | U | | 0.59 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluoroheptanoic Acid (PFHpA) | U | | 0.46 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorohexadecanoic Acid (PFHxDA) | U | | 0.40 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorohexamenesulfonic Acid (PFHxS) | U | | 0.38 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorohexanoic Acid (PFHxA) | U | | 1.2 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorononanesulfonic Acid (PFNS) | U | | 0.52 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorononanoic Acid (PFNA) | U | | 0.91 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorooctadecanoic Acid (PFODA) | U | | 0.68 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorooctanesulfonamide (PFOSA) | U | | 0.74 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorooctanesulfonic Acid (PFOS) | U | | 0.93 | 2.1 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorooctanoic Acid (PFOA) | U | | 0.66 | 2.1 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluoropentanesulfonic Acid (PPeS) | U | | 0.58 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluoropentanoic Acid (PPeA) | U | | 1.3 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorotetradecanoic Acid (PFTeA) | U | | 2.7 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluorotridecanoic Acid (PFTriA) | U | | 0.80 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Perfluoroundecanoic Acid (PFUnA) | U | | 1.0 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| N-ethylperfluoro-1-octanesulfonamide | U | | 1.2 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| N-Ethylperfluorooctanesulfonamidoacetic Acid | U | | 0.65 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| N-Ethylperfluorooctanesulfonamidoethanol | U | | 0.54 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| N-methylperfluoro-1-octanesulfonamide | U | | 0.83 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: Field Blank
Collection Date: 8/18/2021

Work Order: 21081889
Lab ID: 21081889-09
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|--|--------|------|------|--------------|-------|-----------------|-----------------|
| N-Methylperfluorooctanesulfonamidoacetic Acid | U | | 0.67 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| N-Methylperfluorooctanesulfonamidoethanol | U | | 0.50 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Hexafluoropropylene oxide dimer acid (HFPO-DA) | U | | 1.2 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| 4,8-Dioxa-3H-perfluorononanoic Acid (DONA) | U | | 0.59 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| 11Cl-Pf3OUdS | U | | 0.49 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| 9Cl-PF3ONS | U | | 0.47 | 5.2 | ng/L | 1 | 8/23/2021 21:05 |
| Surr: 13C2-FtS 4:2 | 78.5 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C2-FtS 6:2 | 89.2 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C2-FtS 8:2 | 82.9 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C2-PFDA | 88.9 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C2-PFDoA | 79.9 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C2-PFHxA | 79.2 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C2-PFHxDA | 73.5 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C2-PFTeA | 69.2 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C2-PFUra | 90.3 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C3-HFPO-DA | 76.5 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C3-PFBS | 77.7 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C4-PFBA | 80.4 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C4-PFHxA | 67.6 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C4-PFOA | 82.1 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C4-PFOS | 80.7 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C5-PFNA | 73.4 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C5-PFPeA | 77.4 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 13C8-FOSA | 68.5 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: 18O2-PFHxA | 80.1 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: d5-N-EtFOSA | 58.6 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: d5-N-EtFOSAA | 99.9 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: d9-N-EtFOSE | 66.5 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: d3-N-MeFOSA | 58.8 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: d3-N-MeFOSAA | 88.2 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |
| Surr: d7-N-MeFOSE | 67.4 | | | 50-150 | %REC | 1 | 8/23/2021 21:05 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: Trip Blank
Collection Date: 8/18/2021

Work Order: 21081889
Lab ID: 21081889-10
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------------|--------|------|------|-----------------|-------|-----------------|-----------------|
| VOLATILE ORGANIC COMPOUNDS | | | | | | | |
| | | | | Method: SW8260C | | | Analyst: MF |
| 1,1,1,2-Tetrachloroethane | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 00:49 |
| 1,1,1-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 00:49 |
| 1,1,2,2-Tetrachloroethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 00:49 |
| 1,1,2-Trichloroethane | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 00:49 |
| 1,1-Dichloroethane | U | | 0.44 | 1.5 | µg/L | 1 | 8/27/2021 00:49 |
| 1,1-Dichloroethene | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 00:49 |
| 1,1-Dichloropropene | U | | 0.37 | 1.2 | µg/L | 1 | 8/27/2021 00:49 |
| 1,2,3-Trichlorobenzene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 00:49 |
| 1,2,3-Trichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 00:49 |
| 1,2,4-Trichlorobenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 00:49 |
| 1,2,4-Trimethylbenzene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 00:49 |
| 1,2-Dibromo-3-chloropropane | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 00:49 |
| 1,2-Dibromoethane | U | | 0.41 | 1.4 | µg/L | 1 | 8/27/2021 00:49 |
| 1,2-Dichlorobenzene | U | | 0.32 | 1.1 | µg/L | 1 | 8/27/2021 00:49 |
| 1,2-Dichloroethane | U | | 0.44 | 1.4 | µg/L | 1 | 8/27/2021 00:49 |
| 1,2-Dichloropropane | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 00:49 |
| 1,3,5-Trimethylbenzene | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 00:49 |
| 1,3-Dichlorobenzene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 00:49 |
| 1,3-Dichloropropane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 00:49 |
| 1,4-Dichlorobenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 00:49 |
| 2,2-Dichloropropane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 00:49 |
| 2-Butanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 00:49 |
| 2-Chlorotoluene | U | | 0.36 | 1.2 | µg/L | 1 | 8/27/2021 00:49 |
| 4-Chlorotoluene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 00:49 |
| 4-Methyl-2-pentanone | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 00:49 |
| Acetone | U | | 6.2 | 21 | µg/L | 1 | 8/27/2021 00:49 |
| Benzene | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 00:49 |
| Bromobenzene | U | | 0.38 | 1.3 | µg/L | 1 | 8/27/2021 00:49 |
| Bromochloromethane | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 00:49 |
| Bromodichloromethane | U | | 0.49 | 1.6 | µg/L | 1 | 8/27/2021 00:49 |
| Bromoform | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 00:49 |
| Bromomethane | U | | 0.90 | 3.0 | µg/L | 1 | 8/27/2021 00:49 |
| Carbon tetrachloride | U | | 0.40 | 1.4 | µg/L | 1 | 8/27/2021 00:49 |
| Chlorobenzene | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 00:49 |
| Chloroethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 00:49 |
| Chloroform | U | | 0.46 | 1.5 | µg/L | 1 | 8/27/2021 00:49 |
| Chloromethane | U | | 0.83 | 2.8 | µg/L | 1 | 8/27/2021 00:49 |
| cis-1,2-Dichloroethene | U | | 0.42 | 1.4 | µg/L | 1 | 8/27/2021 00:49 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Project: Koeller (47358.003)
Sample ID: Trip Blank
Collection Date: 8/18/2021

Work Order: 21081889
Lab ID: 21081889-10
Matrix: WATER

| Analyses | Result | Qual | MDL | Report Limit | Units | Dilution Factor | Date Analyzed |
|-----------------------------|--------|------|------|--------------|-------|-----------------|-----------------|
| cis-1,3-Dichloropropene | U | | 0.57 | 1.9 | µg/L | 1 | 8/27/2021 00:49 |
| Dibromochloromethane | U | | 0.40 | 1.3 | µg/L | 1 | 8/27/2021 00:49 |
| Dibromomethane | U | | 0.65 | 2.2 | µg/L | 1 | 8/27/2021 00:49 |
| Dichlorodifluoromethane | U | | 0.68 | 2.3 | µg/L | 1 | 8/27/2021 00:49 |
| Ethylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 00:49 |
| Hexachlorobutadiene | U | | 0.56 | 1.9 | µg/L | 1 | 8/27/2021 00:49 |
| Isopropylbenzene | U | | 0.35 | 1.2 | µg/L | 1 | 8/27/2021 00:49 |
| m,p-Xylene | U | | 0.81 | 2.7 | µg/L | 1 | 8/27/2021 00:49 |
| Methyl tert-butyl ether | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 00:49 |
| Methylene chloride | U | | 0.86 | 2.9 | µg/L | 1 | 8/27/2021 00:49 |
| Naphthalene | U | | 0.77 | 2.6 | µg/L | 1 | 8/27/2021 00:49 |
| n-Butylbenzene | U | | 0.34 | 1.1 | µg/L | 1 | 8/27/2021 00:49 |
| n-Propylbenzene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 00:49 |
| o-Xylene | U | | 0.31 | 1.0 | µg/L | 1 | 8/27/2021 00:49 |
| p-Isopropyltoluene | U | | 0.26 | 0.88 | µg/L | 1 | 8/27/2021 00:49 |
| sec-Butylbenzene | U | | 0.30 | 1.0 | µg/L | 1 | 8/27/2021 00:49 |
| Styrene | U | | 0.33 | 1.1 | µg/L | 1 | 8/27/2021 00:49 |
| tert-Butylbenzene | U | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 00:49 |
| Tetrachloroethene | U | | 0.39 | 1.3 | µg/L | 1 | 8/27/2021 00:49 |
| Toluene | U | | 0.45 | 1.5 | µg/L | 1 | 8/27/2021 00:49 |
| trans-1,2-Dichloroethene | U | | 0.48 | 1.6 | µg/L | 1 | 8/27/2021 00:49 |
| trans-1,3-Dichloropropene | U | | 0.38 | 2.7 | µg/L | 1 | 8/27/2021 00:49 |
| Trichloroethene | U | | 0.43 | 1.4 | µg/L | 1 | 8/27/2021 00:49 |
| Trichlorofluoromethane | U | | 0.52 | 1.7 | µg/L | 1 | 8/27/2021 00:49 |
| Vinyl chloride | U | | 0.53 | 1.8 | µg/L | 1 | 8/27/2021 00:49 |
| Xylenes, Total | U | | 0.81 | 4.4 | µg/L | 1 | 8/27/2021 00:49 |
| Surr: 1,2-Dichloroethane-d4 | 99.2 | | | 75-120 | %REC | 1 | 8/27/2021 00:49 |
| Surr: 4-Bromofluorobenzene | 93.8 | | | 80-110 | %REC | 1 | 8/27/2021 00:49 |
| Surr: Dibromofluoromethane | 98.4 | | | 85-115 | %REC | 1 | 8/27/2021 00:49 |
| Surr: Toluene-d8 | 102 | | | 85-110 | %REC | 1 | 8/27/2021 00:49 |

Note: See Qualifiers page for a list of qualifiers and their definitions.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: 182387 | | Instrument ID LCMS1 | | Method: E537 Mod | | | | | | | | | | |
|--------------------------------|--------|-------------------------------|-----|------------------|----------------|------|-----------------------------------|---------------|-------|-----------|------|--|--|--|
| MBLK | | Sample ID: MBLK-182387-182387 | | | Units: ng/L | | Analysis Date: 8/23/2021 06:07 PM | | | | | | | |
| Client ID: | | Run ID: LCMS1_210823B | | | SeqNo: 7690005 | | Prep Date: 8/23/2021 | | DF: 1 | | | | | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | %RPD | RPD Limit | Qual | | | |
| Fluorotelomer Sulphonic Acid | U | 0.94 | 5.0 | | | | | | | | | | | |
| Fluorotelomer Sulphonic Acid | U | 0.66 | 5.0 | | | | | | | | | | | |
| Fluorotelomer Sulphonic Acid | U | 1.1 | 5.0 | | | | | | | | | | | |
| Fluorotelomer Sulphonic Acid | U | 0.9 | 5.0 | | | | | | | | | | | |
| Perfluorobutanesulfonic Acid (| U | 0.35 | 5.0 | | | | | | | | | | | |
| Perfluorobutanoic Acid (PFBA) | U | 2.6 | 5.0 | | | | | | | | | | | |
| Perfluorodecanesulfonic Acid (| U | 1.4 | 5.0 | | | | | | | | | | | |
| Perfluorodecanoic Acid (PFDA) | U | 1.2 | 5.0 | | | | | | | | | | | |
| Perfluorododecanesulfonic Aci | U | 1.4 | 5.0 | | | | | | | | | | | |
| Perfluorododecanoic Acid (PFI) | U | 1.4 | 5.0 | | | | | | | | | | | |
| Perfluoroheptanesulfonic Acid | U | 0.57 | 5.0 | | | | | | | | | | | |
| Perfluoroheptanoic Acid (PFH) | U | 0.44 | 5.0 | | | | | | | | | | | |
| Perfluorohexadecanoic Acid (F | U | 0.38 | 5.0 | | | | | | | | | | | |
| Perfluorohexanesulfonic Acid (| U | 0.37 | 5.0 | | | | | | | | | | | |
| Perfluorohexanoic Acid (PFHx | U | 1.2 | 5.0 | | | | | | | | | | | |
| Perfluorononanesulfonic Acid (| U | 0.5 | 5.0 | | | | | | | | | | | |
| Perfluorononanoic Acid (PFNA) | U | 0.87 | 5.0 | | | | | | | | | | | |
| Perfluoroctadecanoic Acid (P | U | 0.65 | 5.0 | | | | | | | | | | | |
| Perfluoroctanesulfonamide (F | U | 0.71 | 5.0 | | | | | | | | | | | |
| Perfluoroctanesulfonic Acid (I | U | 0.89 | 2.0 | | | | | | | | | | | |
| Perfluoroctanoic Acid (PFOA) | U | 0.63 | 2.0 | | | | | | | | | | | |
| Perfluoropentanesulfonic Acid | U | 0.56 | 5.0 | | | | | | | | | | | |
| Perfluoropentanoic Acid (PPF) | U | 1.3 | 5.0 | | | | | | | | | | | |
| Perfluorotetradecanoic Acid (F | U | 2.6 | 5.0 | | | | | | | | | | | |
| Perfluorotridecanoic Acid (PF1 | U | 0.77 | 5.0 | | | | | | | | | | | |
| Perfluoroundecanoic Acid (PFI) | U | 0.97 | 5.0 | | | | | | | | | | | |
| N-ethylperfluoro-1-octanesulfo | U | 1.2 | 5.0 | | | | | | | | | | | |
| N-Ethylperfluoroctanesulfona | U | 0.63 | 5.0 | | | | | | | | | | | |
| N-Ethylperfluoroctanesulfona | U | 0.52 | 5.0 | | | | | | | | | | | |
| N-methylperfluoro-1-octanesul | U | 0.79 | 5.0 | | | | | | | | | | | |
| N-Methylperfluoroctanesulfur | U | 0.64 | 5.0 | | | | | | | | | | | |
| N-Methylperfluoroctanesulfur | U | 0.48 | 5.0 | | | | | | | | | | | |
| Hexafluoropropylene oxide din | U | 1.2 | 5.0 | | | | | | | | | | | |
| 4,8-Dioxa-3H-perfluorononano | U | 0.56 | 5.0 | | | | | | | | | | | |
| 11CI-Pf3OUdS | U | 0.47 | 5.0 | | | | | | | | | | | |
| 9CI-PF3ONS | U | 0.45 | 5.0 | | | | | | | | | | | |
| Surr: 13C2-FtS 4:2 | 114.4 | 0 | 0 | 149.4 | 0 | 76.6 | 50-150 | 0 | | | | | | |
| Surr: 13C2-FtS 6:2 | 166.4 | 0 | 0 | 152 | 0 | 109 | 50-150 | 0 | | | | | | |
| Surr: 13C2-FtS 8:2 | 135.8 | 0 | 0 | 153.3 | 0 | 88.6 | 50-150 | 0 | | | | | | |
| Surr: 13C2-PFDA | 124.8 | 0 | 0 | 160 | 0 | 78 | 50-150 | 0 | | | | | | |
| Surr: 13C2-PFDa | 173.9 | 0 | 0 | 160 | 0 | 109 | 50-150 | 0 | | | | | | |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: 182387 | Instrument ID LCMS1 | Method: E537 Mod | | | | | | |
|-------------------------|----------------------------|-------------------------|---|-------|---|------|--------|---|
| Surr: 13C2-PFHxA | 124.5 | 0 | 0 | 160 | 0 | 77.8 | 50-150 | 0 |
| Surr: 13C2-PFHxDA | 159.9 | 0 | 0 | 160 | 0 | 99.9 | 50-150 | 0 |
| Surr: 13C2-PFTeA | 140.3 | 0 | 0 | 160 | 0 | 87.7 | 50-150 | 0 |
| Surr: 13C2-PFUnA | 156.6 | 0 | 0 | 160 | 0 | 97.8 | 50-150 | 0 |
| Surr: 13C3-HFPO-DA | 139.2 | 0 | 0 | 160 | 0 | 87 | 50-150 | 0 |
| Surr: 13C3-PFBS | 130.5 | 0 | 0 | 148.8 | 0 | 87.7 | 50-150 | 0 |
| Surr: 13C4-PFBA | 123.5 | 0 | 0 | 160 | 0 | 77.2 | 50-150 | 0 |
| Surr: 13C4-PFHxA | 128 | 0 | 0 | 160 | 0 | 80 | 50-150 | 0 |
| Surr: 13C4-PFOA | 143.4 | 0 | 0 | 160 | 0 | 89.6 | 50-150 | 0 |
| Surr: 13C4-PFOS | 117.6 | 0 | 0 | 152.8 | 0 | 76.9 | 50-150 | 0 |
| Surr: 13C5-PFNA | 130.5 | 0 | 0 | 160 | 0 | 81.6 | 50-150 | 0 |
| Surr: 13C5-PFPeA | 136.6 | 0 | 0 | 160 | 0 | 85.4 | 50-150 | 0 |
| Surr: 13C8-FOSA | 142.8 | 0 | 0 | 160 | 0 | 89.2 | 50-150 | 0 |
| Surr: 18O2-PFHxS | 138.7 | 0 | 0 | 151.2 | 0 | 91.8 | 50-150 | 0 |
| Surr: d5-N-EtFOSA | 98.59 | 0 | 0 | 160 | 0 | 61.6 | 50-150 | 0 |
| Surr: d5-N-EtFOSAA | 179.2 | 0 | 0 | 160 | 0 | 112 | 50-150 | 0 |
| Surr: d9-N-EtFOSE | 125.8 | 0 | 0 | 160 | 0 | 78.6 | 50-150 | 0 |
| Surr: d3-N-MeFOSA | 102.6 | 0 | 0 | 160 | 0 | 64.2 | 50-150 | 0 |
| Surr: d3-N-MeFOSAA | 143.1 | 0 | 0 | 160 | 0 | 89.4 | 50-150 | 0 |
| Surr: d7-N-MeFOSE | 136.5 | 0 | 0 | 160 | 0 | 85.3 | 50-150 | 0 |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

