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1302 North Meridian St., Suite 300 • Indianapolis, Indiana 46202

July 13, 2023

Mr. Frank Deveau
Taft Stettinius & Hollister LLP
One Indiana Square, Suite 3500
Indianapolis, Indiana

**Re: Limited Subsurface Investigation
Jessen Manufacturing
1409 West Beardsley Avenue
Elkhart, Indiana
August Mack Project Number JX1600.740**

Dear Mr. Deveau:

August Mack Environmental, Inc. (August Mack) has completed subsurface investigation activities at the above-referenced property ("subject property"). During a recent Phase I Environmental Site Assessment (ESA) performed by August Mack, (August Mack Project No. JX1208.710) the following recognized environmental conditions (RECs) were identified:

- The long-term industrial use of the subject property involving the use of petroleum products and hazardous substances. Historical records indicate the subject property has been developed with a machine shop since the 1920s.
- The presence and unknown source or extent of chlorinated volatile organic compounds (cVOCs) in groundwater identified along the northern boundary of the subject property potentially originating from the upgradient off-site facilities. Three (3) temporary monitoring wells were installed south of W. Beardsley Avenue along the northern boundary of the subject property in 2008 as part of evaluating the extent of groundwater impacts originating from the CTS Corp facility. Tetrachloroethene (PCE), trichloroethene (TCE), and cis-1,2-dichloroethene (DCE) exceeded their respective Indiana Department of

Environmental Management (IDEM) Risk Based Closure Guide (R2) Human Health Published Levels (PLs) in one or more of the samples

The purpose of this investigation was to determine if subsurface conditions on the subject property have been impacted by the RECs identified in the Phase I ESA. This report was prepared at the request of Mr. Frank Deveau with Taft Stettinius & Hollister LLP and may be relied on by Taft Stettinius & Hollister LLP and Jessen Manufacturing Co., Inc.. Reliance on the information and conclusions presented in this report by any other party(ies) is not authorized by August Mack.

SUBSURFACE INVESTIGATION

August Mack mobilized to the subject property on June 21, 2023, to perform the subsurface investigation activities. Prior to completing the borings, ground penetrating radar (GPR), electromagnetic (EM) locating, and other utility locating tools were utilized to clear all boring locations.

A total of seven (7) soil borings (SB-1 through SB-7) were advanced across the subject property using a Geoprobe® direct push sampling system (Geoprobe®). Information regarding boring locations is provided below and boring locations are depicted on **Figure 1**.

- SB-1 through SB-3 were advanced in loading docks and parking areas on the north portion subject property to evaluate the historical and current operations, as well as groundwater conditions on the upgradient side of the property;
- SB-4 and SB-5 were advanced on the southwestern exterior of the subject property building at a location hydraulically downgradient of the current and historical machine shop operations;
- SB-6 was advanced on the west exterior of the subject property building to evaluate the historical and current operations; and,
- SB-7 was advanced in the center of the subject property building in the vicinity of a former trichloroethylene (TCE) degreaser and central to site operations.

Sampling Methodology

Soil borings SB-1 through SB-6 were advanced to a depth of 17-feet below grade (ft bg) or to groundwater whichever came first, using a Geoprobe® sampling system. SB-7 was only advanced to a depth of 1.5 ft bg due to refusal that was caused by dense sand material. The purpose of the borings was to field screen soils, determine geological

conditions, and collect soil and groundwater samples for laboratory analysis. All soil sample intervals were inspected in the field for odors and staining and screened using a photoionization detector (PID). Field screening results and soil lithological information is provided on soil boring logs included as **Attachment A**.

One (1) soil sample interval from SB-1 through SB-3, SB-6 and SB-7 was selected for laboratory analysis based on field inspection observations and screening results. Samples were generally selected from the highest screened, unsaturated interval of each of those borings. Soil samples were submitted to ENVision Laboratories, Inc. (ENVision) located in Indianapolis, Indiana for laboratory analysis of volatile organic compounds (VOCs), semi-VOCs (SVOCs), Resource Conservation Recovery Act (RCRA) 8 Metals, hexavalent chromium, and polychlorinated biphenyls (PCBs). Soil samples from SB-4 and SB-5 were not submitted for laboratory analysis based on the purpose of these borings to evaluate potential down-gradient groundwater impacts, and based on field screening results.

Soil borings SB1 through SB6 were converted into temporary, 1-inch diameter PVC wells to aid in the collection of groundwater samples. Groundwater samples were collected from the temporary wells using a disposable bailer and submitted to ENVision for analysis of VOCs, SVOCs, and dissolved RCRA 8 metal. Due to shallow refusal that was encountered, August Mack was unable to collect a groundwater sample at SB-7.

August Mack field procedures are provided in **Attachment B**.

Field Observations

Inspection of collected soil samples from the borings revealed that the subsurface geology consists primarily of sand to the depths investigated. Saturated conditions were encountered beginning at approximately 4 ft bg at SB-2, and beginning at approximately 8 ft bg at SB-1, SB-3, SB-4, SB-5, and SB-6.

PID measurements collected from SB-1, SB-3, SB-4, and SB-5 were less than 1.1 parts per million (ppm). At SB-2, PID readings up to 127.6 ppm (6 to 8 ft bg) were encountered. At SB-6, PID readings up to 321.9 ppm (2 to 4 ft bg) were encountered; and at SB-7, PID readings up to 25 ppm (1 to 1.5 ft bg) were encountered. Staining was observed at SB-1 from 1 to 4 ft bg and at SB-6 from 1 to 3 ft bg. An odor was identified at SB-2 from 4 to 12 ft bg.

Soil Analytical Results

The soil analytical results were compared to the Indiana Department of Environmental Management (IDEM) Risk-Based Closure Guide (R2) 2023 Long Term Residential Soil Human Health Published Levels (PLs), Long Term Commercial PLs, and Short Term Excavation PLs. The laboratory analysis revealed the following results:

- One (1) or more VOCs, including various petroleum-related VOCs and cVOCs, were detected above the laboratory reporting limits at SB-1, SB-2, SB-6, and SB-7. All detected concentrations were below their respective R2 PLs.
- One (1) or more SVOCs were detected above the laboratory reporting limits at SB-1, SB-3, SB-6, and SB-7. All detected concentrations are below their respective R2 PLs.
- Metals barium and total chromium were detected in all of the soil samples. Additionally, lead was detected in all of the soil samples except for SB-2 (2-4). However, all detected concentrations are below their respective R2 PLs.
- No other constituents of concern were reported above the laboratory reporting limits, which are below their respective R2 PLs.

The soil analytical results are summarized in **Table 1** and a copy of the laboratory analytical report and chain of custody documentation is included in **Attachment C**.

Groundwater Analytical Results

The groundwater analytical results were compared to the IDEM R2 2023 Long-Term Residential PLs. The laboratory analysis revealed the following results:

- Petroleum VOCs n-butylbenzene, sec-butylbenzene, p-isopropyltolune, and n-propylbenzene were detected above the laboratory reporting limits at SB-2, but below their respective R2 PLs.
- One (1) or more of the following cVOCs were detected above the laboratory reporting limits at SB-1, SB-3, and SB-4: PCE, TCE, 1,1,1-trichloroethane (TCA), and cis,1-2-DCE. All were below their respective R2 PLs except for PCE at SB-1 and SB-4; and TCE at SB-1.
- No other constituents of concern were reported above the laboratory reporting limits, which are below their respective R2 PLs.

The groundwater analytical results are summarized in **Table 2** and a copy of the laboratory analytical report and chain of custody documentation is included in **Attachment C.**

SUMMARY AND CONCLUSION

August Mack has completed subsurface investigation activities at 1409 West Beardsley Avenue, Elkhart, Indiana. A total of seven (7) soil borings were advanced across the subject property in order to determine if subsurface conditions have been impacted by the RECs identified in the recent Phase I ESA completed by August Mack related to current and historical industrial operations on the subject property and, the unknown extent of cVOC impacts in groundwater previously identified along the northern boundary.

During field screening activities, evidence of potential impacts including stained soil, odor, and elevated PID readings were encountered in multiple borings at the subject property. Laboratory analysis of shallow soil samples revealed multiple VOCs (including cVOCs), SVOCs, and metals at concentrations above the laboratory reporting limits. However, none were at concentrations exceeding their respective R2 PLS. Laboratory analysis of the groundwater samples revealed PCE and TCE at concentrations above their respective R2 PLS in borings located on the north and east portions of the subject property. Based on the presumed groundwater flow direction, the extent of groundwater impacts potentially extending off-site is unknown. August Mack notes that the exposure pathway to groundwater at the subject property is currently incomplete since groundwater is not used at the subject property and the subject property is connected to the municipal water supply.

We appreciate the opportunity to provide you with environmental consulting services and trust that this submittal is in accordance with your needs. Please feel free to contact us if you have any questions or comments, or require additional information regarding this project or the project site.