Batch ID: **182387** Instrument ID **LCMS1** Method: **E537 Mod**

| LCS | | Sample ID: LCS-182387-182387 | | | | Units: ng/L | | Analysis Date: 8/23/2021 06:17 PM | | | |
|---|--------|-------------------------------------|-----|---------|-----------------------|--------------------|-----------------------------|--|--------------|-----------|------|
| Client ID: | | Run ID: LCMS1_210823B | | | SeqNo: 7690006 | | Prep Date: 8/23/2021 | | DF: 1 | | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | %RPD | RPD Limit | Qual |
| Fluorotelomer Sulphonic Acid | 36.14 | 0.94 | 5.0 | 29.9 | 0 | 121 | 63-143 | 0 | | | |
| Fluorotelomer Sulphonic Acid | 33.53 | 0.66 | 5.0 | 30.3 | 0 | 111 | 64-140 | 0 | | | |
| Fluorotelomer Sulphonic Acid | 35.03 | 1.1 | 5.0 | 30.7 | 0 | 114 | 67-138 | 0 | | | |
| Fluorotelomer Sulphonic Acid | 52.6 | 0.9 | 5.0 | 30.8 | 0 | 171 | 40-160 | 0 | | S | |
| Perfluorobutanesulfonic Acid (| 30.09 | 0.35 | 5.0 | 28.3 | 0 | 106 | 72-130 | 0 | | | |
| Perfluorobutanoic Acid (PFBA) | 33.92 | 2.6 | 5.0 | 32 | 0 | 106 | 73-129 | 0 | | | |
| Perfluorodecanesulfonic Acid (| 36.78 | 1.4 | 5.0 | 30.8 | 0 | 119 | 53-142 | 0 | | | |
| Perfluorodecanoic Acid (PFDA) | 32.33 | 1.2 | 5.0 | 32 | 0 | 101 | 71-129 | 0 | | | |
| Perfluorododecanesulfonic Acid | 32.51 | 1.4 | 5.0 | 31 | 0 | 105 | 69-134 | 0 | | | |
| Perfluorododecanoic Acid (PFI) | 32.49 | 1.4 | 5.0 | 32 | 0 | 102 | 72-134 | 0 | | | |
| Perfluoroheptanesulfonic Acid | 31.57 | 0.57 | 5.0 | 30.5 | 0 | 104 | 69-134 | 0 | | | |
| Perfluoroheptanoic Acid (PFH _t) | 33.67 | 0.44 | 5.0 | 32 | 0 | 105 | 72-130 | 0 | | | |
| Perfluorohexadecanoic Acid (F | 34.06 | 0.38 | 5.0 | 32 | 0 | 106 | 70-130 | 0 | | | |
| Perfluorohexanesulfonic Acid (| 30.71 | 0.37 | 5.0 | 29.1 | 0 | 106 | 68-131 | 0 | | | |
| Perfluorohexanoic Acid (PFHx) | 34.33 | 1.2 | 5.0 | 32 | 0 | 107 | 72-129 | 0 | | | |
| Perfluorononanesulfonic Acid (| 32.9 | 0.5 | 5.0 | 30.7 | 0 | 107 | 69-127 | 0 | | | |
| Perfluorononanoic Acid (PFNA) | 30.65 | 0.87 | 5.0 | 32 | 0 | 95.8 | 69-130 | 0 | | | |
| Perfluoroctadecanoic Acid (P | 31.67 | 0.65 | 5.0 | 32 | 0 | 99 | 70-130 | 0 | | | |
| Perfluoroctanesulfonamide (F | 34.8 | 0.71 | 5.0 | 32 | 0 | 109 | 67-137 | 0 | | | |
| Perfluoroctanesulfonic Acid (I | 30.59 | 0.89 | 2.0 | 29.7 | 0 | 103 | 65-140 | 0 | | | |
| Perfluoroctanoic Acid (PFOA) | 27.49 | 0.63 | 2.0 | 32 | 0 | 85.9 | 71-133 | 0 | | | |
| Perfluoropentanesulfonic Acid | 30.1 | 0.56 | 5.0 | 30 | 0 | 100 | 71-127 | 0 | | | |
| Perfluoropentanoic Acid (PFPt) | 33.4 | 1.3 | 5.0 | 32 | 0 | 104 | 72-129 | 0 | | | |
| Perfluorotetradecanoic Acid (F | 30.2 | 2.6 | 5.0 | 32 | 0 | 94.4 | 71-132 | 0 | | | |
| Perfluorotridecanoic Acid (PTI) | 36.44 | 0.77 | 5.0 | 32 | 0 | 114 | 65-144 | 0 | | | |
| Perfluoroundecanoic Acid (PFI) | 35.43 | 0.97 | 5.0 | 32 | 0 | 111 | 69-133 | 0 | | | |
| N-ethylperfluoro-1-octanesulfo | 28.88 | 1.2 | 5.0 | 32 | 0 | 90.2 | 70-130 | 0 | | | |
| N-Ethylperfluoroctanesulfona | 29.85 | 0.63 | 5.0 | 32 | 0 | 93.3 | 61-135 | 0 | | | |
| N-Ethylperfluoroctanesulfona | 33.06 | 0.52 | 5.0 | 32 | 0 | 103 | 70-130 | 0 | | | |
| N-methylperfluoro-1-octanesul | 34.44 | 0.79 | 5.0 | 32 | 0 | 108 | 70-130 | 0 | | | |
| N-Methylperfluoroctanesulfuric | 32.68 | 0.64 | 5.0 | 32 | 0 | 102 | 65-136 | 0 | | | |
| N-Methylperfluoroctanesulfuric | 33.32 | 0.48 | 5.0 | 32 | 0 | 104 | 68-141 | 0 | | | |
| Hexafluoropropylene oxide din | 38.071 | 1.2 | 5.0 | 32 | 0 | 119 | 70-130 | 0 | | | |
| 4,8-Dioxa-3H-perfluorononano | 30.58 | 0.56 | 5.0 | 30.1 | 0 | 102 | 70-130 | 0 | | | |
| 11CI-Pf3OUdS | 31.09 | 0.47 | 5.0 | 30.1 | 0 | 103 | 70-130 | 0 | | | |
| 9CI-PF3ONS | 27.49 | 0.45 | 5.0 | 29.8 | 0 | 92.3 | 70-130 | 0 | | | |
| Surr: 13C2-FtS 4:2 | 100.9 | 0 | 0 | 149.4 | 0 | 67.6 | 50-150 | 0 | | | |
| Surr: 13C2-FtS 6:2 | 139.8 | 0 | 0 | 152 | 0 | 92 | 50-150 | 0 | | | |
| Surr: 13C2-FtS 8:2 | 130.2 | 0 | 0 | 153.3 | 0 | 85 | 50-150 | 0 | | | |
| Surr: 13C2-PFDA | 125 | 0 | 0 | 160 | 0 | 78.1 | 50-150 | 0 | | | |
| Surr: 13C2-PFDoA | 154.4 | 0 | 0 | 160 | 0 | 96.5 | 50-150 | 0 | | | |
| Surr: 13C2-PFHxA | 113.6 | 0 | 0 | 160 | 0 | 71 | 50-150 | 0 | | | |
| Surr: 13C2-PFHxDA | 131.4 | 0 | 0 | 160 | 0 | 82.1 | 50-150 | 0 | | | |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: 182387 | Instrument ID LCMS1 | Method: E537 Mod | | | | | | |
|-------------------------------|----------------------------|-------------------------|---|-------|---|------|--------|---|
| Surr: 13C2-PFTeA | 123.4 | 0 | 0 | 160 | 0 | 77.1 | 50-150 | 0 |
| Surr: 13C2-PFUnA | 140.1 | 0 | 0 | 160 | 0 | 87.6 | 50-150 | 0 |
| Surr: 13C3-HFPO-DA | 110.8 | 0 | 0 | 160 | 0 | 69.3 | 50-150 | 0 |
| Surr: 13C3-PFBS | 107.4 | 0 | 0 | 148.8 | 0 | 72.2 | 50-150 | 0 |
| Surr: 13C4-PFBA | 113.1 | 0 | 0 | 160 | 0 | 70.7 | 50-150 | 0 |
| Surr: 13C4-PFH _p A | 127.8 | 0 | 0 | 160 | 0 | 79.8 | 50-150 | 0 |
| Surr: 13C4-PFOA | 129.6 | 0 | 0 | 160 | 0 | 81 | 50-150 | 0 |
| Surr: 13C4-PFOS | 113.9 | 0 | 0 | 152.8 | 0 | 74.6 | 50-150 | 0 |
| Surr: 13C5-PFNA | 121.4 | 0 | 0 | 160 | 0 | 75.9 | 50-150 | 0 |
| Surr: 13C5-PFP _e A | 111.3 | 0 | 0 | 160 | 0 | 69.5 | 50-150 | 0 |
| Surr: 13C8-FOSA | 105.8 | 0 | 0 | 160 | 0 | 66.1 | 50-150 | 0 |
| Surr: 18O2-PFH _x S | 118.1 | 0 | 0 | 151.2 | 0 | 78.1 | 50-150 | 0 |
| Surr: d5-N-EtFOSA | 87.08 | 0 | 0 | 160 | 0 | 54.4 | 50-150 | 0 |
| Surr: d5-N-EtFOSAA | 168.3 | 0 | 0 | 160 | 0 | 105 | 50-150 | 0 |
| Surr: d9-N-EtFOSE | 113.7 | 0 | 0 | 160 | 0 | 71.1 | 50-150 | 0 |
| Surr: d3-N-MeFOSA | 82.78 | 0 | 0 | 160 | 0 | 51.7 | 50-150 | 0 |
| Surr: d3-N-MeFOSAA | 132.1 | 0 | 0 | 160 | 0 | 82.5 | 50-150 | 0 |
| Surr: d7-N-MeFOSE | 128.5 | 0 | 0 | 160 | 0 | 80.3 | 50-150 | 0 |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

Batch ID: **182387** Instrument ID **LCMS1** Method: **E537 Mod**

| MS | Sample ID: 21081457-01BMS | | | | Units: ng/L | | Analysis Date: 8/23/2021 06:28 PM | | | |
|---------------------------------|----------------------------------|------|-----|-----------------------|--------------------|-----------------------------|--|---------------|----------------|------|
| Client ID: | Run ID: LCMS1_210823B | | | SeqNo: 7690007 | | Prep Date: 8/23/2021 | | DF: 1 | | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | RPD %RPD Limit | Qual |
| Fluorotelomer Sulphonic Acid | 30.52 | 0.89 | 4.7 | 28.31 | 0 | 108 | 63-143 | 0 | | |
| Fluorotelomer Sulphonic Acid | 30.45 | 0.63 | 4.7 | 28.69 | 0 | 106 | 64-140 | 0 | | |
| Fluorotelomer Sulphonic Acid | 27.57 | 1.1 | 4.7 | 29.07 | 0 | 94.8 | 67-138 | 0 | | |
| Fluorotelomer Sulphonic Acid | 27.08 | 0.86 | 4.7 | 29.17 | 0 | 92.8 | 40-160 | 0 | | |
| Perfluorobutanesulfonic Acid (| 33.65 | 0.33 | 4.7 | 26.8 | 7.517 | 97.5 | 72-130 | 0 | | |
| Perfluorobutanoic Acid (PFBA) | 19.9 | 2.5 | 4.7 | 30.3 | 0 | 65.7 | 73-129 | 0 | | S |
| Perfluorodecanesulfonic Acid (| 27.6 | 1.3 | 4.7 | 29.17 | 0 | 94.6 | 53-142 | 0 | | |
| Perfluorodecanoic Acid (PFDA) | 31.5 | 1.2 | 4.7 | 30.3 | 1.445 | 99.2 | 71-129 | 0 | | |
| Perfluorododecanesulfonic Acid | 28.69 | 1.4 | 4.7 | 29.36 | 0 | 97.7 | 69-134 | 0 | | |
| Perfluorododecanoic Acid (PFI) | 29.31 | 1.4 | 4.7 | 30.3 | 3.385 | 95.1 | 72-134 | 0 | | |
| Perfluoroheptanesulfonic Acid | 27.94 | 0.54 | 4.7 | 28.88 | 0 | 96.7 | 69-134 | 0 | | |
| Perfluoroheptanoic Acid (PFH) | 32.21 | 0.42 | 4.7 | 30.3 | | | | | | |
| Perfluorohexadecanoic Acid (F | 31.02 | 0.36 | 4.7 | 30.3 | 0 | 102 | 70-130 | 0 | | |
| Perfluorohexanesulfonic Acid (| 29.28 | 0.35 | 4.7 | 27.56 | 1.768 | 99.9 | 68-131 | 0 | | |
| Perfluorohexanoic Acid (PFHx) | 32.03 | 1.1 | 4.7 | 30.3 | 3.287 | 94.9 | 72-129 | 0 | | |
| Perfluorononanesulfonic Acid (| 28.96 | 0.47 | 4.7 | 29.07 | 0 | 99.6 | 69-127 | 0 | | |
| Perfluorononanoic Acid (PFNA) | 30.5 | 0.82 | 4.7 | 30.3 | 1.733 | 94.9 | 69-130 | 0 | | |
| Perfluoroctadecanoic Acid (P | 29.2 | 0.61 | 4.7 | 30.3 | 0 | 96.4 | 70-130 | 0 | | |
| Perfluoroctanesulfonamide (F | 34.64 | 0.67 | 4.7 | 30.3 | 0 | 114 | 67-137 | 0 | | |
| Perfluoroctanesulfonic Acid (I | 71.65 | 0.84 | 1.9 | 28.12 | 44.61 | 96.1 | 65-140 | 0 | | |
| Perfluoroctanoic Acid (PFOA) | 29.58 | 0.6 | 1.9 | 30.3 | 4.586 | 82.5 | 71-133 | 0 | | |
| Perfluoropentanesulfonic Acid | 24.97 | 0.53 | 4.7 | 28.41 | 1.38 | 83 | 71-127 | 0 | | |
| Perfluoropentanoic Acid (PFP) | 28.95 | 1.2 | 4.7 | 30.3 | 2.572 | 87.1 | 72-129 | 0 | | |
| Perfluorotetradecanoic Acid (F | 28.02 | 2.5 | 4.7 | 30.3 | 0 | 92.5 | 71-132 | 0 | | |
| Perfluorotridecanoic Acid (PTI) | 35.41 | 0.73 | 4.7 | 30.3 | 0 | 117 | 65-144 | 0 | | |
| Perfluoroundecanoic Acid (PFI) | 32.56 | 0.92 | 4.7 | 30.3 | 0 | 107 | 69-133 | 0 | | |
| N-ethylperfluoro-1-octanesulfo | 28.85 | 1.1 | 4.7 | 30.3 | 0 | 95.2 | 70-130 | 0 | | |
| N-Ethylperfluoroctanesulfona | 26.12 | 0.59 | 4.7 | 30.3 | 0 | 86.2 | 61-135 | 0 | | |
| N-Ethylperfluoroctanesulfona | 33.52 | 0.49 | 4.7 | 30.3 | 0 | 111 | 70-130 | 0 | | |
| N-methylperfluoro-1-octanesul | 39.27 | 0.75 | 4.7 | 30.3 | 0 | 130 | 70-130 | 0 | | |
| N-Methylperfluoroctanesulfur | 31.64 | 0.61 | 4.7 | 30.3 | 0 | 104 | 65-136 | 0 | | |
| N-Methylperfluoroctanesulfur | 30.54 | 0.46 | 4.7 | 30.3 | 0 | 101 | 68-141 | 0 | | |
| Hexafluoropropylene oxide din | 29.54 | 1.1 | 4.7 | 30.3 | 0 | 97.5 | 70-130 | 0 | | |
| 4,8-Dioxa-3H-perfluorononano | 22.24 | 0.53 | 4.7 | 28.5 | 0 | 78 | 70-130 | 0 | | |
| 11CI-Pf3OUDS | 27.45 | 0.44 | 4.7 | 28.5 | 0 | 96.3 | 70-130 | 0 | | |
| 9CI-PF3ONS | 24.88 | 0.42 | 4.7 | 28.22 | 0 | 88.2 | 70-130 | 0 | | |
| Surr: 13C2-FtS 4:2 | 422.3 | 0 | 0 | 141.5 | 0 | 298 | 50-150 | 0 | | S |
| Surr: 13C2-FtS 6:2 | 560.6 | 0 | 0 | 143.9 | 0 | 389 | 50-150 | 0 | | S |
| Surr: 13C2-FtS 8:2 | 489.6 | 0 | 0 | 145.2 | 0 | 337 | 50-150 | 0 | | S |
| Surr: 13C2-PFDA | 113.2 | 0 | 0 | 151.5 | 0 | 74.7 | 50-150 | 0 | | |
| Surr: 13C2-PFDoA | 131 | 0 | 0 | 151.5 | 0 | 86.4 | 50-150 | 0 | | |
| Surr: 13C2-PFHxA | 72.12 | 0 | 0 | 151.5 | 0 | 47.6 | 50-150 | 0 | | S |
| Surr: 13C2-PFHxDA | 102.4 | 0 | 0 | 151.5 | 0 | 67.6 | 50-150 | 0 | | |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: 182387 | Instrument ID LCMS1 | Method: E537 Mod | | | | | | |
|-------------------------------|----------------------------|-------------------------|---|-------|---|------|--------|---|
| Surr: 13C2-PFTeA | 92.35 | 0 | 0 | 151.5 | 0 | 61 | 50-150 | 0 |
| Surr: 13C2-PFUnA | 138.9 | 0 | 0 | 151.5 | 0 | 91.7 | 50-150 | 0 |
| Surr: 13C3-HFPO-DA | 79.18 | 0 | 0 | 151.5 | 0 | 52.3 | 50-150 | 0 |
| Surr: 13C3-PFBS | 86.92 | 0 | 0 | 140.9 | 0 | 61.7 | 50-150 | 0 |
| Surr: 13C4-PFBA | 86.43 | 0 | 0 | 151.5 | 0 | 57 | 50-150 | 0 |
| Surr: 13C4-PFH _p A | 81.01 | 0 | 0 | 151.5 | 0 | 53.5 | 50-150 | 0 |
| Surr: 13C4-PFOA | 93.92 | 0 | 0 | 151.5 | 0 | 62 | 50-150 | 0 |
| Surr: 13C4-PFOS | 86.19 | 0 | 0 | 144.7 | 0 | 59.6 | 50-150 | 0 |
| Surr: 13C5-PFNA | 98.43 | 0 | 0 | 151.5 | 0 | 65 | 50-150 | 0 |
| Surr: 13C5-PFPeA | 78.01 | 0 | 0 | 151.5 | 0 | 51.5 | 50-150 | 0 |
| Surr: 13C8-FOSA | 81.43 | 0 | 0 | 151.5 | 0 | 53.7 | 50-150 | 0 |
| Surr: 18O2-PFH _x S | 99.98 | 0 | 0 | 143.2 | 0 | 69.8 | 50-150 | 0 |
| Surr: d5-N-EtFOSA | 62.83 | 0 | 0 | 151.5 | 0 | 41.5 | 50-150 | 0 |
| Surr: d5-N-EtFOSAA | 200.3 | 0 | 0 | 151.5 | 0 | 132 | 50-150 | 0 |
| Surr: d9-N-EtFOSE | 72.18 | 0 | 0 | 151.5 | 0 | 47.6 | 50-150 | 0 |
| Surr: d3-N-MeFOSA | 63.45 | 0 | 0 | 151.5 | 0 | 41.9 | 50-150 | 0 |
| Surr: d3-N-MeFOSAA | 143.7 | 0 | 0 | 151.5 | 0 | 94.9 | 50-150 | 0 |
| Surr: d7-N-MeFOSE | 85.53 | 0 | 0 | 151.5 | 0 | 56.5 | 50-150 | 0 |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

Batch ID: **182387** Instrument ID **LCMS1** Method: **E537 Mod**

| MSD | | Sample ID: 21081457-01BMSD | | | | Units: ng/L | | Analysis Date: 8/23/2021 06:38 PM | | | |
|---------------------------------|--------|-----------------------------------|-----|---------|-----------------------|--------------------|-----------------------------|--|--------------|-----------|------|
| Client ID: | | Run ID: LCMS1_210823B | | | SeqNo: 7690008 | | Prep Date: 8/23/2021 | | DF: 1 | | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | %RPD | RPD Limit | Qual |
| Fluorotelomer Sulphonic Acid | 35.13 | 0.96 | 5.1 | 30.51 | 0 | 115 | 63-143 | 30.52 | 14 | 30 | |
| Fluorotelomer Sulphonic Acid | 32.57 | 0.68 | 5.1 | 30.92 | 0 | 105 | 64-140 | 30.45 | 6.72 | 30 | |
| Fluorotelomer Sulphonic Acid | 36.33 | 1.2 | 5.1 | 31.33 | 0 | 116 | 67-138 | 27.57 | 27.4 | 30 | |
| Fluorotelomer Sulphonic Acid | 32.99 | 0.92 | 5.1 | 31.43 | 0 | 105 | 40-160 | 27.08 | 19.7 | 30 | |
| Perfluorobutanesulfonic Acid (| 37.03 | 0.36 | 5.1 | 28.88 | 7.517 | 102 | 72-130 | 33.65 | 9.55 | 30 | |
| Perfluorobutyric Acid (PFBA) | 22.85 | 2.7 | 5.1 | 32.65 | 0 | 70 | 73-129 | 19.9 | 13.8 | 30 | S |
| Perfluorodecanesulfonic Acid | 35.4 | 1.4 | 5.1 | 31.43 | 0 | 113 | 53-142 | 27.6 | 24.8 | 30 | |
| Perfluorodecanoic Acid (PFDA) | 34.08 | 1.3 | 5.1 | 32.65 | 1.445 | 100 | 71-129 | 31.5 | 7.88 | 30 | |
| Perfluorododecanesulfonic Acid | 31.78 | 1.5 | 5.1 | 31.63 | 0 | 100 | 69-134 | 28.69 | 10.2 | 30 | |
| Perfluorododecanoic Acid (PFI) | 30.76 | 1.5 | 5.1 | 32.65 | 0 | 94.2 | 72-134 | 29.31 | 4.83 | 30 | |
| Perfluoroheptanesulfonic Acid | 29.66 | 0.58 | 5.1 | 31.12 | 0 | 95.3 | 69-134 | 27.94 | 5.96 | 30 | |
| Perfluoroheptanoic Acid (PFH) | 37.5 | 0.45 | 5.1 | 32.65 | 3.385 | 104 | 72-130 | 32.21 | 15.2 | 30 | |
| Perfluorohexadecanoic Acid (F | 34.82 | 0.39 | 5.1 | 32.65 | 0 | 107 | 70-130 | 31.02 | 11.5 | 30 | |
| Perfluorohexanesulfonic Acid | 36.18 | 0.38 | 5.1 | 29.69 | 1.768 | 116 | 68-131 | 29.28 | 21.1 | 30 | |
| Perfluorohexanoic Acid (PFHx) | 34.11 | 1.2 | 5.1 | 32.65 | 3.287 | 94.4 | 72-129 | 32.03 | 6.29 | 30 | |
| Perfluorononanesulfonic Acid | 32.17 | 0.51 | 5.1 | 31.33 | 0 | 103 | 69-127 | 28.96 | 10.5 | 30 | |
| Perfluorononanoic Acid (PFNA) | 31.13 | 0.89 | 5.1 | 32.65 | 1.733 | 90 | 69-130 | 30.5 | 2.05 | 30 | |
| Perfluoroctadecanoic Acid (P | 31.28 | 0.66 | 5.1 | 32.65 | 0 | 95.8 | 70-130 | 29.2 | 6.87 | 30 | |
| Perfluoroctanesulfonamide (F | 37.13 | 0.73 | 5.1 | 32.65 | 0 | 114 | 67-137 | 34.64 | 6.94 | 30 | |
| Perfluoroctanesulfonic Acid (I | 71.1 | 0.91 | 2.0 | 30.31 | 44.61 | 87.4 | 65-140 | 71.65 | 0.766 | 30 | |
| Perfluoroctanoic Acid (PFOA) | 29.51 | 0.64 | 2.0 | 32.65 | 4.586 | 76.3 | 71-133 | 29.58 | 0.259 | 30 | |
| Perfluoropentanesulfonic Acid | 28.89 | 0.57 | 5.1 | 30.61 | 1.38 | 89.9 | 71-127 | 24.97 | 14.6 | 30 | |
| Perfluoropentanoic Acid (PFPe) | 32.98 | 1.3 | 5.1 | 32.65 | 2.572 | 93.1 | 72-129 | 28.95 | 13 | 30 | |
| Perfluorotetradecanoic Acid (F | 31.07 | 2.7 | 5.1 | 32.65 | 0 | 95.1 | 71-132 | 28.02 | 10.3 | 30 | |
| Perfluorotridecanoic Acid (PTI) | 36.38 | 0.79 | 5.1 | 32.65 | 0 | 111 | 65-144 | 35.41 | 2.72 | 30 | |
| Perfluoroundecanoic Acid (PFI) | 34.08 | 0.99 | 5.1 | 32.65 | 0 | 104 | 69-133 | 32.56 | 4.56 | 30 | |
| N-ethylperfluoro-1-octanesulfo | 29.91 | 1.2 | 5.1 | 32.65 | 0 | 91.6 | 70-130 | 28.85 | 3.62 | 30 | |
| N-Ethylperfluoroctanesulfona | 27.71 | 0.64 | 5.1 | 32.65 | 0 | 84.9 | 61-135 | 26.12 | 5.91 | 30 | |
| N-Ethylperfluoroctanesulfona | 33.8 | 0.53 | 5.1 | 32.65 | 0 | 104 | 70-130 | 33.52 | 0.845 | 30 | |
| N-methylperfluoro-1-octanesul | 38.21 | 0.81 | 5.1 | 32.65 | 0 | 117 | 70-130 | 39.27 | 2.73 | 30 | |
| N-Methylperfluoroctanesulfuric | 34.98 | 0.66 | 5.1 | 32.65 | 0 | 107 | 65-136 | 31.64 | 10 | 30 | |
| N-Methylperfluoroctanesulfuric | 31.95 | 0.49 | 5.1 | 32.65 | 0 | 97.8 | 68-141 | 30.54 | 4.52 | 30 | |
| Hexafluoropropylene oxide din | 38.271 | 1.2 | 5.1 | 32.65 | 0 | 117 | 70-130 | 29.54 | 25.7 | 30 | |
| 4,8-Dioxa-3H-perfluorononano | 25.3 | 0.57 | 5.1 | 30.71 | 0 | 82.4 | 70-130 | 22.24 | 12.9 | 30 | |
| 11CI-Pf3OUdS | 30.2 | 0.48 | 5.1 | 30.71 | 0 | 98.3 | 70-130 | 27.45 | 9.55 | 30 | |
| 9CI-PF3ONS | 29.19 | 0.46 | 5.1 | 30.41 | 0 | 96 | 70-130 | 24.88 | 15.9 | 30 | |
| Surr: 13C2-FtS 4:2 | 390.2 | 0 | 0 | 152.5 | 0 | 256 | 50-150 | 422.3 | 7.9 | 30 | S |
| Surr: 13C2-FtS 6:2 | 590.1 | 0 | 0 | 155.1 | 0 | 380 | 50-150 | 560.6 | 5.12 | 30 | S |
| Surr: 13C2-FtS 8:2 | 468.5 | 0 | 0 | 156.4 | 0 | 300 | 50-150 | 489.6 | 4.42 | 30 | S |
| Surr: 13C2-PFDA | 111.6 | 0 | 0 | 163.3 | 0 | 68.4 | 50-150 | 113.2 | 1.44 | 30 | |
| Surr: 13C2-PFDoA | 145.2 | 0 | 0 | 163.3 | 0 | 88.9 | 50-150 | 131 | 10.3 | 30 | |
| Surr: 13C2-PFHxA | 78.58 | 0 | 0 | 163.3 | 0 | 48.1 | 50-150 | 72.12 | 8.58 | 30 | |
| Surr: 13C2-PFHxDA | 106.1 | 0 | 0 | 163.3 | 0 | 65 | 50-150 | 102.4 | 3.54 | 30 | |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: 182387 | Instrument ID LCMS1 | Method: E537 Mod | | | | | | | | |
|-------------------------------|----------------------------|-------------------------|---|-------|---|------|--------|-------|-------|----|
| Surr: 13C2-PFTeA | 96.47 | 0 | 0 | 163.3 | 0 | 59.1 | 50-150 | 92.35 | 4.37 | 30 |
| Surr: 13C2-PFUnA | 144.9 | 0 | 0 | 163.3 | 0 | 88.8 | 50-150 | 138.9 | 4.21 | 30 |
| Surr: 13C3-HFPO-DA | 76.02 | 0 | 0 | 163.3 | 0 | 46.6 | 50-150 | 79.18 | 4.08 | 30 |
| Surr: 13C3-PFBS | 87.02 | 0 | 0 | 151.8 | 0 | 57.3 | 50-150 | 86.92 | 0.121 | 30 |
| Surr: 13C4-PFBA | 87.15 | 0 | 0 | 163.3 | 0 | 53.4 | 50-150 | 86.43 | 0.834 | 30 |
| Surr: 13C4-PFH _p A | 77.17 | 0 | 0 | 163.3 | 0 | 47.3 | 50-150 | 81.01 | 4.86 | 30 |
| Surr: 13C4-PFOA | 102.3 | 0 | 0 | 163.3 | 0 | 62.6 | 50-150 | 93.92 | 8.52 | 30 |
| Surr: 13C4-PFOS | 88.92 | 0 | 0 | 155.9 | 0 | 57 | 50-150 | 86.19 | 3.12 | 30 |
| Surr: 13C5-PFNA | 110.7 | 0 | 0 | 163.3 | 0 | 67.8 | 50-150 | 98.43 | 11.7 | 30 |
| Surr: 13C5-PFP _e A | 78.39 | 0 | 0 | 163.3 | 0 | 48 | 50-150 | 78.01 | 0.491 | 30 |
| Surr: 13C8-FOSA | 92.67 | 0 | 0 | 163.3 | 0 | 56.8 | 50-150 | 81.43 | 12.9 | 30 |
| Surr: 18O2-PFH _x S | 103.8 | 0 | 0 | 154.3 | 0 | 67.3 | 50-150 | 99.98 | 3.78 | 30 |
| Surr: d5-N-EtFOSA | 81.57 | 0 | 0 | 163.3 | 0 | 50 | 50-150 | 62.83 | 26 | 30 |
| Surr: d5-N-EtFOSAA | 223.9 | 0 | 0 | 163.3 | 0 | 137 | 50-150 | 200.3 | 11.1 | 30 |
| Surr: d9-N-EtFOSE | 84.93 | 0 | 0 | 163.3 | 0 | 52 | 50-150 | 72.18 | 16.2 | 30 |
| Surr: d3-N-MeFOSA | 84.82 | 0 | 0 | 163.3 | 0 | 52 | 50-150 | 63.45 | 28.8 | 30 |
| Surr: d3-N-MeFOSAA | 149.1 | 0 | 0 | 163.3 | 0 | 91.3 | 50-150 | 143.7 | 3.66 | 30 |
| Surr: d7-N-MeFOSE | 99.27 | 0 | 0 | 163.3 | 0 | 60.8 | 50-150 | 85.53 | 14.9 | 30 |

The following samples were analyzed in this batch:

21081889-01B 21081889-09A

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

Batch ID: R325427B Instrument ID VMS8 Method: SW8260C

| MBLK | Sample ID: 8V-BLKW1-210826-R325427B | | | Units: µg/L | | Analysis Date: 8/27/2021 12:10 AM | | | | | |
|-----------------------------|-------------------------------------|------|-----|----------------|---------------|-----------------------------------|---------------|---------------|----------|-----------|------|
| Client ID: | Run ID: VMS8_210826B | | | SeqNo: 7701424 | | Prep Date: | | DF: 1 | | | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | RPD %RPD | RPD Limit | Qual |
| 1,1,1,2-Tetrachloroethane | U | 0.38 | | 1.3 | | | | | | | |
| 1,1,1-Trichloroethane | U | 0.46 | | 1.5 | | | | | | | |
| 1,1,2,2-Tetrachloroethane | U | 0.4 | | 1.3 | | | | | | | |
| 1,1,2-Trichloroethane | U | 0.46 | | 1.5 | | | | | | | |
| 1,1-Dichloroethane | U | 0.44 | | 1.5 | | | | | | | |
| 1,1-Dichloroethene | U | 0.4 | | 1.4 | | | | | | | |
| 1,1-Dichloropropene | U | 0.37 | | 1.2 | | | | | | | |
| 1,2,3-Trichlorobenzene | U | 0.42 | | 1.4 | | | | | | | |
| 1,2,3-Trichloropropane | U | 0.4 | | 1.3 | | | | | | | |
| 1,2,4-Trichlorobenzene | U | 0.45 | | 1.5 | | | | | | | |
| 1,2,4-Trimethylbenzene | U | 0.45 | | 1.5 | | | | | | | |
| 1,2-Dibromo-3-chloropropane | U | 0.43 | | 1.4 | | | | | | | |
| 1,2-Dibromoethane | U | 0.41 | | 1.4 | | | | | | | |
| 1,2-Dichlorobenzene | U | 0.32 | | 1.1 | | | | | | | |
| 1,2-Dichloroethane | U | 0.44 | | 1.4 | | | | | | | |
| 1,2-Dichloropropene | U | 0.48 | | 1.6 | | | | | | | |
| 1,3,5-Trimethylbenzene | U | 0.65 | | 2.2 | | | | | | | |
| 1,3-Dichlorobenzene | U | 0.33 | | 1.1 | | | | | | | |
| 1,3-Dichloropropane | U | 0.4 | | 1.3 | | | | | | | |
| 1,4-Dichlorobenzene | U | 0.35 | | 1.2 | | | | | | | |
| 2,2-Dichloropropane | U | 0.52 | | 1.7 | | | | | | | |
| 2-Butanone | U | 0.52 | | 1.7 | | | | | | | |
| 2-Chlorotoluene | U | 0.36 | | 1.2 | | | | | | | |
| 4-Chlorotoluene | U | 0.31 | | 1.0 | | | | | | | |
| 4-Methyl-2-pentanone | U | 0.52 | | 1.7 | | | | | | | |
| Acetone | U | 6.2 | | 21 | | | | | | | |
| Benzene | U | 0.46 | | 1.5 | | | | | | | |
| Bromobenzene | U | 0.38 | | 1.3 | | | | | | | |
| Bromochloromethane | U | 0.45 | | 1.5 | | | | | | | |
| Bromodichloromethane | U | 0.49 | | 1.6 | | | | | | | |
| Bromoform | U | 0.56 | | 1.9 | | | | | | | |
| Bromomethane | U | 0.9 | | 3.0 | | | | | | | |
| Carbon tetrachloride | U | 0.4 | | 1.4 | | | | | | | |
| Chlorobenzene | U | 0.4 | | 1.3 | | | | | | | |
| Chloroethane | U | 0.68 | | 2.3 | | | | | | | |
| Chloroform | U | 0.46 | | 1.5 | | | | | | | |
| Chloromethane | U | 0.83 | | 2.8 | | | | | | | |
| cis-1,2-Dichloroethene | U | 0.42 | | 1.4 | | | | | | | |
| cis-1,3-Dichloropropene | U | 0.57 | | 1.9 | | | | | | | |
| Dibromochloromethane | U | 0.4 | | 1.3 | | | | | | | |
| Dibromomethane | U | 0.65 | | 2.2 | | | | | | | |
| Dichlorodifluoromethane | U | 0.68 | | 2.3 | | | | | | | |
| Ethylbenzene | U | 0.34 | | 1.1 | | | | | | | |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: R325427B | Instrument ID VMS8 | Method: SW8260C | | | | | |
|-----------------------------|---------------------------|------------------------|------|----|---|------|--------|
| Hexachlorobutadiene | U | 0.56 | 1.9 | | | | |
| Isopropylbenzene | U | 0.35 | 1.2 | | | | |
| m,p-Xylene | U | 0.81 | 2.7 | | | | |
| Methyl tert-butyl ether | U | 0.45 | 1.5 | | | | |
| Methylene chloride | U | 0.86 | 2.9 | | | | |
| Naphthalene | U | 0.77 | 2.6 | | | | |
| n-Butylbenzene | U | 0.34 | 1.1 | | | | |
| n-Propylbenzene | U | 0.48 | 1.6 | | | | |
| o-Xylene | U | 0.31 | 1.0 | | | | |
| p-Isopropyltoluene | U | 0.26 | 0.88 | | | | |
| sec-Butylbenzene | U | 0.3 | 1.0 | | | | |
| Styrene | U | 0.33 | 1.1 | | | | |
| tert-Butylbenzene | U | 0.39 | 1.3 | | | | |
| Tetrachloroethene | U | 0.39 | 1.3 | | | | |
| Toluene | U | 0.45 | 1.5 | | | | |
| trans-1,2-Dichloroethene | U | 0.48 | 1.6 | | | | |
| trans-1,3-Dichloropropene | U | 0.38 | 2.7 | | | | |
| Trichloroethene | U | 0.43 | 1.4 | | | | |
| Trichlorofluoromethane | U | 0.52 | 1.7 | | | | |
| Vinyl chloride | U | 0.53 | 1.8 | | | | |
| Xylenes, Total | U | 0.81 | 4.4 | | | | |
| Surr: 1,2-Dichloroethane-d4 | 19.7 | 0 | 0 | 20 | 0 | 98.5 | 75-120 |
| Surr: 4-Bromofluorobenzene | 19.15 | 0 | 0 | 20 | 0 | 95.8 | 80-110 |
| Surr: Dibromofluoromethane | 20.25 | 0 | 0 | 20 | 0 | 101 | 85-115 |
| Surr: Toluene-d8 | 20.12 | 0 | 0 | 20 | 0 | 101 | 85-110 |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

QC Page: 10 of 25

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

Batch ID: **R325427B** Instrument ID **VMS8** Method: **SW8260C**

| LCS | Sample ID: 8V-LCSW1-210826-R325427B | | | | Units: µg/L | | Analysis Date: 8/26/2021 11:10 PM | | | |
|-----------------------------|--|------|-----|-----------------------|--------------------|------------|--|---------------|----------------|------|
| Client ID: | Run ID: VMS8_210826B | | | SeqNo: 7701422 | | Prep Date: | | DF: 1 | | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | RPD %RPD Limit | Qual |
| 1,1,1,2-Tetrachloroethane | 17.79 | 0.38 | 1.3 | 20 | 0 | 89 | 73-114 | 0 | 0 | |
| 1,1,1-Trichloroethane | 18.06 | 0.46 | 1.5 | 20 | 0 | 90.3 | 75-130 | 0 | 0 | |
| 1,1,2,2-Tetrachloroethane | 20.07 | 0.4 | 1.3 | 20 | 0 | 100 | 75-130 | 0 | 0 | |
| 1,1,2-Trichloroethane | 18.24 | 0.46 | 1.5 | 20 | 0 | 91.2 | 75-125 | 0 | 0 | |
| 1,1-Dichloroethane | 17 | 0.44 | 1.5 | 20 | 0 | 85 | 68-142 | 0 | 0 | |
| 1,1-Dichloroethene | 18.47 | 0.4 | 1.4 | 20 | 0 | 92.4 | 70-145 | 0 | 0 | |
| 1,1-Dichloropropene | 18.01 | 0.37 | 1.2 | 20 | 0 | 90 | 75-135 | 0 | 0 | |
| 1,2,3-Trichlorobenzene | 18.1 | 0.42 | 1.4 | 20 | 0 | 90.5 | 70-140 | 0 | 0 | |
| 1,2,3-Trichloropropane | 17.51 | 0.4 | 1.3 | 20 | 0 | 87.6 | 75-125 | 0 | 0 | |
| 1,2,4-Trichlorobenzene | 18.53 | 0.45 | 1.5 | 20 | 0 | 92.6 | 70-135 | 0 | 0 | |
| 1,2,4-Trimethylbenzene | 17.06 | 0.45 | 1.5 | 20 | 0 | 85.3 | 75-130 | 0 | 0 | |
| 1,2-Dibromo-3-chloropropane | 17.38 | 0.43 | 1.4 | 20 | 0 | 86.9 | 60-130 | 0 | 0 | |
| 1,2-Dibromoethane | 18.5 | 0.41 | 1.4 | 20 | 0 | 92.5 | 67-155 | 0 | 0 | |
| 1,2-Dichlorobenzene | 16.54 | 0.32 | 1.1 | 20 | 0 | 82.7 | 70-130 | 0 | 0 | |
| 1,2-Dichloroethane | 17.92 | 0.44 | 1.4 | 20 | 0 | 89.6 | 78-125 | 0 | 0 | |
| 1,2-Dichloropropane | 19.26 | 0.48 | 1.6 | 20 | 0 | 96.3 | 75-125 | 0 | 0 | |
| 1,3,5-Trimethylbenzene | 16.68 | 0.65 | 2.2 | 20 | 0 | 83.4 | 75-130 | 0 | 0 | |
| 1,3-Dichlorobenzene | 18.18 | 0.33 | 1.1 | 20 | 0 | 90.9 | 75-130 | 0 | 0 | |
| 1,3-Dichloropropane | 18.71 | 0.4 | 1.3 | 20 | 0 | 93.6 | 75-125 | 0 | 0 | |
| 1,4-Dichlorobenzene | 16.65 | 0.35 | 1.2 | 20 | 0 | 83.2 | 75-130 | 0 | 0 | |
| 2,2-Dichloropropane | 19.12 | 0.52 | 1.7 | 20 | 0 | 95.6 | 43-150 | 0 | 0 | |
| 2-Butanone | 19.42 | 0.52 | 1.7 | 20 | 0 | 97.1 | 55-150 | 0 | 0 | |
| 2-Chlorotoluene | 16.49 | 0.36 | 1.2 | 20 | 0 | 82.4 | 76-117 | 0 | 0 | |
| 4-Chlorotoluene | 16.4 | 0.31 | 1.0 | 20 | 0 | 82 | 80-125 | 0 | 0 | |
| 4-Methyl-2-pentanone | 18.49 | 0.52 | 1.7 | 20 | 0 | 92.4 | 77-178 | 0 | 0 | |
| Acetone | 19.19 | 6.2 | 21 | 20 | 0 | 96 | 60-160 | 0 | 0 | J |
| Benzene | 18.09 | 0.46 | 1.5 | 20 | 0 | 90.4 | 70-130 | 0 | 0 | |
| Bromobenzene | 18.05 | 0.38 | 1.3 | 20 | 0 | 90.2 | 80-125 | 0 | 0 | |
| Bromochloromethane | 17.28 | 0.45 | 1.5 | 20 | 0 | 86.4 | 72-141 | 0 | 0 | |
| Bromodichloromethane | 18.22 | 0.49 | 1.6 | 20 | 0 | 91.1 | 75-125 | 0 | 0 | |
| Bromoform | 14.95 | 0.56 | 1.9 | 20 | 0 | 74.8 | 60-125 | 0 | 0 | |
| Bromomethane | 19.24 | 0.9 | 3.0 | 20 | 0 | 96.2 | 30-185 | 0 | 0 | |
| Carbon tetrachloride | 16.62 | 0.4 | 1.4 | 20 | 0 | 83.1 | 65-140 | 0 | 0 | |
| Chlorobenzene | 18.37 | 0.4 | 1.3 | 20 | 0 | 91.8 | 80-120 | 0 | 0 | |
| Chloroethane | 17.62 | 0.68 | 2.3 | 20 | 0 | 88.1 | 31-172 | 0 | 0 | |
| Chloroform | 17.07 | 0.46 | 1.5 | 20 | 0 | 85.4 | 66-135 | 0 | 0 | |
| Chloromethane | 14.98 | 0.83 | 2.8 | 20 | 0 | 74.9 | 46-148 | 0 | 0 | |
| cis-1,2-Dichloroethene | 17.69 | 0.42 | 1.4 | 20 | 0 | 88.4 | 75-134 | 0 | 0 | |
| cis-1,3-Dichloropropene | 16.56 | 0.57 | 1.9 | 20 | 0 | 82.8 | 70-130 | 0 | 0 | |
| Dibromochloromethane | 16.35 | 0.4 | 1.3 | 20 | 0 | 81.8 | 60-115 | 0 | 0 | |
| Dibromomethane | 18.33 | 0.65 | 2.2 | 20 | 0 | 91.6 | 79-126 | 0 | 0 | |
| Dichlorodifluoromethane | 18.63 | 0.68 | 2.3 | 20 | 0 | 93.2 | 10-180 | 0 | 0 | |
| Ethylbenzene | 16.64 | 0.34 | 1.1 | 20 | 0 | 83.2 | 76-123 | 0 | 0 | |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: R325427B | Instrument ID VMS8 | Method: SW8260C | | | | | | |
|-----------------------------|---------------------------|------------------------|------|----|---|------|--------|---|
| Hexachlorobutadiene | 18.91 | 0.56 | 1.9 | 20 | 0 | 94.6 | 70-155 | 0 |
| Isopropylbenzene | 18.83 | 0.35 | 1.2 | 20 | 0 | 94.2 | 80-127 | 0 |
| m,p-Xylene | 33.59 | 0.81 | 2.7 | 40 | 0 | 84 | 75-130 | 0 |
| Methyl tert-butyl ether | 20.47 | 0.45 | 1.5 | 20 | 0 | 102 | 68-129 | 0 |
| Methylene chloride | 18.17 | 0.86 | 2.9 | 20 | 0 | 90.8 | 72-125 | 0 |
| Naphthalene | 17.07 | 0.77 | 2.6 | 20 | 0 | 85.4 | 55-160 | 0 |
| n-Butylbenzene | 16.19 | 0.34 | 1.1 | 20 | 0 | 81 | 75-145 | 0 |
| n-Propylbenzene | 16.73 | 0.48 | 1.6 | 20 | 0 | 83.6 | 76-116 | 0 |
| o-Xylene | 17.18 | 0.31 | 1.0 | 20 | 0 | 85.9 | 76-127 | 0 |
| p-Isopropyltoluene | 16.66 | 0.26 | 0.88 | 20 | 0 | 83.3 | 61-164 | 0 |
| sec-Butylbenzene | 17.27 | 0.3 | 1.0 | 20 | 0 | 86.4 | 80-134 | 0 |
| Styrene | 18.56 | 0.33 | 1.1 | 20 | 0 | 92.8 | 79-117 | 0 |
| tert-Butylbenzene | 17.38 | 0.39 | 1.3 | 20 | 0 | 86.9 | 70-130 | 0 |
| Tetrachloroethene | 17.69 | 0.39 | 1.3 | 20 | 0 | 88.4 | 68-166 | 0 |
| Toluene | 17.71 | 0.45 | 1.5 | 20 | 0 | 88.6 | 76-125 | 0 |
| trans-1,2-Dichloroethene | 18.08 | 0.48 | 1.6 | 20 | 0 | 90.4 | 80-140 | 0 |
| trans-1,3-Dichloropropene | 18.22 | 0.38 | 2.7 | 20 | 0 | 91.1 | 56-132 | 0 |
| Trichloroethene | 16.19 | 0.43 | 1.4 | 20 | 0 | 81 | 77-125 | 0 |
| Trichlorofluoromethane | 18.65 | 0.52 | 1.7 | 20 | 0 | 93.2 | 60-140 | 0 |
| Vinyl chloride | 18.22 | 0.53 | 1.8 | 20 | 0 | 91.1 | 50-136 | 0 |
| Xylenes, Total | 50.77 | 0.81 | 4.4 | 60 | 0 | 84.6 | 76-127 | 0 |
| Surr: 1,2-Dichloroethane-d4 | 19.36 | 0 | 0 | 20 | 0 | 96.8 | 75-120 | 0 |
| Surr: 4-Bromofluorobenzene | 20.2 | 0 | 0 | 20 | 0 | 101 | 80-110 | 0 |
| Surr: Dibromofluoromethane | 19.91 | 0 | 0 | 20 | 0 | 99.6 | 85-115 | 0 |
| Surr: Toluene-d8 | 20.13 | 0 | 0 | 20 | 0 | 101 | 85-110 | 0 |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