Sincerely,

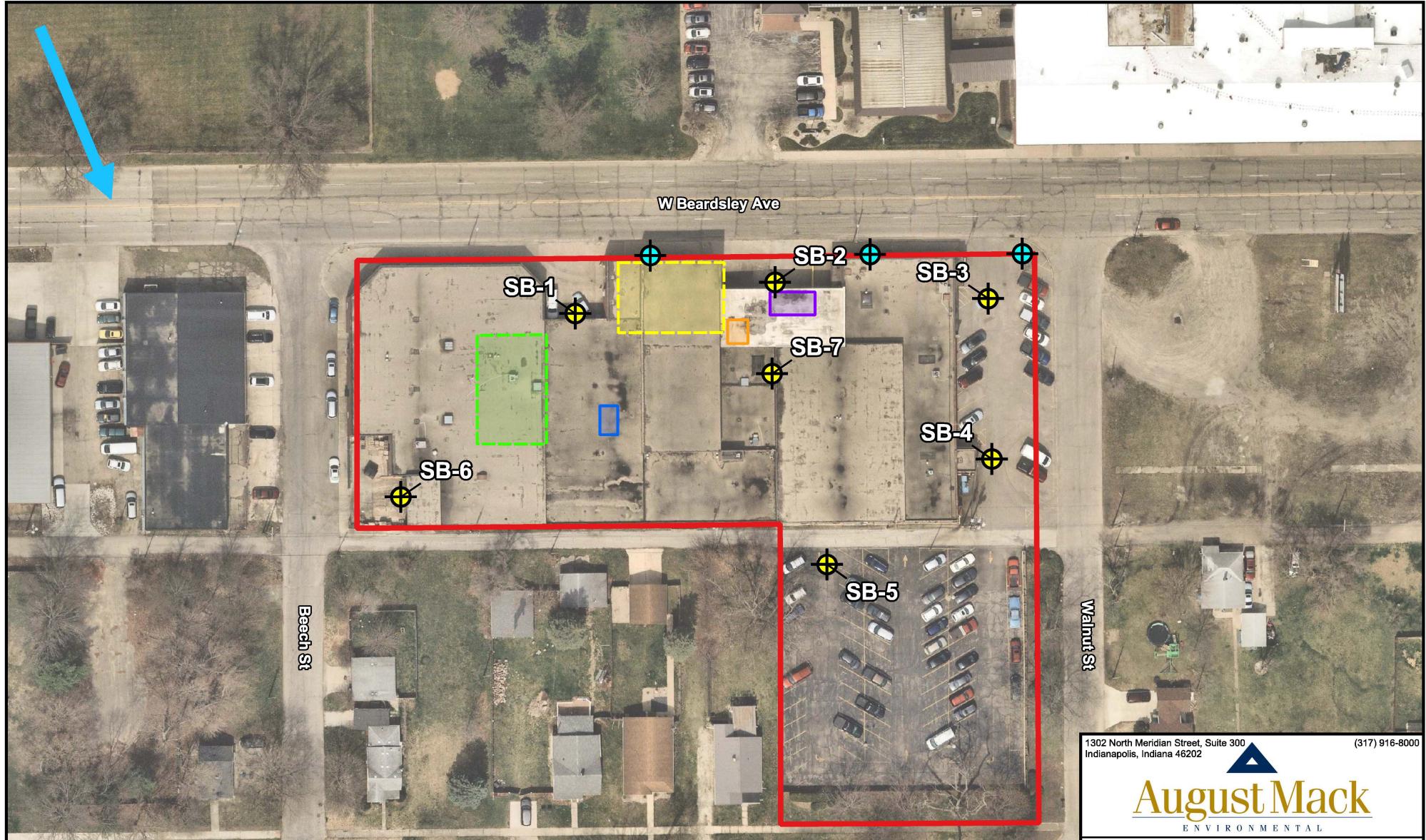
Amanda O'Connor

Samantha O'Connor
Due Diligence Manager

Tyler Zschiedrich

Tyler Zschiedrich
Sr. Due Diligence Manager

Figures



Legend

| | |
|------------------------------------------------------------------------------------------------------------------------------------------------------|--------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| Subject Property | Mineral Spirits AST |
| Former Heating Oil USTs (approximate location) | Former Groundwater Sample Location with VOCs above IDEM PLs (approximate location) |
| Former Lubricant and Cutting Oil USTs (approximate location) | Soil Boring |
| Cutting Oil AST | |
| Hydraulic Oil and Cutting Oil ASTs | → Presumed Groundwater Flow Direction (SSE) |

Maple Row

Nearmap Aerial Imagery: April 9, 2023

1302 North Meridian Street, Suite 300
Indianapolis, Indiana 46202

(317) 916-8000



August Mack
ENVIRONMENTAL

Jessen Manufacturing

1409 West Beardsley Avenue
Elkhart, Indiana 46515

Sample Location Map

| | | |
|-------------------------|----------------|---------------------------------------------------------------------------------------|
| PROJECT NO.: JX1600.740 | |  |
| DATE: 06/23/2023 | SCALE: 1:1,000 | |
| DRAWN BY: ES | FIGURE: 1 |  |

Tables

SUMMARY OF SOIL ANALYTICAL DATA
JESSEN MANUFACTURING

| August Mack ENVIRONMENTAL | Sample ID: | IDEM 2023 LONG TERM RESIDENTIAL SOIL PLs (*) | IDEM 2023 LONG TERM COMMERCIAL SOIL PLs (**) | IDEM 2023 SHORT TERM EXCAVATION SOIL PLs (#) | SB-1 | SB-2 | SB-3 | SB-6 | SB-7 |
|---------------------------------------------------------------------|---------------|----------------------------------------------------------|----------------------------------------------------------|----------------------------------------------------------|------------|------------|------------|------------|------------|
| | Depth - feet: | | | | 2.0-4.0 | 2.0-4.0 | 2.0-4.0 | 2.0-4.0 | 1.0-1.5 |
| | Sample Date: | | | | 06/21/2023 | 06/21/2023 | 06/21/2023 | 06/21/2023 | 06/21/2023 |
| VOLATILE ORGANIC COMPOUNDS (VOCs) VIA USEPA METHOD 8260 | | | | | | | | | |
| n-Butylbenzene | NE | NE | 100 | <0.0059 | <0.0052 | <0.0060 | <0.0055 | 0.198 | |
| sec-Butylbenzene | NE | NE | 100 | <0.0059 | <0.0052 | <0.0060 | 0.0137 | 0.161 | |
| 1,1-Dichloroethane (11DCA) | NE | NE | 2,000 | <0.0059 | <0.0052 | <0.0060 | <0.0055 | 0.0111 | |
| cis-1,2-Dichloroethene (c12DCE) | NE | NE | 1,000 | <0.0059 | <0.0052 | <0.0060 | <0.0055 | 0.357 | |
| Ethylbenzene | NE | NE | 500 | <0.0059 | <0.0052 | <0.0060 | <0.0055 | 0.0285 | |
| Isopropylbenzene (Cumene) | NE | NE | 300 | <0.0059 | <0.0052 | <0.0060 | <0.0055 | 0.0482 | |
| p-Isopropyltoluene | NE | NE | NE | <0.0059 | <0.0052 | <0.0060 | 0.0872 | 0.168 | |
| n-Propylbenzene | NE | NE | 300 | <0.0059 | <0.0052 | <0.0060 | 0.00720 | 0.132 | |
| Tetrachloroethylene (PCE) | NE | NE | 200 | 0.00599 | 0.0188 | <0.0060 | <0.0055 | 0.817 | |
| Trichloroethylene (TCE) | NE | NE | 100 | <0.0059 | <0.0052 | <0.0060 | <0.0055 | 1.50 | |
| 1,2,4-Trimethylbenzene | NE | NE | 200 | <0.0059 | <0.0052 | <0.0060 | 0.211 | 1.67 | |
| 1,3,5-Trimethylbenzene | NE | NE | 200 | <0.0059 | <0.0052 | <0.0060 | 0.622 | 0.580 | |
| Xylene (M&P) | NE | NE | 300 | <0.0059 | <0.0052 | <0.0060 | 0.0113 | 0.0815 | |
| Xylene (Ortho) | NE | NE | 300 | <0.0059 | <0.0052 | <0.0060 | 0.131 | 0.116 | |
| Xylene (Total) | NE | NE | 300 | <0.012 | <0.010 | <0.012 | 0.142 | 0.197 | |
| All Other Analyzed VOCs | Varies | Varies | Varies | BRL | BRL | BRL | BRL | BRL | |
| SEMOVOLATILE ORGANIC COMPOUNDS (SVOCs) VIA USEPA METHOD 8270 | | | | | | | | | |
| Benzo(a)anthracene | 20.0 | 200 | 10,000 | 0.988 | <0.35 | <0.40 | <0.37 | <0.41 | |
| Benzo(a)pyrene | 2.00 | 20.0 | 500 | 1.01 | <0.069 | <0.079 | <0.073 | <0.081 | |
| Benzo(b)fluoranthene | 20.0 | 200 | 10,000 | 1.06 | <0.35 | <0.40 | <0.37 | <0.41 | |
| Benzo(g,h,i)perylene | NE | NE | NE | 0.947 | <0.35 | <0.40 | <0.37 | <0.41 | |
| Benzo(k)fluoranthene | 200 | 2,000 | 100,000 | 0.411 | <0.35 | <0.40 | <0.37 | <0.41 | |
| Chrysene | 2,000 | 20,000 | 100,000 | 1.03 | <0.35 | <0.40 | <0.37 | <0.41 | |
| Fluoranthene | 3,000 | 30,000 | 70,000 | 1.80 | <0.35 | 0.762 | <0.37 | <0.41 | |
| Indeno(1,2,3-cd)pyrene | 20.0 | 200 | 10,000 | 0.885 | <0.35 | <0.40 | <0.37 | <0.41 | |
| 1-Methylnaphthalene | 300 | 400 | 400 | <0.39 | <0.35 | <0.40 | <0.37 | 0.519 | |
| 2-Methylnaphthalene | 300 | 3,000 | 7,000 | <0.39 | <0.35 | <0.40 | <0.37 | 0.684 | |
| Naphthalene | 30 | 90 | 3,000 | <0.078 | <0.069 | <0.079 | 0.893 | 1.00 | |
| Phenanthrene | NE | NE | NE | 0.791 | <0.35 | 0.668 | <0.37 | <0.41 | |
| Pyrene | 3,000 | 20,000 | 50,000 | 1.54 | <0.35 | 0.678 | <0.37 | <0.41 | |
| All Other Analyzed SVOCs | Varies | Varies | Varies | BRL | BRL | BRL | BRL | BRL | |
| METALS VIA USEPA METHODS 6010/7196/7471 | | | | | | | | | |
| Barium | 20,000 | 100,000 | 100,000 | 31 | 4.7 | 17 | 196 | 58 | |
| Chromium (Total) | NE | NE | NE | 8.8 | 7.3 | 4.0 | 43 | 17 | |
| Lead | 400 | 800 | 1,000 | 11 | <2.1 | 8.9 | 36 | 95 | |
| All Other Analyzed Metals | Varies | Varies | Varies | BRL | BRL | BRL | BRL | BRL | |
| POLYCHLORINATED BIPHENYLS (PCBs) VIA USEPA METHOD 8082 | | | | | | | | | |
| All Analyzed PCBs | Varies | Varies | Varies | BRL | BRL | BRL | BRL | BRL | |

Abbreviations & Notes

BRL = Below Laboratory Reporting Limits

IDEM = Indiana Department of Environmental Management

NE = Not Established

PLs = Human Health Published Levels

R2 = Risk-Based Closure Guide

USEPA = United States Environmental Protection Agency

The following denote the symbol and color of screening level exceedances:

* = At or Above 2023 IDEM R2 Long Term Residential Soil PLs

** = At or Above 2023 IDEM R2 Long Term Commercial Soil PLs

= At or Above 2023 IDEM R2 Short Term Excavation Soil PLs

Results and IDEM PLs are reported in milligrams per kilogram (mg/kg).

IDEM PLs are based on the IDEM R2, Table 1: Human Health Published Levels with updates.

SUMMARY OF GROUNDWATER ANALYTICAL DATA
JESSEN MANUFACTURING

|  August Mack <small>ENVIRONMENTAL</small> | Sample ID: | 2023 IDEM R2 LONG TERM RESIDENTIAL GROUNDWATER PLs (^) | SB-1-GW | SB-2-GW | SB-3-GW | SB-4-GW | SB-5-GW | SB-6-GW |
|-----------------------------------------------------------------------------------------------------------------------------------------|---------------|-----------------------------------------------------------------------------------|------------|------------|------------|------------|------------|------------|
| | Depth - feet: | | 6-16 | 2-12 | 7-17 | 6-16 | 6-16 | 6-16 |
| | Sample Date: | | 06/21/2023 | 06/21/2023 | 06/21/2023 | 06/21/2023 | 06/21/2023 | 06/21/2023 |
| VOLATILE ORGANIC COMPOUNDS (VOCs) VIA USEPA METHOD 8260 | | | | | | | | |
| n-Butylbenzene | 1,000 | <5.0 | 24.0 | <5.0 | <5.0 | <5.0 | <5.0 | <5.0 |
| sec-Butylbenzene | 2,000 | <5.0 | 13.2 | <5.0 | <5.0 | <5.0 | <5.0 | <5.0 |
| cis-1,2-Dichloroethene (c12DCE) | 70 | <5.0 | <5.0 | 12.2 | 11.7 | <5.0 | <5.0 | <5.0 |
| p-Isopropyltoluene | NE | <5.0 | 5.59 | <5.0 | <5.0 | <5.0 | <5.0 | <5.0 |
| n-Propylbenzene | 700 | <5.0 | 7.33 | <5.0 | <5.0 | <5.0 | <5.0 | <5.0 |
| Tetrachloroethene (PCE) | 5 | 9.31 ^ | <5.0 | <5.0 | 15.1 ^ | <5.0 | <5.0 | <5.0 |
| 1,1,1-Trichloroethane (111TCA) | 200 | 5.95 | <5.0 | <5.0 | <5.0 | <5.0 | <5.0 | <5.0 |
| Trichloroethene (TCE) | 5 | 8.05 ^ | <5.0 | <5.0 | <5.0 | <5.0 | <5.0 | <5.0 |
| All Other Analyzed VOCs | Varies | BRL | BRL | BRL | BRL | BRL | BRL | BRL |
| SEMIVOLATILE ORGANIC COMPOUNDS (SVOCs) VIA USEPA METHOD 8270/8270 SIM | | | | | | | | |
| All Analyzed SVOCs | Varies | BRL | BRL | BRL | BRL | BRL | BRL | BRL |
| METALS VIA USEPA METHODS 6010/7470 | | | | | | | | |
| All Analyzed Metals | Varies | BRL | BRL | BRL | BRL | BRL | BRL | BRL |

Abbreviations & Notes

BRL = Below Laboratory Reporting Limits

IDEML = Indiana Department of Environmental Management

NE = Not Established

PLs = Human Health Published Levels

R2 = Risk-Based Closure Guide

SIM = Selected Ion Monitoring

USEPA = United States Environmental Protection Agency

The following denote the symbol and color of screening level exceedances:

^ = At or Above 2023 IDEM R2 Long Term Residential Groundwater PLs

Results and IDEM PLs are reported in micrograms per liter ($\mu\text{g}/\text{L}$).

IDEML PLs are based on the IDEM R2, Table 1: Human Health Published Levels with updates.

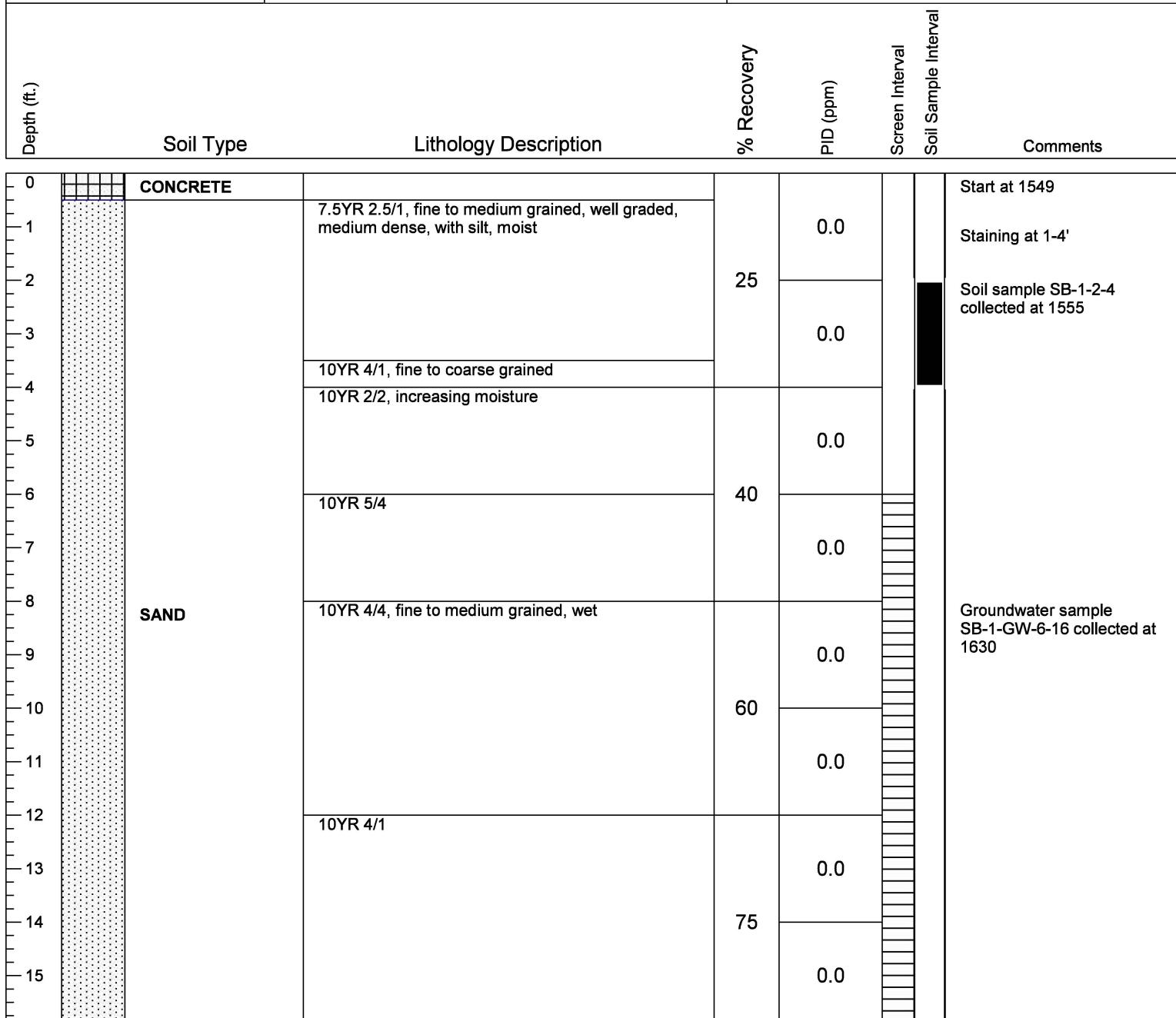
ATTACHMENT A

Soil Boring Logs



| | |
|------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|------------------------------------------|
| Project Number: JX1600.740 Project Name: Jessen Manufacturing Site Address: 1409 W Beardsley Ave. City, State: Elkhart, IN Boring Location: NW Exterior Northing*: Not Measured (NM) Surface Elevation: NM | Date Drilled: 6/21/2023 |
| | Personnel: A. Hicks |
| | Driller: Terracon - C. Bradshaw |
| | Driller License: 4231 |
| | Drilling Method: Direct Push - Dual Tube |