Batch ID: **R325427B** Instrument ID **VMS8** Method: **SW8260C**

| MS | Sample ID: 21081675-01A MS | | | | Units: µg/L | | Analysis Date: 8/27/2021 07:26 AM | | |
|-----------------------------|-----------------------------------|-----|-------|-----------------------|--------------------|------------|--|----------------|---------------------|
| Client ID: | Run ID: VMS8_210826B | | | SeqNo: 7701446 | | Prep Date: | | DF: 100 | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | RPD %RPD Limit Qual |
| 1,1,1,2-Tetrachloroethane | 1950 | 38 | 130 | 2000 | 0 | 97.5 | 73-114 | 0 | |
| 1,1,1-Trichloroethane | 1988 | 46 | 150 | 2000 | 0 | 99.4 | 75-130 | 0 | |
| 1,1,2,2-Tetrachloroethane | 2066 | 40 | 130 | 2000 | 0 | 103 | 75-130 | 0 | |
| 1,1,2-Trichloroethane | 1975 | 46 | 150 | 2000 | 0 | 98.8 | 75-125 | 0 | |
| 1,1-Dichloroethane | 1882 | 44 | 150 | 2000 | 0 | 94.1 | 68-142 | 0 | |
| 1,1-Dichloroethene | 2155 | 40 | 140 | 2000 | 0 | 108 | 70-145 | 0 | |
| 1,1-Dichloropropene | 1902 | 37 | 120 | 2000 | 0 | 95.1 | 75-135 | 0 | |
| 1,2,3-Trichlorobenzene | 1752 | 42 | 140 | 2000 | 0 | 87.6 | 70-140 | 0 | |
| 1,2,3-Trichloropropane | 1856 | 40 | 130 | 2000 | 0 | 92.8 | 75-125 | 0 | |
| 1,2,4-Trichlorobenzene | 1813 | 45 | 150 | 2000 | 0 | 90.6 | 70-135 | 0 | |
| 1,2,4-Trimethylbenzene | 1880 | 45 | 150 | 2000 | 0 | 94 | 75-130 | 0 | |
| 1,2-Dibromo-3-chloropropane | 1627 | 43 | 140 | 2000 | 0 | 81.4 | 60-130 | 0 | |
| 1,2-Dibromoethane | 1982 | 41 | 140 | 2000 | 0 | 99.1 | 67-155 | 0 | |
| 1,2-Dichlorobenzene | 1763 | 32 | 110 | 2000 | 0 | 88.2 | 70-130 | 0 | |
| 1,2-Dichloroethane | 1943 | 44 | 140 | 2000 | 0 | 97.2 | 78-125 | 0 | |
| 1,2-Dichloropropene | 2086 | 48 | 160 | 2000 | 0 | 104 | 75-125 | 0 | |
| 1,3,5-Trimethylbenzene | 1851 | 65 | 220 | 2000 | 0 | 92.6 | 75-130 | 0 | |
| 1,3-Dichlorobenzene | 1986 | 33 | 110 | 2000 | 0 | 99.3 | 75-130 | 0 | |
| 1,3-Dichloropropane | 1996 | 40 | 130 | 2000 | 0 | 99.8 | 75-125 | 0 | |
| 1,4-Dichlorobenzene | 1802 | 35 | 120 | 2000 | 0 | 90.1 | 75-130 | 0 | |
| 2,2-Dichloropropane | 1767 | 52 | 170 | 2000 | 0 | 88.4 | 43-150 | 0 | |
| 2-Butanone | 2041 | 52 | 170 | 2000 | 0 | 102 | 55-150 | 0 | |
| 2-Chlorotoluene | 1825 | 36 | 120 | 2000 | 0 | 91.2 | 76-117 | 0 | |
| 4-Chlorotoluene | 1854 | 31 | 100 | 2000 | 0 | 92.7 | 80-125 | 0 | |
| 4-Methyl-2-pentanone | 2664 | 52 | 170 | 2000 | 0 | 133 | 77-178 | 0 | |
| Acetone | 1929 | 620 | 2,100 | 2000 | 0 | 96.4 | 60-160 | 0 | J |
| Benzene | 1998 | 46 | 150 | 2000 | 0 | 99.9 | 70-130 | 0 | |
| Bromobenzene | 1965 | 38 | 130 | 2000 | 0 | 98.2 | 80-125 | 0 | |
| Bromochloromethane | 2107 | 45 | 150 | 2000 | 0 | 105 | 72-141 | 0 | |
| Bromodichloromethane | 1988 | 49 | 160 | 2000 | 0 | 99.4 | 75-125 | 0 | |
| Bromoform | 1601 | 56 | 190 | 2000 | 0 | 80 | 60-125 | 0 | |
| Bromomethane | 3326 | 90 | 300 | 2000 | 0 | 166 | 30-185 | 0 | |
| Carbon tetrachloride | 1935 | 40 | 140 | 2000 | 0 | 96.8 | 65-140 | 0 | |
| Chlorobenzene | 2010 | 40 | 130 | 2000 | 0 | 100 | 80-120 | 0 | |
| Chloroethane | 3797 | 68 | 230 | 2000 | 0 | 190 | 31-172 | 0 | S |
| Chloroform | 1918 | 46 | 150 | 2000 | 0 | 95.9 | 66-135 | 0 | |
| Chloromethane | 1412 | 83 | 280 | 2000 | 0 | 70.6 | 46-148 | 0 | |
| cis-1,2-Dichloroethene | 1912 | 42 | 140 | 2000 | 0 | 95.6 | 75-134 | 0 | |
| cis-1,3-Dichloropropene | 1801 | 57 | 190 | 2000 | 0 | 90 | 70-130 | 0 | |
| Dibromochloromethane | 1796 | 40 | 130 | 2000 | 0 | 89.8 | 60-115 | 0 | |
| Dibromomethane | 1976 | 65 | 220 | 2000 | 0 | 98.8 | 79-126 | 0 | |
| Dichlorodifluoromethane | 1753 | 68 | 230 | 2000 | 0 | 87.6 | 10-180 | 0 | |
| Ethylbenzene | 1851 | 34 | 110 | 2000 | 0 | 92.6 | 76-123 | 0 | |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: R325427B | Instrument ID VMS8 | Method: SW8260C | | | | | | |
|------------------------------------|---------------------------|------------------------|-----|------|----|------|--------|---|
| Hexachlorobutadiene | 2033 | 56 | 190 | 2000 | 0 | 102 | 70-155 | 0 |
| Isopropylbenzene | 2144 | 35 | 120 | 2000 | 0 | 107 | 80-127 | 0 |
| m,p-Xylene | 3812 | 81 | 270 | 4000 | 0 | 95.3 | 75-130 | 0 |
| Methyl tert-butyl ether | 2498 | 45 | 150 | 2000 | 0 | 125 | 68-129 | 0 |
| Methylene chloride | 1980 | 86 | 290 | 2000 | 0 | 99 | 72-125 | 0 |
| Naphthalene | 1687 | 77 | 260 | 2000 | 0 | 84.4 | 55-160 | 0 |
| n-Butylbenzene | 1788 | 34 | 110 | 2000 | 0 | 89.4 | 75-145 | 0 |
| n-Propylbenzene | 1877 | 48 | 160 | 2000 | 0 | 93.8 | 76-116 | 0 |
| o-Xylene | 1897 | 31 | 100 | 2000 | 0 | 94.8 | 76-127 | 0 |
| p-Isopropyltoluene | 1899 | 26 | 88 | 2000 | 0 | 95 | 61-164 | 0 |
| sec-Butylbenzene | 1954 | 30 | 100 | 2000 | 0 | 97.7 | 80-134 | 0 |
| Styrene | 2057 | 33 | 110 | 2000 | 0 | 103 | 79-117 | 0 |
| tert-Butylbenzene | 2000 | 39 | 130 | 2000 | 0 | 100 | 70-130 | 0 |
| Tetrachloroethene | 1881 | 39 | 130 | 2000 | 26 | 92.8 | 68-166 | 0 |
| Toluene | 1927 | 45 | 150 | 2000 | 0 | 96.4 | 76-125 | 0 |
| trans-1,2-Dichloroethene | 2044 | 48 | 160 | 2000 | 0 | 102 | 80-140 | 0 |
| trans-1,3-Dichloropropene | 1894 | 38 | 270 | 2000 | 0 | 94.7 | 56-132 | 0 |
| Trichloroethene | 1814 | 43 | 140 | 2000 | 0 | 90.7 | 77-125 | 0 |
| Trichlorofluoromethane | 2136 | 52 | 170 | 2000 | 0 | 107 | 60-140 | 0 |
| Vinyl chloride | 1872 | 53 | 180 | 2000 | 0 | 93.6 | 50-136 | 0 |
| Xylenes, Total | 5709 | 81 | 440 | 6000 | 0 | 95.2 | 76-127 | 0 |
| <i>Surr: 1,2-Dichloroethane-d4</i> | 1953 | 0 | 0 | 2000 | 0 | 97.6 | 75-120 | 0 |
| <i>Surr: 4-Bromofluorobenzene</i> | 2075 | 0 | 0 | 2000 | 0 | 104 | 80-110 | 0 |
| <i>Surr: Dibromofluoromethane</i> | 2026 | 0 | 0 | 2000 | 0 | 101 | 85-115 | 0 |
| <i>Surr: Toluene-d8</i> | 2013 | 0 | 0 | 2000 | 0 | 101 | 85-110 | 0 |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

Batch ID: **R325427B** Instrument ID **VMS8** Method: **SW8260C**

| MSD | | Sample ID: 21081675-01A MSD | | | | Units: µg/L | | Analysis Date: 8/27/2021 07:46 AM | | | |
|-----------------------------|--------|------------------------------------|-------|---------|-----------------------|--------------------|---------------|--|----------------|-----------|------|
| Client ID: | | Run ID: VMS8_210826B | | | SeqNo: 7701447 | | Prep Date: | | DF: 100 | | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | %RPD | RPD Limit | Qual |
| 1,1,1,2-Tetrachloroethane | 1926 | 38 | 130 | 2000 | 0 | 96.3 | 73-114 | 1950 | 1.24 | 30 | |
| 1,1,1-Trichloroethane | 2024 | 46 | 150 | 2000 | 0 | 101 | 75-130 | 1988 | 1.79 | 30 | |
| 1,1,2,2-Tetrachloroethane | 1982 | 40 | 130 | 2000 | 0 | 99.1 | 75-130 | 2066 | 4.15 | 30 | |
| 1,1,2-Trichloroethane | 1918 | 46 | 150 | 2000 | 0 | 95.9 | 75-125 | 1975 | 2.93 | 30 | |
| 1,1-Dichloroethane | 1902 | 44 | 150 | 2000 | 0 | 95.1 | 68-142 | 1882 | 1.06 | 30 | |
| 1,1-Dichloroethene | 2086 | 40 | 140 | 2000 | 0 | 104 | 70-145 | 2155 | 3.25 | 30 | |
| 1,1-Dichloropropene | 1974 | 37 | 120 | 2000 | 0 | 98.7 | 75-135 | 1902 | 3.72 | 30 | |
| 1,2,3-Trichlorobenzene | 1767 | 42 | 140 | 2000 | 0 | 88.4 | 70-140 | 1752 | 0.853 | 30 | |
| 1,2,3-Trichloropropane | 1745 | 40 | 130 | 2000 | 0 | 87.2 | 75-125 | 1856 | 6.16 | 30 | |
| 1,2,4-Trichlorobenzene | 1785 | 45 | 150 | 2000 | 0 | 89.2 | 70-135 | 1813 | 1.56 | 30 | |
| 1,2,4-Trimethylbenzene | 1805 | 45 | 150 | 2000 | 0 | 90.2 | 75-130 | 1880 | 4.07 | 30 | |
| 1,2-Dibromo-3-chloropropane | 1632 | 43 | 140 | 2000 | 0 | 81.6 | 60-130 | 1627 | 0.307 | 30 | |
| 1,2-Dibromoethane | 1952 | 41 | 140 | 2000 | 0 | 97.6 | 67-155 | 1982 | 1.53 | 30 | |
| 1,2-Dichlorobenzene | 1727 | 32 | 110 | 2000 | 0 | 86.4 | 70-130 | 1763 | 2.06 | 30 | |
| 1,2-Dichloroethane | 1925 | 44 | 140 | 2000 | 0 | 96.2 | 78-125 | 1943 | 0.931 | 30 | |
| 1,2-Dichloropropane | 2088 | 48 | 160 | 2000 | 0 | 104 | 75-125 | 2086 | 0.0958 | 30 | |
| 1,3,5-Trimethylbenzene | 1814 | 65 | 220 | 2000 | 0 | 90.7 | 75-130 | 1851 | 2.02 | 30 | |
| 1,3-Dichlorobenzene | 1946 | 33 | 110 | 2000 | 0 | 97.3 | 75-130 | 1986 | 2.03 | 30 | |
| 1,3-Dichloropropane | 1917 | 40 | 130 | 2000 | 0 | 95.8 | 75-125 | 1996 | 4.04 | 30 | |
| 1,4-Dichlorobenzene | 1799 | 35 | 120 | 2000 | 0 | 90 | 75-130 | 1802 | 0.167 | 30 | |
| 2,2-Dichloropropane | 1752 | 52 | 170 | 2000 | 0 | 87.6 | 43-150 | 1767 | 0.853 | 30 | |
| 2-Butanone | 2072 | 52 | 170 | 2000 | 0 | 104 | 55-150 | 2041 | 1.51 | 30 | |
| 2-Chlorotoluene | 1750 | 36 | 120 | 2000 | 0 | 87.5 | 76-117 | 1825 | 4.2 | 30 | |
| 4-Chlorotoluene | 1787 | 31 | 100 | 2000 | 0 | 89.4 | 80-125 | 1854 | 3.68 | 30 | |
| 4-Methyl-2-pentanone | 2669 | 52 | 170 | 2000 | 0 | 133 | 77-178 | 2664 | 0.188 | 30 | |
| Acetone | 1917 | 620 | 2,100 | 2000 | 0 | 95.8 | 60-160 | 1929 | 0 | 30 | J |
| Benzene | 1985 | 46 | 150 | 2000 | 0 | 99.2 | 70-130 | 1998 | 0.653 | 30 | |
| Bromobenzene | 1908 | 38 | 130 | 2000 | 0 | 95.4 | 80-125 | 1965 | 2.94 | 30 | |
| Bromochloromethane | 2070 | 45 | 150 | 2000 | 0 | 104 | 72-141 | 2107 | 1.77 | 30 | |
| Bromodichloromethane | 1999 | 49 | 160 | 2000 | 0 | 100 | 75-125 | 1988 | 0.552 | 30 | |
| Bromoform | 1596 | 56 | 190 | 2000 | 0 | 79.8 | 60-125 | 1601 | 0.313 | 30 | |
| Bromomethane | 3286 | 90 | 300 | 2000 | 0 | 164 | 30-185 | 3326 | 1.21 | 30 | |
| Carbon tetrachloride | 1936 | 40 | 140 | 2000 | 0 | 96.8 | 65-140 | 1935 | 0.0517 | 30 | |
| Chlorobenzene | 1959 | 40 | 130 | 2000 | 0 | 98 | 80-120 | 2010 | 2.57 | 30 | |
| Chloroethane | 3719 | 68 | 230 | 2000 | 0 | 186 | 31-172 | 3797 | 2.08 | 30 | S |
| Chloroform | 1868 | 46 | 150 | 2000 | 0 | 93.4 | 66-135 | 1918 | 2.64 | 30 | |
| Chloromethane | 1368 | 83 | 280 | 2000 | 0 | 68.4 | 46-148 | 1412 | 3.17 | 30 | |
| cis-1,2-Dichloroethene | 1905 | 42 | 140 | 2000 | 0 | 95.2 | 75-134 | 1912 | 0.367 | 30 | |
| cis-1,3-Dichloropropene | 1773 | 57 | 190 | 2000 | 0 | 88.6 | 70-130 | 1801 | 1.57 | 30 | |
| Dibromochloromethane | 1773 | 40 | 130 | 2000 | 0 | 88.6 | 60-115 | 1796 | 1.29 | 30 | |
| Dibromomethane | 1955 | 65 | 220 | 2000 | 0 | 97.8 | 79-126 | 1976 | 1.07 | 30 | |
| Dichlorodifluoromethane | 1648 | 68 | 230 | 2000 | 0 | 82.4 | 10-180 | 1753 | 6.17 | 30 | |
| Ethylbenzene | 1803 | 34 | 110 | 2000 | 0 | 90.2 | 76-123 | 1851 | 2.63 | 30 | |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

QC Page: 15 of 25

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: R325427B | Instrument ID VMS8 | Method: SW8260C | | | | | | | | |
|------------------------------------|---------------------------|------------------------|-----|------|----|------|--------|------|-------|----|
| Hexachlorobutadiene | 1948 | 56 | 190 | 2000 | 0 | 97.4 | 70-155 | 2033 | 4.27 | 30 |
| Isopropylbenzene | 2043 | 35 | 120 | 2000 | 0 | 102 | 80-127 | 2144 | 4.82 | 30 |
| m,p-Xylene | 3659 | 81 | 270 | 4000 | 0 | 91.5 | 75-130 | 3812 | 4.1 | 30 |
| Methyl tert-butyl ether | 2415 | 45 | 150 | 2000 | 0 | 121 | 68-129 | 2498 | 3.38 | 30 |
| Methylene chloride | 1956 | 86 | 290 | 2000 | 0 | 97.8 | 72-125 | 1980 | 1.22 | 30 |
| Naphthalene | 1684 | 77 | 260 | 2000 | 0 | 84.2 | 55-160 | 1687 | 0.178 | 30 |
| n-Butylbenzene | 1763 | 34 | 110 | 2000 | 0 | 88.2 | 75-145 | 1788 | 1.41 | 30 |
| n-Propylbenzene | 1787 | 48 | 160 | 2000 | 0 | 89.4 | 76-116 | 1877 | 4.91 | 30 |
| o-Xylene | 1809 | 31 | 100 | 2000 | 0 | 90.4 | 76-127 | 1897 | 4.75 | 30 |
| p-Isopropyltoluene | 1851 | 26 | 88 | 2000 | 0 | 92.6 | 61-164 | 1899 | 2.56 | 30 |
| sec-Butylbenzene | 1866 | 30 | 100 | 2000 | 0 | 93.3 | 80-134 | 1954 | 4.61 | 30 |
| Styrene | 1974 | 33 | 110 | 2000 | 0 | 98.7 | 79-117 | 2057 | 4.12 | 30 |
| tert-Butylbenzene | 1868 | 39 | 130 | 2000 | 0 | 93.4 | 70-130 | 2000 | 6.83 | 30 |
| Tetrachloroethene | 1923 | 39 | 130 | 2000 | 26 | 94.8 | 68-166 | 1881 | 2.21 | 30 |
| Toluene | 1885 | 45 | 150 | 2000 | 0 | 94.2 | 76-125 | 1927 | 2.2 | 30 |
| trans-1,2-Dichloroethene | 1970 | 48 | 160 | 2000 | 0 | 98.5 | 80-140 | 2044 | 3.69 | 30 |
| trans-1,3-Dichloropropene | 1848 | 38 | 270 | 2000 | 0 | 92.4 | 56-132 | 1894 | 2.46 | 30 |
| Trichloroethene | 1744 | 43 | 140 | 2000 | 0 | 87.2 | 77-125 | 1814 | 3.93 | 30 |
| Trichlorofluoromethane | 2190 | 52 | 170 | 2000 | 0 | 110 | 60-140 | 2136 | 2.5 | 30 |
| Vinyl chloride | 1847 | 53 | 180 | 2000 | 0 | 92.4 | 50-136 | 1872 | 1.34 | 30 |
| Xylenes, Total | 5468 | 81 | 440 | 6000 | 0 | 91.1 | 76-127 | 5709 | 4.31 | 30 |
| <i>Surr: 1,2-Dichloroethane-d4</i> | 1997 | 0 | 0 | 2000 | 0 | 99.8 | 75-120 | 1953 | 2.23 | 30 |
| <i>Surr: 4-Bromofluorobenzene</i> | 1915 | 0 | 0 | 2000 | 0 | 95.8 | 80-110 | 2075 | 8.02 | 30 |
| <i>Surr: Dibromofluoromethane</i> | 2062 | 0 | 0 | 2000 | 0 | 103 | 85-115 | 2026 | 1.76 | 30 |
| <i>Surr: Toluene-d8</i> | 1990 | 0 | 0 | 2000 | 0 | 99.5 | 85-110 | 2013 | 1.15 | 30 |