SB-1

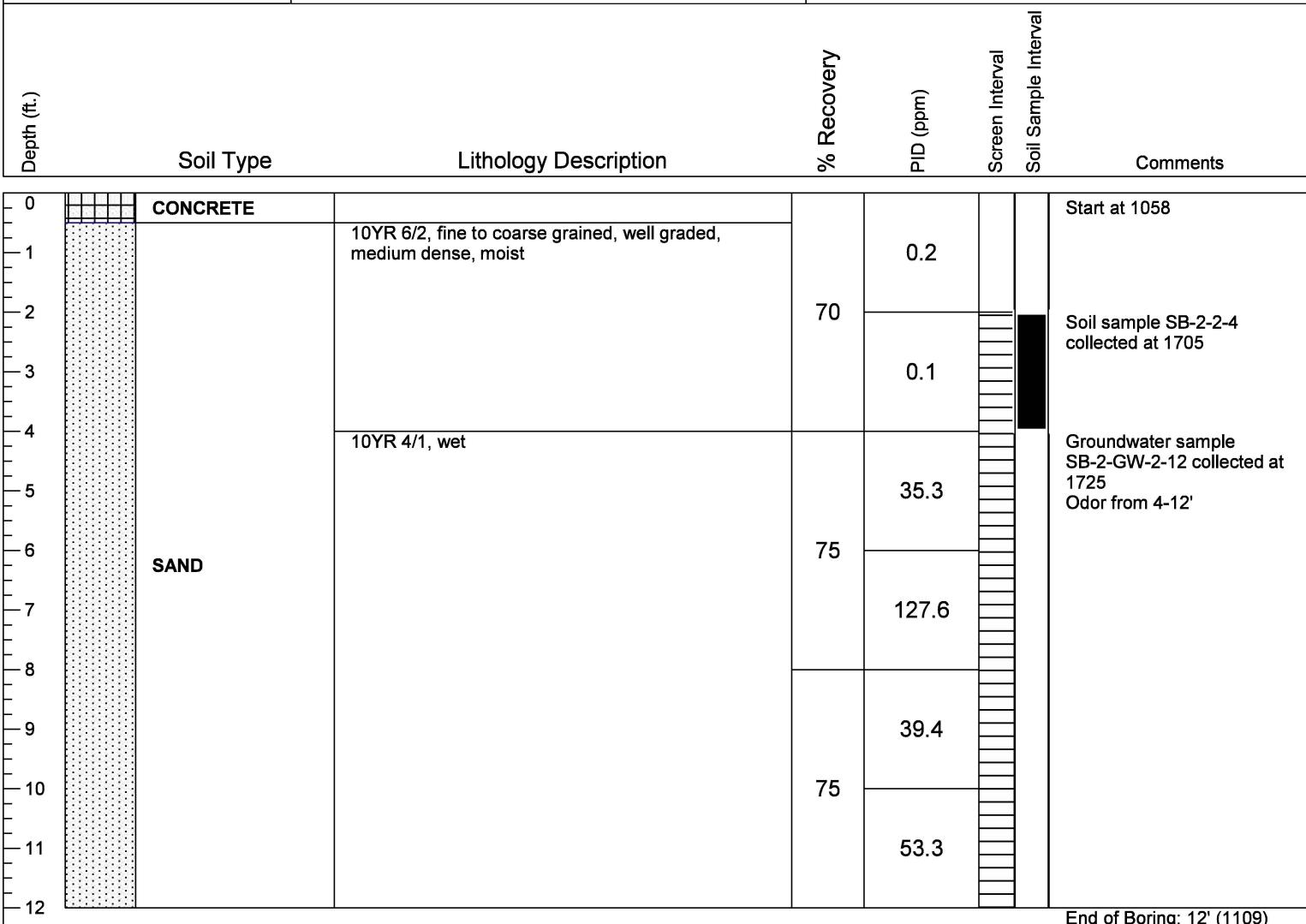


* = Northing, Easting, and Surface Elevation (State Plane or UTM) are estimated, unless specified in the report to have been surveyed.



| | |
|-----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|------------------------------------------|
| Project Number: JX1600.740 Project Name: Jessen Manufacturing Site Address: 1409 W Beardsley Ave. City, State: Elkhart, IN Boring Location: N Exterior Northing*: Not Measured (NM) Surface Elevation: NM | Date Drilled: 6/21/2023 |
| | Personnel: A. Hicks |
| | Driller: Terracon - C. Bradshaw |
| | Driller License: 4231 |
| | Drilling Method: Direct Push - Dual Tube |

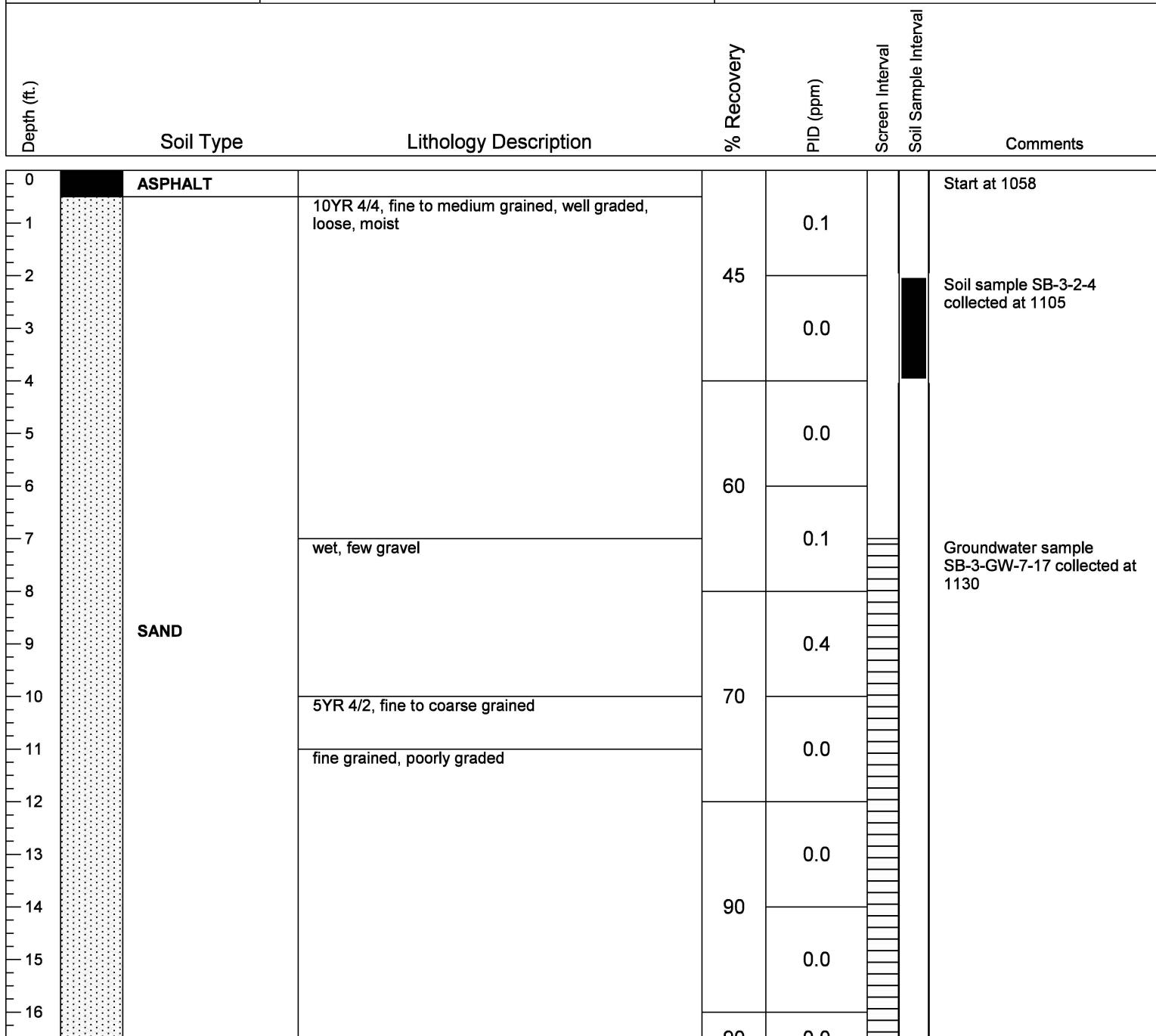
SB-2



* = Northing, Easting, and Surface Elevation (State Plane or UTM) are estimated, unless specified in the report to have been surveyed.

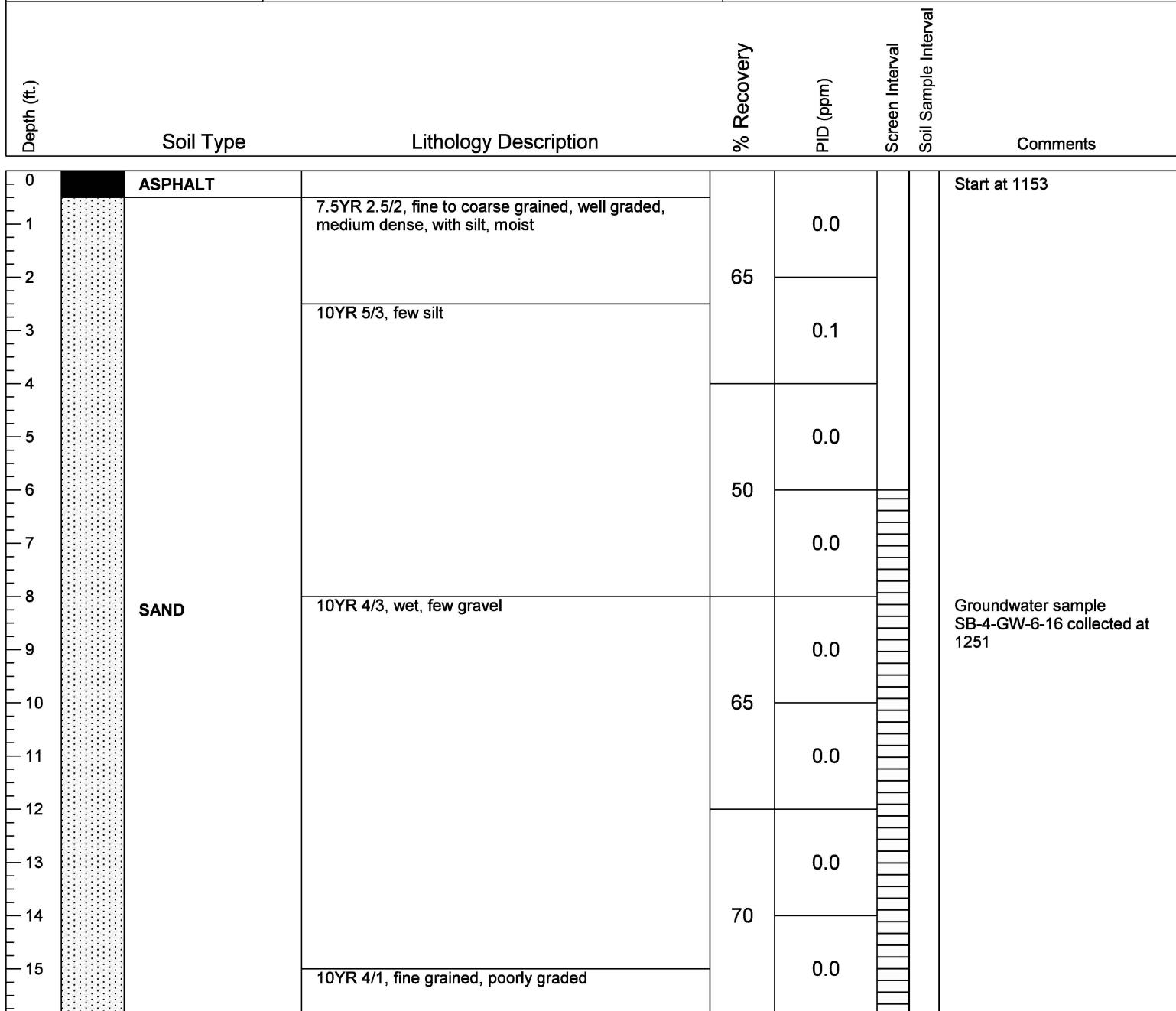


| | | |
|------|-------------------------------------|------------------------------------------|
| | Project Number: JX1600.740 | Date Drilled: 6/21/2023 |
| | Project Name: Jessen Manufacturing | Personnel: A. Hicks |
| | Site Address: 1409 W Beardsley Ave. | Driller: Terracon - C. Bradshaw |
| | City, State: Elkhart, IN | Driller License: 4231 |
| | Boring Location: NE Exterior | Drilling Method: Direct Push - Dual Tube |
| SB-3 | Northing*: Not Measured (NM) | Easting*: NM |
| | Surface Elevation: NM | GW Sample Method: Bailer |



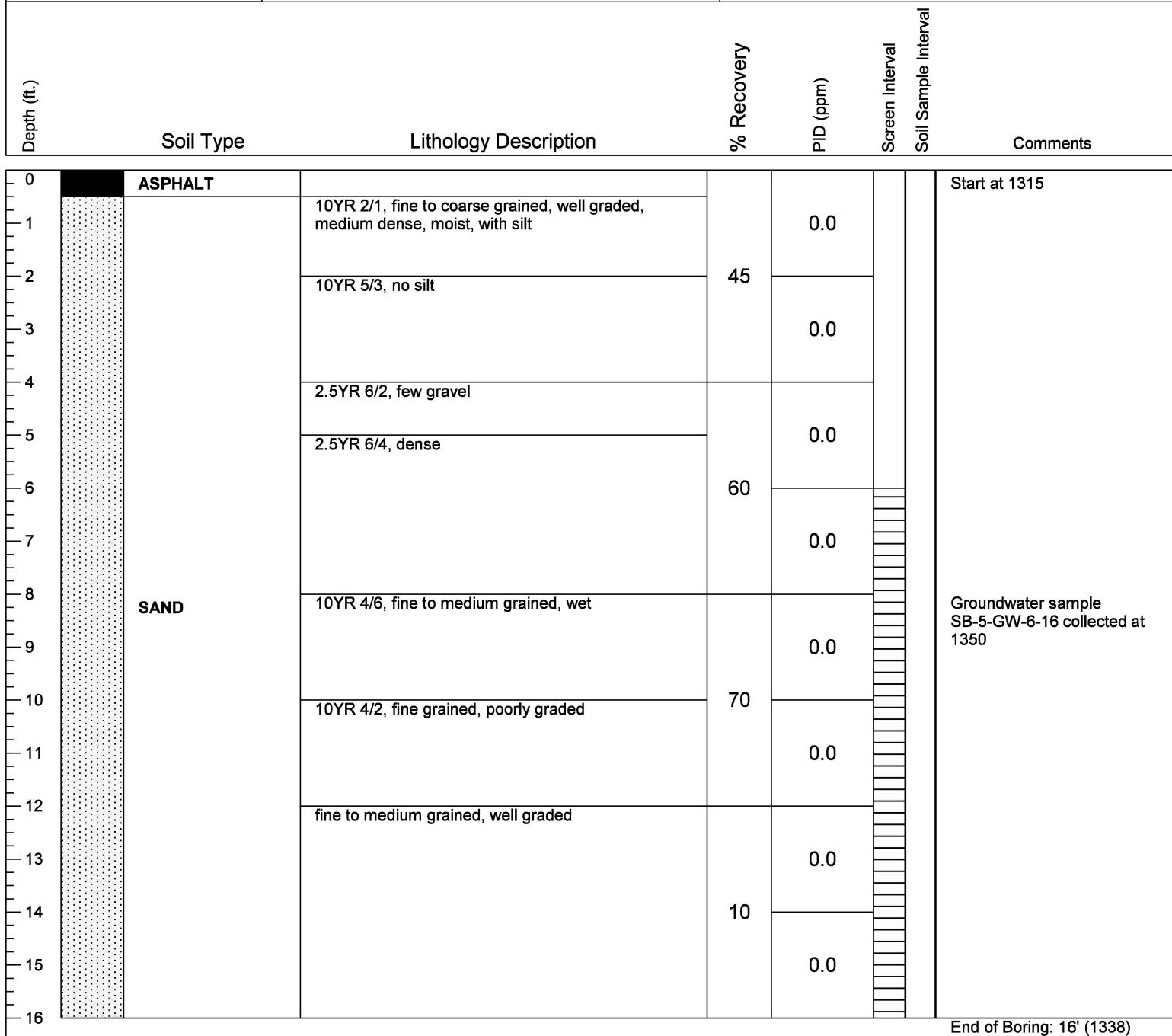
* = Northing, Easting, and Surface Elevation (State Plane or UTM) are estimated, unless specified in the report to have been surveyed.

| | | | | |
|--------------------------------------------------------------------------------------------------------------------------|--------------------|-----------------------|-------------------|-------------------------|
|  August Mack ENVIRONMENTAL | Project Number: | JX1600.740 | Date Drilled: | 6/21/2023 |
| | Project Name: | Jessen Manufacturing | Personnel: | A. Hicks |
| | Site Address: | 1409 W Beardsley Ave. | Driller: | Terracon - C. Bradshaw |
| | City, State: | Elkhart, IN | Driller License: | 4231 |
| | Boring Location: | E Exterior | Drilling Method: | Direct Push - Dual Tube |
| SB-4 | Northing*: | Not Measured (NM) | Easting*: | NM |
| | Surface Elevation: | NM | GW Sample Method: | Bailer |



* = Northing, Easting, and Surface Elevation (State Plane or UTM) are estimated, unless specified in the report to have been surveyed.