The following samples were analyzed in this batch:

| | | |
|--------------|--------------|--------------|
| 21081889-01A | 21081889-02A | 21081889-03A |
| 21081889-04A | 21081889-05A | 21081889-06A |
| 21081889-07A | 21081889-08A | 21081889-10A |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

QC Page: 16 of 25

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

Batch ID: R325443a Instrument ID VMS8 Method: SW8260C

| MBLK | Sample ID: 8V-BLKW1-210827-R325443a | | | Units: µg/L | | Analysis Date: 8/27/2021 02:43 PM | | | | | |
|-----------------------------|-------------------------------------|------|-----|----------------|---------------|-----------------------------------|---------------|---------------|------|-----------|------|
| Client ID: | Run ID: VMS8_210827A | | | SeqNo: 7704532 | | Prep Date: | | DF: 1 | | | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | %RPD | RPD Limit | Qual |
| 1,1,1,2-Tetrachloroethane | U | 0.38 | | 1.3 | | | | | | | |
| 1,1,1-Trichloroethane | U | 0.46 | | 1.5 | | | | | | | |
| 1,1,2,2-Tetrachloroethane | U | 0.4 | | 1.3 | | | | | | | |
| 1,1,2-Trichloroethane | U | 0.46 | | 1.5 | | | | | | | |
| 1,1-Dichloroethane | U | 0.44 | | 1.5 | | | | | | | |
| 1,1-Dichloroethene | U | 0.4 | | 1.4 | | | | | | | |
| 1,1-Dichloropropene | U | 0.37 | | 1.2 | | | | | | | |
| 1,2,3-Trichlorobenzene | U | 0.42 | | 1.4 | | | | | | | |
| 1,2,3-Trichloropropane | U | 0.4 | | 1.3 | | | | | | | |
| 1,2,4-Trichlorobenzene | U | 0.45 | | 1.5 | | | | | | | |
| 1,2,4-Trimethylbenzene | U | 0.45 | | 1.5 | | | | | | | |
| 1,2-Dibromo-3-chloropropane | U | 0.43 | | 1.4 | | | | | | | |
| 1,2-Dibromoethane | U | 0.41 | | 1.4 | | | | | | | |
| 1,2-Dichlorobenzene | U | 0.32 | | 1.1 | | | | | | | |
| 1,2-Dichloroethane | U | 0.44 | | 1.4 | | | | | | | |
| 1,2-Dichloropropane | U | 0.48 | | 1.6 | | | | | | | |
| 1,3,5-Trimethylbenzene | U | 0.65 | | 2.2 | | | | | | | |
| 1,3-Dichlorobenzene | U | 0.33 | | 1.1 | | | | | | | |
| 1,3-Dichloropropane | U | 0.4 | | 1.3 | | | | | | | |
| 1,4-Dichlorobenzene | U | 0.35 | | 1.2 | | | | | | | |
| 2,2-Dichloropropane | U | 0.52 | | 1.7 | | | | | | | |
| 2-Butanone | U | 0.52 | | 1.7 | | | | | | | |
| 2-Chlorotoluene | U | 0.36 | | 1.2 | | | | | | | |
| 4-Chlorotoluene | U | 0.31 | | 1.0 | | | | | | | |
| 4-Methyl-2-pentanone | U | 0.52 | | 1.7 | | | | | | | |
| Acetone | U | 6.2 | | 21 | | | | | | | |
| Benzene | U | 0.46 | | 1.5 | | | | | | | |
| Bromobenzene | U | 0.38 | | 1.3 | | | | | | | |
| Bromochloromethane | U | 0.45 | | 1.5 | | | | | | | |
| Bromodichloromethane | U | 0.49 | | 1.6 | | | | | | | |
| Bromoform | U | 0.56 | | 1.9 | | | | | | | |
| Bromomethane | U | 0.9 | | 3.0 | | | | | | | |
| Carbon tetrachloride | U | 0.4 | | 1.4 | | | | | | | |
| Chlorobenzene | U | 0.4 | | 1.3 | | | | | | | |
| Chloroethane | U | 0.68 | | 2.3 | | | | | | | |
| Chloroform | U | 0.46 | | 1.5 | | | | | | | |
| Chloromethane | U | 0.83 | | 2.8 | | | | | | | |
| cis-1,2-Dichloroethene | U | 0.42 | | 1.4 | | | | | | | |
| cis-1,3-Dichloropropene | U | 0.57 | | 1.9 | | | | | | | |
| Dibromochloromethane | U | 0.4 | | 1.3 | | | | | | | |
| Dibromomethane | U | 0.65 | | 2.2 | | | | | | | |
| Dichlorodifluoromethane | U | 0.68 | | 2.3 | | | | | | | |
| Ethylbenzene | U | 0.34 | | 1.1 | | | | | | | |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: R325443a | Instrument ID VMS8 | Method: SW8260C | | | | | |
|-----------------------------|--------------------|-----------------|------|----|---|------|--------|
| Hexachlorobutadiene | U | 0.56 | 1.9 | | | | |
| Isopropylbenzene | U | 0.35 | 1.2 | | | | |
| m,p-Xylene | U | 0.81 | 2.7 | | | | |
| Methyl tert-butyl ether | U | 0.45 | 1.5 | | | | |
| Methylene chloride | U | 0.86 | 2.9 | | | | |
| Naphthalene | U | 0.77 | 2.6 | | | | |
| n-Butylbenzene | U | 0.34 | 1.1 | | | | |
| n-Propylbenzene | U | 0.48 | 1.6 | | | | |
| o-Xylene | U | 0.31 | 1.0 | | | | |
| p-Isopropyltoluene | U | 0.26 | 0.88 | | | | |
| sec-Butylbenzene | U | 0.3 | 1.0 | | | | |
| Styrene | U | 0.33 | 1.1 | | | | |
| tert-Butylbenzene | U | 0.39 | 1.3 | | | | |
| Tetrachloroethene | U | 0.39 | 1.3 | | | | |
| Toluene | U | 0.45 | 1.5 | | | | |
| trans-1,2-Dichloroethene | U | 0.48 | 1.6 | | | | |
| trans-1,3-Dichloropropene | U | 0.38 | 2.7 | | | | |
| Trichloroethene | U | 0.43 | 1.4 | | | | |
| Trichlorofluoromethane | U | 0.52 | 1.7 | | | | |
| Vinyl chloride | U | 0.53 | 1.8 | | | | |
| Xylenes, Total | U | 0.81 | 4.4 | | | | |
| Surr: 1,2-Dichloroethane-d4 | 20.3 | 0 | 0 | 20 | 0 | 102 | 75-120 |
| Surr: 4-Bromofluorobenzene | 19.55 | 0 | 0 | 20 | 0 | 97.8 | 80-110 |
| Surr: Dibromofluoromethane | 20.45 | 0 | 0 | 20 | 0 | 102 | 85-115 |
| Surr: Toluene-d8 | 19.9 | 0 | 0 | 20 | 0 | 99.5 | 85-110 |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

Batch ID: **R325443a** Instrument ID **VMS8** Method: **SW8260C**

| LCS | Sample ID: 8V-LCSW1-210827-R325443a | | | | Units: µg/L | | Analysis Date: 8/27/2021 01:44 PM | | | |
|-----------------------------|--|------|-----|-----------------------|--------------------|------------|--|---------------|----------------|------|
| Client ID: | Run ID: VMS8_210827A | | | SeqNo: 7704530 | | Prep Date: | | DF: 1 | | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | RPD %RPD Limit | Qual |
| 1,1,1,2-Tetrachloroethane | 18.07 | 0.38 | 1.3 | 20 | 0 | 90.4 | 73-114 | 0 | 0 | |
| 1,1,1-Trichloroethane | 17.95 | 0.46 | 1.5 | 20 | 0 | 89.8 | 75-130 | 0 | 0 | |
| 1,1,2,2-Tetrachloroethane | 20.56 | 0.4 | 1.3 | 20 | 0 | 103 | 75-130 | 0 | 0 | |
| 1,1,2-Trichloroethane | 18.45 | 0.46 | 1.5 | 20 | 0 | 92.2 | 75-125 | 0 | 0 | |
| 1,1-Dichloroethane | 17.3 | 0.44 | 1.5 | 20 | 0 | 86.5 | 68-142 | 0 | 0 | |
| 1,1-Dichloroethene | 17.42 | 0.4 | 1.4 | 20 | 0 | 87.1 | 70-145 | 0 | 0 | |
| 1,1-Dichloropropene | 17.4 | 0.37 | 1.2 | 20 | 0 | 87 | 75-135 | 0 | 0 | |
| 1,2,3-Trichlorobenzene | 17.63 | 0.42 | 1.4 | 20 | 0 | 88.2 | 70-140 | 0 | 0 | |
| 1,2,3-Trichloropropane | 17.72 | 0.4 | 1.3 | 20 | 0 | 88.6 | 75-125 | 0 | 0 | |
| 1,2,4-Trichlorobenzene | 17.88 | 0.45 | 1.5 | 20 | 0 | 89.4 | 70-135 | 0 | 0 | |
| 1,2,4-Trimethylbenzene | 17.43 | 0.45 | 1.5 | 20 | 0 | 87.2 | 75-130 | 0 | 0 | |
| 1,2-Dibromo-3-chloropropane | 16.59 | 0.43 | 1.4 | 20 | 0 | 83 | 60-130 | 0 | 0 | |
| 1,2-Dibromoethane | 19.68 | 0.41 | 1.4 | 20 | 0 | 98.4 | 67-155 | 0 | 0 | |
| 1,2-Dichlorobenzene | 16.79 | 0.32 | 1.1 | 20 | 0 | 84 | 70-130 | 0 | 0 | |
| 1,2-Dichloroethane | 18.69 | 0.44 | 1.4 | 20 | 0 | 93.4 | 78-125 | 0 | 0 | |
| 1,2-Dichloropropene | 20.22 | 0.48 | 1.6 | 20 | 0 | 101 | 75-125 | 0 | 0 | |
| 1,3,5-Trimethylbenzene | 17.23 | 0.65 | 2.2 | 20 | 0 | 86.2 | 75-130 | 0 | 0 | |
| 1,3-Dichlorobenzene | 18.58 | 0.33 | 1.1 | 20 | 0 | 92.9 | 75-130 | 0 | 0 | |
| 1,3-Dichloropropane | 19.03 | 0.4 | 1.3 | 20 | 0 | 95.2 | 75-125 | 0 | 0 | |
| 1,4-Dichlorobenzene | 17.33 | 0.35 | 1.2 | 20 | 0 | 86.6 | 75-130 | 0 | 0 | |
| 2,2-Dichloropropane | 21.9 | 0.52 | 1.7 | 20 | 0 | 110 | 43-150 | 0 | 0 | |
| 2-Butanone | 18.4 | 0.52 | 1.7 | 20 | 0 | 92 | 55-150 | 0 | 0 | |
| 2-Chlorotoluene | 16.8 | 0.36 | 1.2 | 20 | 0 | 84 | 76-117 | 0 | 0 | |
| 4-Chlorotoluene | 16.92 | 0.31 | 1.0 | 20 | 0 | 84.6 | 80-125 | 0 | 0 | |
| 4-Methyl-2-pentanone | 26.68 | 0.52 | 1.7 | 20 | 0 | 133 | 77-178 | 0 | 0 | |
| Acetone | 18.53 | 6.2 | 21 | 20 | 0 | 92.6 | 60-160 | 0 | 0 | J |
| Benzene | 18.98 | 0.46 | 1.5 | 20 | 0 | 94.9 | 70-130 | 0 | 0 | |
| Bromobenzene | 18.32 | 0.38 | 1.3 | 20 | 0 | 91.6 | 80-125 | 0 | 0 | |
| Bromochloromethane | 19.34 | 0.45 | 1.5 | 20 | 0 | 96.7 | 72-141 | 0 | 0 | |
| Bromodichloromethane | 18.54 | 0.49 | 1.6 | 20 | 0 | 92.7 | 75-125 | 0 | 0 | |
| Bromoform | 15.07 | 0.56 | 1.9 | 20 | 0 | 75.4 | 60-125 | 0 | 0 | |
| Bromomethane | 18.55 | 0.9 | 3.0 | 20 | 0 | 92.8 | 30-185 | 0 | 0 | |
| Carbon tetrachloride | 16.09 | 0.4 | 1.4 | 20 | 0 | 80.4 | 65-140 | 0 | 0 | |
| Chlorobenzene | 18.76 | 0.4 | 1.3 | 20 | 0 | 93.8 | 80-120 | 0 | 0 | |
| Chloroethane | 17.38 | 0.68 | 2.3 | 20 | 0 | 86.9 | 31-172 | 0 | 0 | |
| Chloroform | 17.23 | 0.46 | 1.5 | 20 | 0 | 86.2 | 66-135 | 0 | 0 | |
| Chloromethane | 12.45 | 0.83 | 2.8 | 20 | 0 | 62.2 | 46-148 | 0 | 0 | |
| cis-1,2-Dichloroethene | 18.83 | 0.42 | 1.4 | 20 | 0 | 94.2 | 75-134 | 0 | 0 | |
| cis-1,3-Dichloropropene | 17.96 | 0.57 | 1.9 | 20 | 0 | 89.8 | 70-130 | 0 | 0 | |
| Dibromochloromethane | 16.93 | 0.4 | 1.3 | 20 | 0 | 84.6 | 60-115 | 0 | 0 | |
| Dibromomethane | 18.56 | 0.65 | 2.2 | 20 | 0 | 92.8 | 79-126 | 0 | 0 | |
| Dichlorodifluoromethane | 13.02 | 0.68 | 2.3 | 20 | 0 | 65.1 | 10-180 | 0 | 0 | |
| Ethylbenzene | 16.67 | 0.34 | 1.1 | 20 | 0 | 83.4 | 76-123 | 0 | 0 | |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: R325443a | Instrument ID VMS8 | Method: SW8260C | | | | | | |
|-----------------------------|--------------------|-----------------|------|----|---|------|--------|---|
| Hexachlorobutadiene | 19.21 | 0.56 | 1.9 | 20 | 0 | 96 | 70-155 | 0 |
| Isopropylbenzene | 19.07 | 0.35 | 1.2 | 20 | 0 | 95.4 | 80-127 | 0 |
| m,p-Xylene | 34.55 | 0.81 | 2.7 | 40 | 0 | 86.4 | 75-130 | 0 |
| Methyl tert-butyl ether | 18.05 | 0.45 | 1.5 | 20 | 0 | 90.2 | 68-129 | 0 |
| Methylene chloride | 18.3 | 0.86 | 2.9 | 20 | 0 | 91.5 | 72-125 | 0 |
| Naphthalene | 16.52 | 0.77 | 2.6 | 20 | 0 | 82.6 | 55-160 | 0 |
| n-Butylbenzene | 16.49 | 0.34 | 1.1 | 20 | 0 | 82.4 | 75-145 | 0 |
| n-Propylbenzene | 16.94 | 0.48 | 1.6 | 20 | 0 | 84.7 | 76-116 | 0 |
| o-Xylene | 17.41 | 0.31 | 1.0 | 20 | 0 | 87 | 76-127 | 0 |
| p-Isopropyltoluene | 17.44 | 0.26 | 0.88 | 20 | 0 | 87.2 | 61-164 | 0 |
| sec-Butylbenzene | 18.01 | 0.3 | 1.0 | 20 | 0 | 90 | 80-134 | 0 |
| Styrene | 19.38 | 0.33 | 1.1 | 20 | 0 | 96.9 | 79-117 | 0 |
| tert-Butylbenzene | 17.75 | 0.39 | 1.3 | 20 | 0 | 88.8 | 70-130 | 0 |
| Tetrachloroethene | 17.31 | 0.39 | 1.3 | 20 | 0 | 86.6 | 68-166 | 0 |
| Toluene | 17.95 | 0.45 | 1.5 | 20 | 0 | 89.8 | 76-125 | 0 |
| trans-1,2-Dichloroethene | 17.97 | 0.48 | 1.6 | 20 | 0 | 89.8 | 80-140 | 0 |
| trans-1,3-Dichloropropene | 19.06 | 0.38 | 2.7 | 20 | 0 | 95.3 | 56-132 | 0 |
| Trichloroethene | 16.41 | 0.43 | 1.4 | 20 | 0 | 82 | 77-125 | 0 |
| Trichlorofluoromethane | 14.97 | 0.52 | 1.7 | 20 | 0 | 74.8 | 60-140 | 0 |
| Vinyl chloride | 14.52 | 0.53 | 1.8 | 20 | 0 | 72.6 | 50-136 | 0 |
| Xylenes, Total | 51.96 | 0.81 | 4.4 | 60 | 0 | 86.6 | 76-127 | 0 |
| Surr: 1,2-Dichloroethane-d4 | 19.52 | 0 | 0 | 20 | 0 | 97.6 | 75-120 | 0 |
| Surr: 4-Bromofluorobenzene | 20.3 | 0 | 0 | 20 | 0 | 102 | 80-110 | 0 |
| Surr: Dibromofluoromethane | 20.05 | 0 | 0 | 20 | 0 | 100 | 85-115 | 0 |
| Surr: Toluene-d8 | 20.33 | 0 | 0 | 20 | 0 | 102 | 85-110 | 0 |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

QC Page: 20 of 25

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

Batch ID: **R325443a** Instrument ID **VMS8** Method: **SW8260C**

| MS | Sample ID: 21082011-02A MS | | | | Units: µg/L | | Analysis Date: 8/27/2021 11:11 PM | | | |
|-----------------------------|-----------------------------------|-----|-----------------------------|---------|-----------------------|------|--|---------------|----------------|------|
| | Client ID: | | Run ID: VMS8_210827A | | SeqNo: 7704557 | | Prep Date: | | DF: 20 | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | RPD %RPD Limit | Qual |
| 1,1,1,2-Tetrachloroethane | 381.8 | 7.6 | 26 | 400 | 0 | 95.4 | 73-114 | 0 | | H |
| 1,1,1-Trichloroethane | 413.6 | 9.2 | 30 | 400 | 0 | 103 | 75-130 | 0 | | H |
| 1,1,2,2-Tetrachloroethane | 405.4 | 8 | 27 | 400 | 0 | 101 | 75-130 | 0 | | H |
| 1,1,2-Trichloroethane | 392.6 | 9.2 | 31 | 400 | 0 | 98.2 | 75-125 | 0 | | H |
| 1,1-Dichloroethane | 385.8 | 8.8 | 29 | 400 | 0 | 96.4 | 68-142 | 0 | | H |
| 1,1-Dichloroethene | 456 | 8 | 27 | 400 | 0 | 114 | 70-145 | 0 | | H |
| 1,1-Dichloropropene | 389.2 | 7.4 | 25 | 400 | 0 | 97.3 | 75-135 | 0 | | H |
| 1,2,3-Trichlorobenzene | 288.8 | 8.4 | 28 | 400 | 0 | 72.2 | 70-140 | 0 | | H |
| 1,2,3-Trichloropropane | 351.4 | 8 | 26 | 400 | 0 | 87.8 | 75-125 | 0 | | H |
| 1,2,4-Trichlorobenzene | 333.2 | 9 | 30 | 400 | 0 | 83.3 | 70-135 | 0 | | H |
| 1,2,4-Trimethylbenzene | 386.8 | 9 | 30 | 400 | 0 | 96.7 | 75-130 | 0 | | H |
| 1,2-Dibromo-3-chloropropane | 284.2 | 8.6 | 29 | 400 | 0 | 71 | 60-130 | 0 | | H |
| 1,2-Dibromoethane | 396.8 | 8.2 | 27 | 400 | 0 | 99.2 | 67-155 | 0 | | H |
| 1,2-Dichlorobenzene | 354.4 | 6.4 | 21 | 400 | 0 | 88.6 | 70-130 | 0 | | H |
| 1,2-Dichloroethane | 378.6 | 8.8 | 29 | 400 | 0 | 94.6 | 78-125 | 0 | | H |
| 1,2-Dichloropropane | 425.4 | 9.6 | 32 | 400 | 0 | 106 | 75-125 | 0 | | H |
| 1,3,5-Trimethylbenzene | 372.4 | 13 | 43 | 400 | 0 | 93.1 | 75-130 | 0 | | H |
| 1,3-Dichlorobenzene | 396.2 | 6.6 | 22 | 400 | 0 | 99 | 75-130 | 0 | | H |
| 1,3-Dichloropropane | 401 | 8 | 26 | 400 | 0 | 100 | 75-125 | 0 | | H |
| 1,4-Dichlorobenzene | 359 | 7 | 23 | 400 | 0 | 89.8 | 75-130 | 0 | | H |
| 2,2-Dichloropropane | 444 | 10 | 34 | 400 | 0 | 111 | 43-150 | 0 | | H |
| 2-Butanone | 428.2 | 10 | 35 | 400 | 0 | 107 | 55-150 | 0 | | H |
| 2-Chlorotoluene | 368.2 | 7.2 | 24 | 400 | 0 | 92 | 76-117 | 0 | | H |
| 4-Chlorotoluene | 370.4 | 6.2 | 20 | 400 | 0 | 92.6 | 80-125 | 0 | | H |
| 4-Methyl-2-pentanone | 522.4 | 10 | 35 | 400 | 0 | 131 | 77-178 | 0 | | H |
| Acetone | 390 | 120 | 410 | 400 | 7 | 95.8 | 60-160 | 0 | | JH |
| Benzene | 408.8 | 9.2 | 30 | 400 | 0 | 102 | 70-130 | 0 | | H |
| Bromobenzene | 388.8 | 7.6 | 25 | 400 | 0 | 97.2 | 80-125 | 0 | | H |
| Bromochloromethane | 436.8 | 9 | 30 | 400 | 0 | 109 | 72-141 | 0 | | H |
| Bromodichloromethane | 384 | 9.8 | 33 | 400 | 0 | 96 | 75-125 | 0 | | H |
| Bromoform | 310 | 11 | 37 | 400 | 0 | 77.5 | 60-125 | 0 | | H |
| Bromomethane | 651 | 18 | 60 | 400 | 0 | 163 | 30-185 | 0 | | H |
| Carbon tetrachloride | 381.8 | 8 | 27 | 400 | 0 | 95.4 | 65-140 | 0 | | H |
| Chlorobenzene | 399.6 | 8 | 27 | 400 | 0 | 99.9 | 80-120 | 0 | | H |
| Chloroethane | 773.6 | 14 | 45 | 400 | 0 | 193 | 31-172 | 0 | | SH |
| Chloroform | 379.2 | 9.2 | 31 | 400 | 0 | 94.8 | 66-135 | 0 | | H |
| Chloromethane | 317.4 | 17 | 55 | 400 | 0 | 79.4 | 46-148 | 0 | | H |
| cis-1,2-Dichloroethene | 408.2 | 8.4 | 28 | 400 | 0 | 102 | 75-134 | 0 | | H |
| cis-1,3-Dichloropropene | 368.2 | 11 | 38 | 400 | 0 | 92 | 70-130 | 0 | | H |
| Dibromochloromethane | 343.2 | 8 | 26 | 400 | 0 | 85.8 | 60-115 | 0 | | H |
| Dibromomethane | 390.6 | 13 | 43 | 400 | 0 | 97.6 | 79-126 | 0 | | H |
| Dichlorodifluoromethane | 435 | 14 | 45 | 400 | 0 | 109 | 10-180 | 0 | | H |
| Ethylbenzene | 391 | 6.8 | 22 | 400 | 0 | 97.8 | 76-123 | 0 | | H |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: R325443a | Instrument ID VMS8 | Method: SW8260C | | | | | | | |
|-----------------------------|--------------------|-----------------|----|------|---|------|--------|---|----|
| Hexachlorobutadiene | 372 | 11 | 37 | 400 | 0 | 93 | 70-155 | 0 | H |
| Isopropylbenzene | 420.8 | 7 | 23 | 400 | 0 | 105 | 80-127 | 0 | H |
| m,p-Xylene | 833.6 | 16 | 54 | 800 | 0 | 104 | 75-130 | 0 | H |
| Methyl tert-butyl ether | 541.6 | 9 | 30 | 400 | 0 | 135 | 68-129 | 0 | SH |
| Methylene chloride | 411 | 17 | 58 | 400 | 0 | 103 | 72-125 | 0 | H |
| Naphthalene | 263 | 15 | 51 | 400 | 0 | 65.8 | 55-160 | 0 | H |
| n-Butylbenzene | 364 | 6.8 | 22 | 400 | 0 | 91 | 75-145 | 0 | H |
| n-Propylbenzene | 370.6 | 9.6 | 32 | 400 | 0 | 92.6 | 76-116 | 0 | H |
| o-Xylene | 398 | 6.2 | 21 | 400 | 0 | 99.5 | 76-127 | 0 | H |
| p-Isopropyltoluene | 374.4 | 5.2 | 18 | 400 | 0 | 93.6 | 61-164 | 0 | H |
| sec-Butylbenzene | 381.8 | 6 | 20 | 400 | 0 | 95.4 | 80-134 | 0 | H |
| Styrene | 413.8 | 6.6 | 22 | 400 | 0 | 103 | 79-117 | 0 | H |
| tert-Butylbenzene | 381 | 7.8 | 26 | 400 | 0 | 95.2 | 70-130 | 0 | H |
| Tetrachloroethene | 395.8 | 7.8 | 26 | 400 | 0 | 99 | 68-166 | 0 | H |
| Toluene | 397.4 | 9 | 30 | 400 | 0 | 99.4 | 76-125 | 0 | H |
| trans-1,2-Dichloroethene | 425 | 9.6 | 32 | 400 | 0 | 106 | 80-140 | 0 | H |
| trans-1,3-Dichloropropene | 387.8 | 7.6 | 55 | 400 | 0 | 97 | 56-132 | 0 | H |
| Trichloroethene | 358.4 | 8.6 | 29 | 400 | 0 | 89.6 | 77-125 | 0 | H |
| Trichlorofluoromethane | 456 | 10 | 34 | 400 | 0 | 114 | 60-140 | 0 | H |
| Vinyl chloride | 400 | 11 | 35 | 400 | 0 | 100 | 50-136 | 0 | H |
| Xylenes, Total | 1232 | 16 | 89 | 1200 | 0 | 103 | 76-127 | 0 | H |
| Surr: 1,2-Dichloroethane-d4 | 392 | 0 | 0 | 400 | 0 | 98 | 75-120 | 0 | |
| Surr: 4-Bromofluorobenzene | 405.2 | 0 | 0 | 400 | 0 | 101 | 80-110 | 0 | |
| Surr: Dibromofluoromethane | 403 | 0 | 0 | 400 | 0 | 101 | 85-115 | 0 | |
| Surr: Toluene-d8 | 403 | 0 | 0 | 400 | 0 | 101 | 85-110 | 0 | |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

Batch ID: **R325443a** Instrument ID **VMS8** Method: **SW8260C**

| MSD | | Sample ID: 21082011-02A MSD | | | | Units: µg/L | | Analysis Date: 8/27/2021 11:30 PM | | | |
|-----------------------------|--------|------------------------------------|-----|---------|---------------|-----------------------|---------------|--|-------|---------------|------|
| Client ID: | | Run ID: VMS8_210827A | | | | SeqNo: 7704558 | | Prep Date: | | DF: 20 | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | %RPD | RPD Limit | Qual |
| 1,1,1,2-Tetrachloroethane | 392 | 7.6 | 26 | 400 | 0 | 98 | 73-114 | 381.8 | 2.64 | 30 | H |
| 1,1,1-Trichloroethane | 419.4 | 9.2 | 30 | 400 | 0 | 105 | 75-130 | 413.6 | 1.39 | 30 | H |
| 1,1,2,2-Tetrachloroethane | 416.6 | 8 | 27 | 400 | 0 | 104 | 75-130 | 405.4 | 2.73 | 30 | H |
| 1,1,2-Trichloroethane | 405 | 9.2 | 31 | 400 | 0 | 101 | 75-125 | 392.6 | 3.11 | 30 | H |
| 1,1-Dichloroethane | 397.6 | 8.8 | 29 | 400 | 0 | 99.4 | 68-142 | 385.8 | 3.01 | 30 | H |
| 1,1-Dichloroethene | 466.2 | 8 | 27 | 400 | 0 | 117 | 70-145 | 456 | 2.21 | 30 | H |
| 1,1-Dichloropropene | 413.2 | 7.4 | 25 | 400 | 0 | 103 | 75-135 | 389.2 | 5.98 | 30 | H |
| 1,2,3-Trichlorobenzene | 293 | 8.4 | 28 | 400 | 0 | 73.2 | 70-140 | 288.8 | 1.44 | 30 | H |
| 1,2,3-Trichloropropane | 360.6 | 8 | 26 | 400 | 0 | 90.2 | 75-125 | 351.4 | 2.58 | 30 | H |
| 1,2,4-Trichlorobenzene | 349.6 | 9 | 30 | 400 | 0 | 87.4 | 70-135 | 333.2 | 4.8 | 30 | H |
| 1,2,4-Trimethylbenzene | 389.4 | 9 | 30 | 400 | 0 | 97.4 | 75-130 | 386.8 | 0.67 | 30 | H |
| 1,2-Dibromo-3-chloropropane | 300 | 8.6 | 29 | 400 | 0 | 75 | 60-130 | 284.2 | 5.41 | 30 | H |
| 1,2-Dibromoethane | 408.2 | 8.2 | 27 | 400 | 0 | 102 | 67-155 | 396.8 | 2.83 | 30 | H |
| 1,2-Dichlorobenzene | 368.2 | 6.4 | 21 | 400 | 0 | 92 | 70-130 | 354.4 | 3.82 | 30 | H |
| 1,2-Dichloroethane | 403.8 | 8.8 | 29 | 400 | 0 | 101 | 78-125 | 378.6 | 6.44 | 30 | H |
| 1,2-Dichloropropene | 443.4 | 9.6 | 32 | 400 | 0 | 111 | 75-125 | 425.4 | 4.14 | 30 | H |
| 1,3,5-Trimethylbenzene | 383.6 | 13 | 43 | 400 | 0 | 95.9 | 75-130 | 372.4 | 2.96 | 30 | H |
| 1,3-Dichlorobenzene | 409.2 | 6.6 | 22 | 400 | 0 | 102 | 75-130 | 396.2 | 3.23 | 30 | H |
| 1,3-Dichloropropane | 401.6 | 8 | 26 | 400 | 0 | 100 | 75-125 | 401 | 0.15 | 30 | H |
| 1,4-Dichlorobenzene | 378 | 7 | 23 | 400 | 0 | 94.5 | 75-130 | 359 | 5.16 | 30 | H |
| 2,2-Dichloropropane | 457.4 | 10 | 34 | 400 | 0 | 114 | 43-150 | 444 | 2.97 | 30 | H |
| 2-Butanone | 439.4 | 10 | 35 | 400 | 0 | 110 | 55-150 | 428.2 | 2.58 | 30 | H |
| 2-Chlorotoluene | 375 | 7.2 | 24 | 400 | 0 | 93.8 | 76-117 | 368.2 | 1.83 | 30 | H |
| 4-Chlorotoluene | 375.6 | 6.2 | 20 | 400 | 0 | 93.9 | 80-125 | 370.4 | 1.39 | 30 | H |
| 4-Methyl-2-pentanone | 566.8 | 10 | 35 | 400 | 0 | 142 | 77-178 | 522.4 | 8.15 | 30 | H |
| Acetone | 416.6 | 120 | 410 | 400 | 7 | 102 | 60-160 | 390 | 6.6 | 30 | H |
| Benzene | 419.8 | 9.2 | 30 | 400 | 0 | 105 | 70-130 | 408.8 | 2.66 | 30 | H |
| Bromobenzene | 399.8 | 7.6 | 25 | 400 | 0 | 100 | 80-125 | 388.8 | 2.79 | 30 | H |
| Bromochloromethane | 463.6 | 9 | 30 | 400 | 0 | 116 | 72-141 | 436.8 | 5.95 | 30 | H |
| Bromodichloromethane | 416.2 | 9.8 | 33 | 400 | 0 | 104 | 75-125 | 384 | 8.05 | 30 | H |
| Bromoform | 316.6 | 11 | 37 | 400 | 0 | 79.2 | 60-125 | 310 | 2.11 | 30 | H |
| Bromomethane | 668.6 | 18 | 60 | 400 | 0 | 167 | 30-185 | 651 | 2.67 | 30 | H |
| Carbon tetrachloride | 401.2 | 8 | 27 | 400 | 0 | 100 | 65-140 | 381.8 | 4.96 | 30 | H |
| Chlorobenzene | 417.4 | 8 | 27 | 400 | 0 | 104 | 80-120 | 399.6 | 4.36 | 30 | H |
| Chloroethane | 835.6 | 14 | 45 | 400 | 0 | 209 | 31-172 | 773.6 | 7.71 | 30 | SH |
| Chloroform | 395.2 | 9.2 | 31 | 400 | 0 | 98.8 | 66-135 | 379.2 | 4.13 | 30 | H |
| Chloromethane | 336.6 | 17 | 55 | 400 | 0 | 84.2 | 46-148 | 317.4 | 5.87 | 30 | H |
| cis-1,2-Dichloroethene | 410.6 | 8.4 | 28 | 400 | 0 | 103 | 75-134 | 408.2 | 0.586 | 30 | H |
| cis-1,3-Dichloropropene | 382.4 | 11 | 38 | 400 | 0 | 95.6 | 70-130 | 368.2 | 3.78 | 30 | H |
| Dibromochloromethane | 356 | 8 | 26 | 400 | 0 | 89 | 60-115 | 343.2 | 3.66 | 30 | H |
| Dibromomethane | 414.6 | 13 | 43 | 400 | 0 | 104 | 79-126 | 390.6 | 5.96 | 30 | H |
| Dichlorodifluoromethane | 435.8 | 14 | 45 | 400 | 0 | 109 | 10-180 | 435 | 0.184 | 30 | H |
| Ethylbenzene | 386 | 6.8 | 22 | 400 | 0 | 96.5 | 76-123 | 391 | 1.29 | 30 | H |