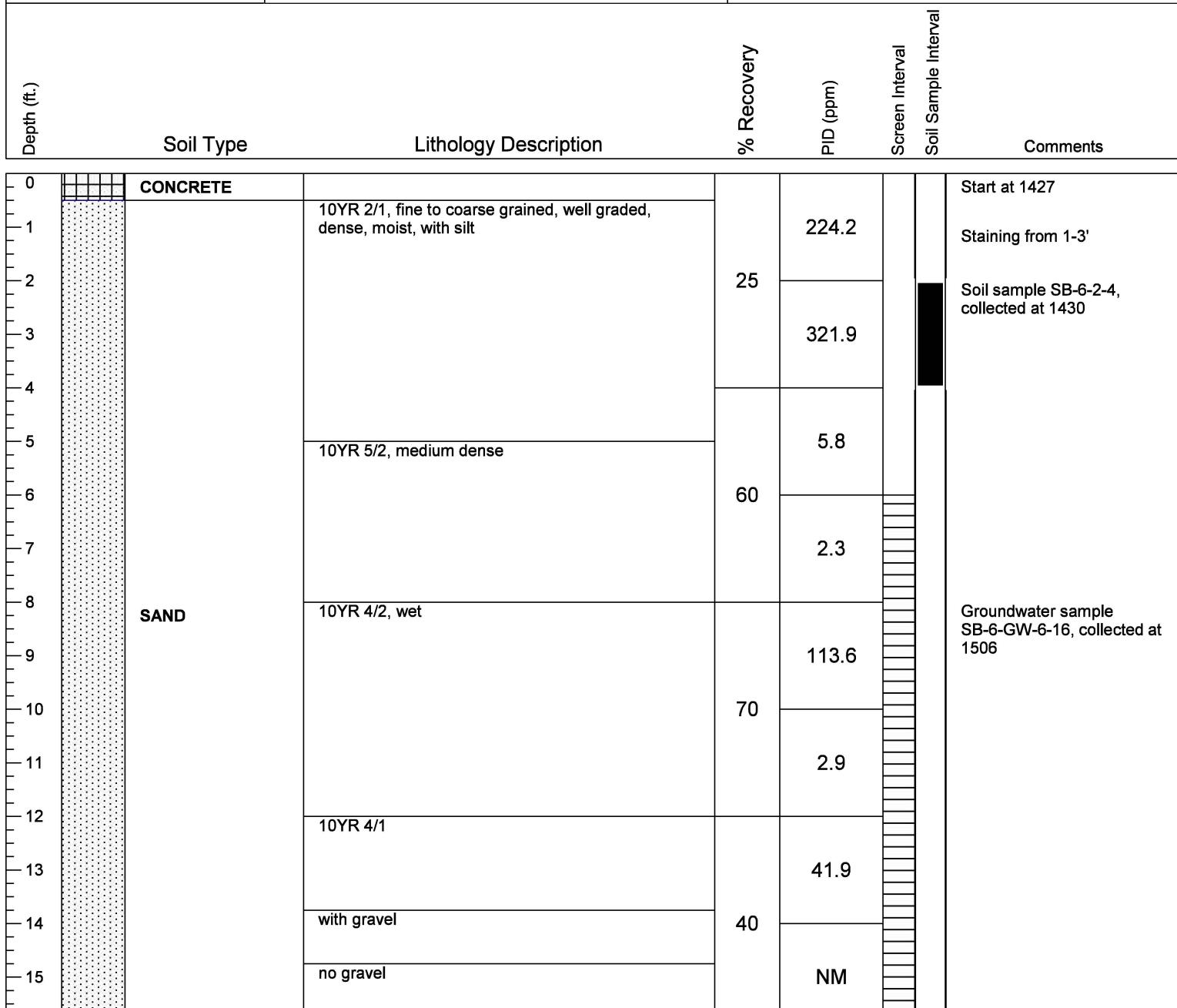
| | | | | |
|--------------------------------------------------------------------------------------------------------------------------------------|--------------------|-----------------------|-------------------|-------------------------|
|  August Mack <small>ENVIRONMENTAL</small> | Project Number: | JX1600.740 | Date Drilled: | 6/21/2023 |
| | Project Name: | Jessen Manufacturing | Personnel: | A. Hicks |
| | Site Address: | 1409 W Beardsley Ave. | Driller: | Terracon - C. Bradshaw |
| | City, State: | Elkhart, IN | Driller License: | 4231 |
| | Boring Location: | S Exterior | Drilling Method: | Direct Push - Dual Tube |
| | Northing*: | Not Measured (NM) | Easting*: | NM |
| | Surface Elevation: | NM | GW Sample Method: | Bailer |





| | |
|-----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|------------------------------------------|
| Project Number: JX1600.740 Project Name: Jessen Manufacturing Site Address: 1409 W Beardsley Ave. City, State: Elkhart, IN Boring Location: W Exterior Northing*: Not Measured (NM) Surface Elevation: NM | Date Drilled: 6/21/2023 |
| | Personnel: A. Hicks |
| | Driller: Terracon - C. Bradshaw |
| | Driller License: 4231 |
| | Drilling Method: Direct Push - Dual Tube |

SB-6



* = Northing, Easting, and Surface Elevation (State Plane or UTM) are estimated, unless specified in the report to have been surveyed.



| | | |
|-------------------------------------------------------------------------------------------------------------------------------------|-------------------|-------------------------|
| Project Number: JX1600.740 Project Name: Jessen Manufacturing Site Address: 1409 W Beardsley Ave. City, State: Elkhart, IN | Date Drilled: | 6/21/2023 |
| | Personnel: | A. Hicks |
| | Driller: | Terracon - F. Smith |
| | Driller License: | 4231 |
| | Drilling Method: | Direct Push - Dual Tube |
| SB-7 Boring Location: Central Exterior Northing*: Not Measured (NM) Surface Elevation: NM | Easting*: | NM |
| | GW Sample Method: | Not Applicable |

| Depth (ft.) | Soil Type | Lithology Description | % Recovery | PID (ppm) | Screen Interval | Soil Sample Interval | Comments |
|-------------|-----------|---------------------------------------------------------------------|------------|-----------|-----------------|----------------------|-----------------------------------------------------------------------------------|
| 0 | | CONCRETE | | | | | Start at 1240 Soil sample SB-7-1-1.5, collected at 1525 Odor from 1-1.5' |
| 1 | | SAND 10YR 2/1, fine to coarse grained, well graded, dense, moist | 100 | 25.0 | | | Refusal: 1.5' (1426) |

ATTACHMENT B

Field Procedures

Soil Sampling Activities

Soil borings were advanced using a Geoprobe® Direct Push Dual-Tube Sampling System (Geoprobe®). Soil borings were advanced to the desired depth required for the investigation. Soil samples were collected continuously from each boring location by using the dual-tube tooling, which includes a disposable acetate sample liner. The sampler was recovered with a 4-foot soil sample collected within an acetate liner inside the barrel. A new acetate liner was used for each sample collected. All reusable equipment that contacted the soil samples was decontaminated with a Liquinox® solution and rinsed with water between each sample collection.

Upon retrieving the 4-foot sections of soil, the samples were divided into 2-foot sections and inspected in the field for evidence of contamination (odors, staining, etc.). Each sample was also screened in the field by headspace analysis using a MiniRae® photoionization detector (PID). All samples were submitted to ENVision for laboratory analysis of VOCs, SVOCs, RCRA 8 Metals, PCBs and Hexavalent Chromium using standard United States Environmental Protection Agency (U.S. EPA) SW-846 analytical methods. SB-5 was also submitted for RCRA 8 Metal analysis.

Groundwater Sampling Activities

Temporary 1-inch groundwater sampling piezometers were installed at SB-1 through SB-6 to allow groundwater to collect for sampling purposes. Groundwater samples were collected from the temporary groundwater sampling points using a bailer. The groundwater samples were transferred to clean, labeled sample containers (provided by the laboratory) and placed on ice in a cooler for preservation in the field. Groundwater samples were submitted to ENVision for laboratory analysis of VOCs, SVOCs and dissolved RCRA 8 Metals using standard United States Environmental Protection Agency (U.S. EPA) SW-846 analytical methods.

Site Restoration Activities

Upon completion of the field sampling activities, the boreholes were abandoned by manually pouring soil cuttings and bentonite into the boring. Study site restoration was completed by patching the surface materials to match pre-investigation conditions.

ATTACHMENT C

Laboratory Results



ENVision Laboratories, Inc.
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Mr. Tyler Zschiedrich
August Mack Environmental
1302 North Meridian Street, Suite 300
Indianapolis, IN 46202

July 7, 2023

ENVision Project Number: 2023-1274
Client Project Name: JX1600

Dear Mr. Zschiedrich,

Please find the attached analytical report for the samples received June 22, 2023. All test methods performed were fully compliant with local, state, and federal EPA methods unless otherwise noted. The project was analyzed as requested on the enclosed chain of custody record. Please review the comments section for additional information about your results or Quality Control data.

The reference for the preservation technique utilized by ENVision Laboratories for Volatile Organics in soil may be found on Table A.1 (p. 42) of Method 5035A: Closed-System Purge-and-Trap and Extraction for Volatile Organics in Soil and Waste Samples, July 2002, Draft Revision 1.

Feel free to contact me if you have any questions or comments regarding your analytical report or service.

Thank you for your business. ENVision Laboratories looks forward to working with you on your next project.

Yours Sincerely,

A handwritten signature in black ink that reads "Cheryl A. Crum".

Cheryl A. Crum
Director of Project Management
ENVision Laboratories, Inc.



Analytical Report

ENVision Laboratories, Inc.
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Client Name: AUGUST MACK ENVIRONMENTAL
Project ID: JX1600
Client Project Manager: TYLER ZSCHIEDRICH
ENVision Project Number: 2023-1274
Analytical Method: EPA 8260
Prep Method: EPA 5030B
Analytical Batch: 062523VW
Client Sample ID: SB-1-GW-6-16 **Sample Collection Date/Time:** 6/21/23 16:30
Envision Sample Number: 23-12287 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: water

| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------|------------------------------|-------------------------------|--------------|
| Acetone | < 100 | 100 | |
| Acrolein | < 1 | 1 | |
| Acrylonitrile | < 0.45 | 1 | 1 |
| Benzene | < 5 | 5 | |
| Bromobenzene | < 5 | 5 | |
| Bromochloromethane | < 5 | 5 | |
| Bromodichloromethane | < 5 | 5 | |
| Bromoform | < 5 | 5 | |
| Bromomethane | < 5 | 5 | |
| n-Butanol | < 50 | 50 | |
| 2-Butanone (MEK) | < 10 | 10 | |
| n-Butylbenzene | < 5 | 5 | |
| sec-Butylbenzene | < 5 | 5 | |
| tert-Butylbenzene | < 5 | 5 | |
| Carbon Disulfide | < 5 | 5 | |
| Carbon Tetrachloride | < 5 | 5 | |
| Chlorobenzene | < 5 | 5 | |
| Chloroethane | < 5 | 5 | |
| 2-Chloroethylvinylether | < 50 | 50 | |
| Chloroform | < 5 | 5 | |
| Chloromethane | < 5 | 5 | |
| 2-Chlorotoluene | < 5 | 5 | |
| 4-Chlorotoluene | < 5 | 5 | |
| 1,2-Dibromo-3-chloropropane | < 1 | 1 | |
| Dibromochloromethane | < 5 | 5 | |
| 1,2-Dibromoethane (EDB) | < 1 | 1 | |
| Dibromomethane | < 5 | 5 | |
| 1,2-Dichlorobenzene | < 5 | 5 | |
| 1,3-Dichlorobenzene | < 5 | 5 | |
| 1,4-Dichlorobenzene | < 5 | 5 | |
| trans-1,4-Dichloro-2-butene | < 1 | 1 | |
| Dichlorodifluoromethane | < 5 | 5 | |



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8260 continued...

| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------------|------------------------------|-------------------------------|--------------|
| 1,1-Dichloroethane | < 5 | 5 | |
| 1,2-Dichloroethane | < 5 | 5 | |
| 1,1-Dichloroethene | < 5 | 5 | |
| cis-1,2-Dichloroethene | < 5 | 5 | |
| trans-1,2-Dichloroethene | < 5 | 5 | |
| 1,2-Dichloropropane | < 5 | 5 | |
| 1,3-Dichloropropane | < 5 | 5 | |
| 2,2-Dichloropropane | < 5 | 5 | |
| 1,1-Dichloropropene | < 5 | 5 | |
| 1,3-Dichloropropene | < 4.1 | 4.1 | |
| Ethylbenzene | < 5 | 5 | |
| Ethyl methacrylate | < 100 | 100 | |
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| n-Hexane | < 10 | 10 | |
| 2-Hexanone | < 10 | 10 | |
| Iodomethane | < 10 | 10 | |
| Isopropylbenzene (Cumene) | < 5 | 5 | |
| p-Isopropyltoluene | < 5 | 5 | |
| Methylene chloride | < 5 | 5 | |
| 4-Methyl-2-pentanone (MIBK) | < 10 | 10 | |
| Methyl-tert-butyl-ether | < 5 | 5 | |
| n-Propylbenzene | < 5 | 5 | |
| Styrene | < 5 | 5 | |
| 1,1,1,2-Tetrachloroethane | < 5 | 5 | |
| 1,1,2,2-Tetrachloroethane | < 0.66 | 1 | 1 |
| Tetrachloroethene | 9.31 | 5 | |
| Toluene | < 5 | 5 | |
| 1,2,3-Trichlorobenzene | < 5 | 5 | |
| 1,2,4-Trichlorobenzene | < 5 | 5 | |
| 1,1,1-Trichloroethane | 5.95 | 5 | |
| 1,1,2-Trichloroethane | < 5 | 5 | |
| Trichloroethene | 8.05 | 5 | |
| Trichlorofluoromethane | < 5 | 5 | |
| 1,2,3-Trichloropropane | < 1 | 1 | |
| 1,2,4-Trimethylbenzene | < 5 | 5 | |
| 1,3,5-Trimethylbenzene | < 5 | 5 | |
| Vinyl acetate | < 10 | 10 | |
| Vinyl chloride | < 2 | 2 | |
| Xylene, M&P | < 5 | 5 | |
| Xylene, Ortho | < 5 | 5 | |
| Xylene (Total) | < 10 | 10 | |
| Dibromofluoromethane (surrogate) | 95% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 97% | | |
| Toluene-d8 (surrogate) | 109% | | |
| 4-bromofluorobenzene (surrogate) | 103% | | |
| Analysis Date/Time: | 6-25-23/14:23 | | |
| Analyst Initials | tjg | | |



Analytical Report

ENVision Laboratories, Inc.
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Client Name: AUGUST MACK ENVIRONMENTAL
Project ID: JX1600
Client Project Manager: TYLER ZSCHIEDRICH
ENVision Project Number: 2023-1274
Analytical Method: EPA 8270 BNA/PAH-SIM
Prep Method: EPA 3520C
BNA Analytical Batch: 062823BW

Client Sample ID: SB-1-GW-6-16 **Sample Collection Date/Time:** 6/21/23 16:30
Envision Sample Number: 23-12287 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: water

| <u>BNA Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------|------------------------------|-------------------------------|--------------|
| Aniline | < 10 | 10 | |
| Benzoic Acid | < 50 | 50 | |
| Benzyl Alcohol | < 20 | 20 | |
| 4-Bromophenylphenyl ether | < 10 | 10 | |
| Butylbenzylphthalate | < 10 | 10 | |
| Carbazole | < 20 | 20 | |
| 4-Chloro-3-methylphenol | < 20 | 20 | |
| 4-Chloroaniline | < 3.2 | 3.2 | |
| bis(2-Chloroethoxy)methane | < 10 | 10 | |
| bis(2-Chloroethyl)ether | < 0.12 | 1.0 | 1 |
| bis(2-Chloroisopropyl)ether | < 10 | 10 | |
| 2-Chloronaphthalene | < 10 | 10 | |
| 2-Chlorophenol | < 10 | 10 | |
| 4-Chlorophenylphenyl ether | < 10 | 10 | |
| Dibenzofuran | < 10 | 10 | |
| 1,2-Dichlorobenzene | < 10 | 10 | |
| 1,3-Dichlorobenzene | < 10 | 10 | |
| 1,4-Dichlorobenzene | < 10 | 10 | |
| 3,3-Dichlorobenzidine | < 1.1 | 1.1 | |
| 2,4-Dichlorophenol | < 10 | 10 | |
| Diethylphthalate | < 10 | 10 | |
| 2,4-Dimethylphenol | < 10 | 10 | |
| Dimethylphthalate | < 10 | 10 | |
| Di-n-butylphthalate | < 10 | 10 | |
| 4,6-Dinitro-2-methylphenol | < 1.2 | 1.2 | |
| 2,4-Dinitrophenol | < 30 | 50 | |
| 2,4-Dinitrotoluene | < 2.0 | 2 | |
| 2,6-Dinitrotoluene | < 10 | 10 | |
| Di-n-octylphthalate | < 10 | 10 | |
| bis(2-Ethylhexyl)phthalate | < 5 | 5 | |
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| Hexachlorobenzene | < 1.0 | 1 | |
| Hexachlorocyclopentadiene | < 25 | 25 | |
| Hexachloroethane | < 5.1 | 5.1 | |



Analytical Report

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| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|----------------------------------|------------------------------|-------------------------------|--------------|
| Isophorone | < 10 | 10 | |
| 2-Methylphenol (o-Cresol) | < 10 | 10 | |
| 3&4-Methylphenol | < 20 | 20 | |
| 2-Nitroaniline | < 50 | 50 | |
| 3-Nitroaniline | < 50 | 50 | |
| 4-Nitroaniline | < 33 | 33 | |
| Nitrobenzene | < 1.2 | 1.2 | |
| 2-Nitrophenol | < 10 | 10 | |
| 4-Nitrophenol | < 50 | 50 | |
| N-Nitroso-di-n-propylamine | < 0.093 | 1.0 | 1 |
| N-Nitrosodiphenylamine | < 10 | 10 | |
| Pentachlorophenol | < 1.0 | 1 | |
| Phenol | < 10 | 10 | |
| 1,2,4-Trichlorobenzene | < 10 | 10 | |
| 2,4,5-Trichlorophenol | < 10 | 10 | |
| 2,4,6-Trichlorophenol | < 9 | 9 | |
| 2-Fluorophenol (surrogate) | 106% | | |
| Phenol-d6 (surrogate) | 49% | | |
| Nitrobenzene-d5 (surrogate) | 81% | | |
| 2-Fluorobiphenyl (surrogate) | 75% | | |
| 2,4,6-Tribromophenol (surrogate) | 97% | | |
| p-Terphenyl-d14 (surrogate) | 92% | | |
| Analysis Date/Time: | 06-28-23/19:15 | | |
| Analyst Initials: | gjd | | |
| Date Extracted: | 6/28/23 | | |
| Initial Sample Volume: | 1000 mL | | |
| Final Volume: | 1.0 mL | | |

PAH-SIM Analytical Batch: 062323PW2

| <u>PAH-SIM Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|--------------------------|------------------------------|-------------------------------|--------------|
| Acenaphthene | < 1.0 | 1.0 | |
| Acenaphthylene | < 1.0 | 1.0 | |
| Anthracene | < 0.10 | 0.10 | |
| Benzo(a)anthracene | < 0.10 | 0.10 | |
| Benzo(a)pyrene | < 0.10 | 0.10 | |
| Benzo(b)fluoranthene | < 0.10 | 0.10 | |
| Benzo(g,h,i)perylene | < 0.10 | 0.10 | |
| Benzo(k)fluoranthene | < 0.10 | 0.10 | |
| Chrysene | < 0.10 | 0.10 | |
| Dibenzo(a,h)anthracene | < 0.029 | 0.029 | |
| Fluoranthene | < 1.0 | 1.0 | |
| Fluorene | < 1.0 | 1.0 | |
| Indeno(1,2,3-cd)pyrene | < 0.022 | 0.022 | |
| 1-methylnaphthalene | < 1.0 | 1.0 | |
| 2-methylnaphthalene | < 1.0 | 1.0 | |
| Naphthalene | < 1.0 | 1.0 | |
| Phenanthrene | < 1.0 | 1.0 | |
| Pyrene | < 1.0 | 1.0 | |
| Analysis Date/Time: | 06-24-23/03:02 | | |
| Analyst Initials | gjd | | |