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

| Batch ID: R325443a | Instrument ID VMS8 | Method: SW8260C | | | | | | | | | |
|-----------------------------|--------------------|-----------------|----|------|---|------|--------|-------|-------|----|----|
| Hexachlorobutadiene | 383.2 | 11 | 37 | 400 | 0 | 95.8 | 70-155 | 372 | 2.97 | 30 | H |
| Isopropylbenzene | 433.8 | 7 | 23 | 400 | 0 | 108 | 80-127 | 420.8 | 3.04 | 30 | H |
| m,p-Xylene | 806.6 | 16 | 54 | 800 | 0 | 101 | 75-130 | 833.6 | 3.29 | 30 | H |
| Methyl tert-butyl ether | 554 | 9 | 30 | 400 | 0 | 138 | 68-129 | 541.6 | 2.26 | 30 | SH |
| Methylene chloride | 426.6 | 17 | 58 | 400 | 0 | 107 | 72-125 | 411 | 3.72 | 30 | H |
| Naphthalene | 275.6 | 15 | 51 | 400 | 0 | 68.9 | 55-160 | 263 | 4.68 | 30 | H |
| n-Butylbenzene | 380.2 | 6.8 | 22 | 400 | 0 | 95 | 75-145 | 364 | 4.35 | 30 | H |
| n-Propylbenzene | 381.6 | 9.6 | 32 | 400 | 0 | 95.4 | 76-116 | 370.6 | 2.92 | 30 | H |
| o-Xylene | 397.2 | 6.2 | 21 | 400 | 0 | 99.3 | 76-127 | 398 | 0.201 | 30 | H |
| p-Isopropyltoluene | 393 | 5.2 | 18 | 400 | 0 | 98.2 | 61-164 | 374.4 | 4.85 | 30 | H |
| sec-Butylbenzene | 392.2 | 6 | 20 | 400 | 0 | 98 | 80-134 | 381.8 | 2.69 | 30 | H |
| Styrene | 424.2 | 6.6 | 22 | 400 | 0 | 106 | 79-117 | 413.8 | 2.48 | 30 | H |
| tert-Butylbenzene | 390.8 | 7.8 | 26 | 400 | 0 | 97.7 | 70-130 | 381 | 2.54 | 30 | H |
| Tetrachloroethene | 392.6 | 7.8 | 26 | 400 | 0 | 98.2 | 68-166 | 395.8 | 0.812 | 30 | H |
| Toluene | 406.4 | 9 | 30 | 400 | 0 | 102 | 76-125 | 397.4 | 2.24 | 30 | H |
| trans-1,2-Dichloroethene | 431.8 | 9.6 | 32 | 400 | 0 | 108 | 80-140 | 425 | 1.59 | 30 | H |
| trans-1,3-Dichloropropene | 396.2 | 7.6 | 55 | 400 | 0 | 99 | 56-132 | 387.8 | 2.14 | 30 | H |
| Trichloroethene | 376.4 | 8.6 | 29 | 400 | 0 | 94.1 | 77-125 | 358.4 | 4.9 | 30 | H |
| Trichlorofluoromethane | 463 | 10 | 34 | 400 | 0 | 116 | 60-140 | 456 | 1.52 | 30 | H |
| Vinyl chloride | 415.6 | 11 | 35 | 400 | 0 | 104 | 50-136 | 400 | 3.83 | 30 | H |
| Xylenes, Total | 1204 | 16 | 89 | 1200 | 0 | 100 | 76-127 | 1232 | 2.28 | 30 | H |
| Surr: 1,2-Dichloroethane-d4 | 394.4 | 0 | 0 | 400 | 0 | 98.6 | 75-120 | 392 | 0.61 | 30 | |
| Surr: 4-Bromofluorobenzene | 398 | 0 | 0 | 400 | 0 | 99.5 | 80-110 | 405.2 | 1.79 | 30 | |
| Surr: Dibromofluoromethane | 404 | 0 | 0 | 400 | 0 | 101 | 85-115 | 403 | 0.248 | 30 | |
| Surr: Toluene-d8 | 402.4 | 0 | 0 | 400 | 0 | 101 | 85-110 | 403 | 0.149 | 30 | |

The following samples were analyzed in this batch:

| 21081889-01A

21081889-02A

21081889-08A

Note: See Qualifiers Page for a list of Qualifiers and their explanation.

QC Page: 24 of 25

Client: Gannett Fleming, Inc.
Work Order: 21081889
Project: Koeller (47358.003)

QC BATCH REPORT

Batch ID: R325619 Instrument ID VMS9 Method: SW8260B

| MBLK | | Sample ID: 9V-BLKW2-210830-R325619 | | | | Units: µg/L | | Analysis Date: 8/30/2021 04:44 PM | | | |
|------------------|--------|------------------------------------|-----|---------|---------------|----------------|---------------|-----------------------------------|------|-----------|------|
| Client ID: | | Run ID: VMS9_210830A | | | | SeqNo: 7709833 | | Prep Date: | | DF: 1 | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | %RPD | RPD Limit | Qual |
| 1,4-Dioxane | U | 0.44 | 1.0 | | | | | | | | |
| Surr: Toluene-d8 | 4.42 | 0 | 0 | 5 | 0 | 88.4 | 74-124 | 0 | 0 | | |
| LCS | | Sample ID: 9V-LCSW2-210830-R325619 | | | | Units: µg/L | | Analysis Date: 8/30/2021 03:57 PM | | | |
| Client ID: | | Run ID: VMS9_210830A | | | | SeqNo: 7709832 | | Prep Date: | | DF: 1 | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | %RPD | RPD Limit | Qual |
| 1,4-Dioxane | 37.55 | 0.44 | 1.0 | 40 | 0 | 93.9 | 70-130 | 0 | 0 | | |
| Surr: Toluene-d8 | 4.29 | 0 | 0 | 5 | 0 | 85.8 | 74-124 | 0 | 0 | | |
| MS | | Sample ID: 21082492-02A MS | | | | Units: µg/L | | Analysis Date: 8/30/2021 06:33 PM | | | |
| Client ID: | | Run ID: VMS9_210830A | | | | SeqNo: 7709839 | | Prep Date: | | DF: 1 | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | %RPD | RPD Limit | Qual |
| 1,4-Dioxane | 35.1 | 0.44 | 1.0 | 40 | 0 | 87.8 | 70-130 | 0 | 0 | | |
| Surr: Toluene-d8 | 4.21 | 0 | 0 | 5 | 0 | 84.2 | 74-124 | 0 | 0 | | |
| DUP | | Sample ID: 21082492-01A DUP | | | | Units: µg/L | | Analysis Date: 8/30/2021 06:17 PM | | | |
| Client ID: | | Run ID: VMS9_210830A | | | | SeqNo: 7709838 | | Prep Date: | | DF: 1 | |
| Analyte | Result | MDL | PQL | SPK Val | SPK Ref Value | %REC | Control Limit | RPD Ref Value | %RPD | RPD Limit | Qual |
| 1,4-Dioxane | U | 0.44 | 1.0 | 0 | 0 | 0 | | 0.62 | 0 | 30 | |
| Surr: Toluene-d8 | 4.63 | 0 | 0 | 5 | 0 | 92.6 | 74-124 | 4.4 | 5.09 | 30 | |

The following samples were analyzed in this batch:

21081889-01A

Note: See Qualifiers Page for a list of Qualifiers and their explanation.



Cincinnati, OH

+1 513 733 5336

Everett, WA

+1 425 356 2600

Fort Collins, CO

+1 970 490 1511

Holland, MI

+1 616 399 6070

Chain of Custody Form

Houston, TX

+1 281 530 5656

Middletown, PA

+1 717 944 5541

Spring City, PA

+1 610 948 4903

Salt Lake City, UT

+1 801 266 7700

South Charleston, WV

+1 304 356 3168

Page 1 of 1

COC ID: 231189

ALS Project Manager: JAD ALS Work Order #: 21081889

| Customer Information | | Project Information | | Parameter/Method Request for Analysis | | | | | | | | | | | |
|----------------------|--|---------------------|------------------------|---------------------------------------|----------------------|--|--|--|--|--|--|--|--|--|--|
| Purchase Order | <u>Z021 ALS</u> | Project Name | <u>Koeller</u> | A | VOLCs | | | | | | | | | | |
| Work Order | | Project Number | <u>47358.003</u> | B | 1,4-Dioxane 8260 SIM | | | | | | | | | | |
| Company Name | Gannett Fleming, Inc | Bill To Company | Gannett Fleming, Inc | C | PFAS 537 Mod. | | | | | | | | | | |
| Send Report To | <u>Tony Miller</u> 8040 Excelsior Drive | Invoice Attn | Accounts Payable | D | | | | | | | | | | | |
| Address | Suite 303 | Address | 8040 Excelsior Drive | E | | | | | | | | | | | |
| City/State/Zip | Madison, WI 53717-1338 | City/State/Zip | Madison, WI 53717-1338 | F | | | | | | | | | | | |
| Phone | (608) 836-1500 | Phone | (608) 836-1500 | G | | | | | | | | | | | |
| Fax | | Fax | | H | | | | | | | | | | | |
| e-Mail Address | <u>awmiller@gfnet.com</u> | e-Mail Address | | I | | | | | | | | | | | |
| | | | | J | | | | | | | | | | | |

| No. | Sample Description | Date | Time | Matrix | Pres. | # Bottles | A | B | C | D | E | F | G | H | I | J | Hold |
|-----|---------------------------|----------------|-------------|-----------|------------|-----------|----------|----------|----------|---|----------|---|---|---|---|---|------|
| 1 | MW-1 | <u>8/18/21</u> | <u>1400</u> | <u>6W</u> | <u>HCl</u> | <u>6</u> | <u>X</u> | <u>X</u> | <u>X</u> | | | | | | | | |
| 2 | MW-2 | | <u>1755</u> | | | <u>3</u> | | | | | | | | | | | |
| 3 | MW-3 | | <u>1325</u> | | | | | | | | | | | | | | |
| 4 | MW-4 | | <u>1140</u> | | | | | | | | | | | | | | |
| 5 | MW-5 | | <u>1025</u> | | | | | | | | | | | | | | |
| 6 | MW-6 | | <u>1210</u> | | | | | | | | | | | | | | |
| 7 | MW-7 | | <u>1105</u> | | | | | | | | | | | | | | |
| 8 | MW-1 Pup | | <u>1405</u> | | | <u>2</u> | | | | | | | | | | | |
| 9 | MW Field Blunk | | | <u>W</u> | | <u>2</u> | | | | | <u>X</u> | | | | | | |
| 10 | Tripp Blunk | | | <u>W</u> | <u>HCl</u> | <u>3</u> | <u>X</u> | | | | | | | | | | |

| | | | |
|--------------------------------|-----------------|--|-------------------|
| Sampler(s) Please Print & Sign | Shipment Method | Required Turnaround Time: (Check Box) | Results Due Date: |
| <u>Marcus Mussey</u> | <u>FedEx</u> | <input type="checkbox"/> Std 10 WK Days <input type="checkbox"/> 5 WK Days <input checked="" type="checkbox"/> Other <input type="checkbox"/> 2 WK Days <input type="checkbox"/> 24 Hour | |

| | | | | |
|---------------------------------------|----------------------|-------------------|---------------------------|--------|
| Relinquished by: <u>Marcus Mussey</u> | Date: <u>8/19/21</u> | Time: <u>9:30</u> | Received by: <u>FedEx</u> | Notes: |
|---------------------------------------|----------------------|-------------------|---------------------------|--------|

| | | | | | | |
|-------------------------------|----------------------|-------------------|--|-----------|--------------|-----------------------------------|
| Relinquished by: <u>FedEx</u> | Date: <u>8/20/21</u> | Time: <u>0830</u> | Received by (Laboratory): <u>[Signature]</u> | Cooler ID | Cooler Temp. | QC Package: (Check One Box Below) |
|-------------------------------|----------------------|-------------------|--|-----------|--------------|-----------------------------------|

| | | | | | | |
|-----------------------------------|----------------------|-------------------|---|------------|------------|--|
| Logged by (Laboratory): <u>KC</u> | Date: <u>8/20/21</u> | Time: <u>1445</u> | Checked by (Laboratory): <u>[Signature]</u> | <u>123</u> | <u>47°</u> | <input type="checkbox"/> Level II Std QC <input type="checkbox"/> TRRP CheckList |
|-----------------------------------|----------------------|-------------------|---|------------|------------|--|

| | | | | | | |
|--|--|--|--|--|--|--|
| Preservative Key: 1-HCl 2-HNO ₃ 3-H ₂ SO ₄ 4-NaOH 5-Na ₂ S ₂ O ₃ 6-NaHSO ₄ 7-Other 8-4°C 9-5035 | | | | | | <input type="checkbox"/> Level III Std QC <input type="checkbox"/> TRRP Level IV |
|--|--|--|--|--|--|--|

| | | | | | | |
|--|--|--|--|--|--|---|
| | | | | | | <input type="checkbox"/> Level IV SW846/CLP |
|--|--|--|--|--|--|---|

| | | | | | | |
|--|--|--|--|--|--|--------------------------------|
| | | | | | | <input type="checkbox"/> Other |
|--|--|--|--|--|--|--------------------------------|

- Note: 1. Any changes must be made in writing once samples and COC Form have been submitted to ALS Environmental.
 2. Unless otherwise agreed in a formal contract, services provided by ALS Environmental are expressly limited to the terms and conditions stated on the reverse.
 3. The Chain of Custody is a legal document. All information must be completed accurately.

Sample Receipt ChecklistClient Name: GANNETTFLEMING - WIDate/Time Received: 20-Aug-21 08:30Work Order: 21081889Received by: KRWChecklist completed by Keith Werenka
eSignature

20-Aug-21

Date

Reviewed by: Jodi Blawie
eSignature

20-Aug-21

Date

Matrices: WaterCarrier name: FedEx

| | | | |
|---|---|--|---|
| Shipping container/cooler in good condition? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | Not Present <input type="checkbox"/> |
| Custody seals intact on shipping container/cooler? | Yes <input type="checkbox"/> | No <input type="checkbox"/> | Not Present <input checked="" type="checkbox"/> |
| Custody seals intact on sample bottles? | Yes <input type="checkbox"/> | No <input type="checkbox"/> | Not Present <input checked="" type="checkbox"/> |
| Chain of custody present? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| Chain of custody signed when relinquished and received? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| Chain of custody agrees with sample labels? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| Samples in proper container/bottle? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| Sample containers intact? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| Sufficient sample volume for indicated test? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| All samples received within holding time? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| Container/Temp Blank temperature in compliance? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| Sample(s) received on ice? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | |
| Temperature(s)/Thermometer(s): | <u>4.4/5.4 C</u> <input type="checkbox"/> <u>IR3</u> <input type="checkbox"/> | | |
| Cooler(s)/Kit(s): | <input type="checkbox"/> | | |
| Date/Time sample(s) sent to storage: | <u>8/20/2021 2:45:00 PM</u> <input type="checkbox"/> | | |
| Water - VOA vials have zero headspace? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | No VOA vials submitted <input type="checkbox"/> |
| Water - pH acceptable upon receipt? | Yes <input checked="" type="checkbox"/> | No <input type="checkbox"/> | N/A <input type="checkbox"/> |
| pH adjusted? | Yes <input type="checkbox"/> | No <input checked="" type="checkbox"/> | N/A <input type="checkbox"/> |
| pH adjusted by: | <input type="checkbox"/> <u>-</u> | | |

Login Notes:

Client Contacted:

Date Contacted:

Person Contacted:

Contacted By:

Regarding:

Comments:

CorrectiveAction:

ATTACHMENT B

CHART SHOWING PCE CONCENTRATIONS/TREND LINES &
DEPTH TO WATER MEASURED IN MW-1

PCE CONCENTRATIONS AND DEPTH TO
WATER TABLE MEASURED IN MW-1
OCTOBER 2006 - AUGUST 2021