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 6010

Prep Method: EPA 3010A

Client Sample ID: SB-1-GW-6-16

Sample Collection Date/Time: 6/21/23 16:30

Envision Sample Number: 23-12287

Sample Received Date/Time: 6/22/23 9:25

Sample Matrix: water

| Compounds | Sample Results (ug/L) | Reporting Limit (ug/L) | Flags |
|--------------------|------------------------------|-------------------------------|--------------|
| Arsenic-Dissolved | < 10 | 10 | |
| Barium-Dissolved | < 100 | 100 | |
| Cadmium-Dissolved | < 5 | 5 | |
| Chromium-Dissolved | < 10 | 10 | |
| Lead-Dissolved | < 10 | 10 | |
| Selenium-Dissolved | < 10 | 10 | |
| Silver-Dissolved | < 50 | 50 | |

ICP Analysis Date/Time: 6-27-23/14:05

Analyst Initials: gjd

Date Digested: 6/27/23

Initial Sample Volume: 50 mL

Final Volume: 50 mL

Analytical Batch: 062723icp

Analytical & Prep Method: EPA 7470

| Compounds | Sample Results (ug/L) | Reporting Limit (ug/L) | Flags |
|-------------------|------------------------------|-------------------------------|--------------|
| Mercury-Dissolved | < 2 | 2 | |

Hg Analysis Date/Time: 6/28/23/11:35

Hg Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Volume: 50 mL

Final Volume: 50 mL

Analytical Batch: 062823hg



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL
Project ID: JX1600
Client Project Manager: TYLER ZSCHIEDRICH
ENVision Project Number: 2023-1274
Analytical Method: EPA 8260
Prep Method: EPA 5030B
Analytical Batch: 062523VW
Client Sample ID: SB-2-GW-2-12 **Sample Collection Date/Time:** 6/21/23 17:25
Envision Sample Number: 23-12288 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: water

| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------|------------------------------|-------------------------------|--------------|
| Acetone | < 100 | 100 | |
| Acrolein | < 1 | 1 | |
| Acrylonitrile | < 0.45 | 1 | 1 |
| Benzene | < 5 | 5 | |
| Bromobenzene | < 5 | 5 | |
| Bromochloromethane | < 5 | 5 | |
| Bromodichloromethane | < 5 | 5 | |
| Bromoform | < 5 | 5 | |
| Bromomethane | < 5 | 5 | |
| n-Butanol | < 50 | 50 | |
| 2-Butanone (MEK) | < 10 | 10 | |
| n-Butylbenzene | 24.0 | 5 | |
| sec-Butylbenzene | 13.2 | 5 | |
| tert-Butylbenzene | < 5 | 5 | |
| Carbon Disulfide | < 5 | 5 | |
| Carbon Tetrachloride | < 5 | 5 | |
| Chlorobenzene | < 5 | 5 | |
| Chloroethane | < 5 | 5 | |
| 2-Chloroethylvinylether | < 50 | 50 | |
| Chloroform | < 5 | 5 | |
| Chloromethane | < 5 | 5 | |
| 2-Chlorotoluene | < 5 | 5 | |
| 4-Chlorotoluene | < 5 | 5 | |
| 1,2-Dibromo-3-chloropropane | < 1 | 1 | |
| Dibromochloromethane | < 5 | 5 | |
| 1,2-Dibromoethane (EDB) | < 1 | 1 | |
| Dibromomethane | < 5 | 5 | |
| 1,2-Dichlorobenzene | < 5 | 5 | |
| 1,3-Dichlorobenzene | < 5 | 5 | |
| 1,4-Dichlorobenzene | < 5 | 5 | |
| trans-1,4-Dichloro-2-butene | < 1 | 1 | |
| Dichlorodifluoromethane | < 5 | 5 | |



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8260 continued...

| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------------|------------------------------|-------------------------------|--------------|
| 1,1-Dichloroethane | < 5 | 5 | |
| 1,2-Dichloroethane | < 5 | 5 | |
| 1,1-Dichloroethene | < 5 | 5 | |
| cis-1,2-Dichloroethene | < 5 | 5 | |
| trans-1,2-Dichloroethene | < 5 | 5 | |
| 1,2-Dichloropropane | < 5 | 5 | |
| 1,3-Dichloropropane | < 5 | 5 | |
| 2,2-Dichloropropane | < 5 | 5 | |
| 1,1-Dichloropropene | < 5 | 5 | |
| 1,3-Dichloropropene | < 4.1 | 4.1 | |
| Ethylbenzene | < 5 | 5 | |
| Ethyl methacrylate | < 100 | 100 | |
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| n-Hexane | < 10 | 10 | |
| 2-Hexanone | < 10 | 10 | |
| Iodomethane | < 10 | 10 | |
| Isopropylbenzene (Cumene) | < 5 | 5 | |
| p-Isopropyltoluene | 5.59 | 5 | |
| Methylene chloride | < 5 | 5 | |
| 4-Methyl-2-pentanone (MIBK) | < 10 | 10 | |
| Methyl-tert-butyl-ether | < 5 | 5 | |
| n-Propylbenzene | 7.33 | 5 | |
| Styrene | < 5 | 5 | |
| 1,1,1,2-Tetrachloroethane | < 5 | 5 | |
| 1,1,2,2-Tetrachloroethane | < 0.66 | 1 | 1 |
| Tetrachloroethene | < 5 | 5 | |
| Toluene | < 5 | 5 | |
| 1,2,3-Trichlorobenzene | < 5 | 5 | |
| 1,2,4-Trichlorobenzene | < 5 | 5 | |
| 1,1,1-Trichloroethane | < 5 | 5 | |
| 1,1,2-Trichloroethane | < 5 | 5 | |
| Trichloroethene | < 5 | 5 | |
| Trichlorofluoromethane | < 5 | 5 | |
| 1,2,3-Trichloropropane | < 1 | 1 | |
| 1,2,4-Trimethylbenzene | < 5 | 5 | |
| 1,3,5-Trimethylbenzene | < 5 | 5 | |
| Vinyl acetate | < 10 | 10 | |
| Vinyl chloride | < 2 | 2 | |
| Xylene, M&P | < 5 | 5 | |
| Xylene, Ortho | < 5 | 5 | |
| Xylene (Total) | < 10 | 10 | |
| Dibromofluoromethane (surrogate) | 90% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 94% | | |
| Toluene-d8 (surrogate) | 106% | | |
| 4-bromofluorobenzene (surrogate) | 97% | | |
| Analysis Date/Time: | 6-25-23/14:39 | | |
| Analyst Initials | tjg | | |



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL
Project ID: JX1600
Client Project Manager: TYLER ZSCHIEDRICH
ENVision Project Number: 2023-1274
Analytical Method: EPA 8270 BNA/PAH-SIM
Prep Method: EPA 3520C
BNA Analytical Batch: 062823BW

Client Sample ID: SB-2-GW-2-12 **Sample Collection Date/Time:** 6/21/23 17:25
Envision Sample Number: 23-12288 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: water

| <u>BNA Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------|------------------------------|-------------------------------|--------------|
| Aniline | < 10 | 10 | |
| Benzoic Acid | < 50 | 50 | |
| Benzyl Alcohol | < 20 | 20 | |
| 4-Bromophenylphenyl ether | < 10 | 10 | |
| Butylbenzylphthalate | < 10 | 10 | |
| Carbazole | < 20 | 20 | |
| 4-Chloro-3-methylphenol | < 20 | 20 | |
| 4-Chloroaniline | < 3.2 | 3.2 | |
| bis(2-Chloroethoxy)methane | < 10 | 10 | |
| bis(2-Chloroethyl)ether | < 0.12 | 1.0 | 1 |
| bis(2-Chloroisopropyl)ether | < 10 | 10 | |
| 2-Chloronaphthalene | < 10 | 10 | |
| 2-Chlorophenol | < 10 | 10 | |
| 4-Chlorophenylphenyl ether | < 10 | 10 | |
| Dibenzofuran | < 10 | 10 | |
| 1,2-Dichlorobenzene | < 10 | 10 | |
| 1,3-Dichlorobenzene | < 10 | 10 | |
| 1,4-Dichlorobenzene | < 10 | 10 | |
| 3,3-Dichlorobenzidine | < 1.1 | 1.1 | |
| 2,4-Dichlorophenol | < 10 | 10 | |
| Diethylphthalate | < 10 | 10 | |
| 2,4-Dimethylphenol | < 10 | 10 | |
| Dimethylphthalate | < 10 | 10 | |
| Di-n-butylphthalate | < 10 | 10 | |
| 4,6-Dinitro-2-methylphenol | < 1.2 | 1.2 | |
| 2,4-Dinitrophenol | < 30 | 50 | |
| 2,4-Dinitrotoluene | < 2.0 | 2 | |
| 2,6-Dinitrotoluene | < 10 | 10 | |
| Di-n-octylphthalate | < 10 | 10 | |
| bis(2-Ethylhexyl)phthalate | < 5 | 5 | |
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| Hexachlorobenzene | < 1.0 | 1 | |
| Hexachlorocyclopentadiene | < 25 | 25 | |
| Hexachloroethane | < 5.1 | 5.1 | |



Analytical Report

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8270 Continued...

| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|----------------------------------|------------------------------|-------------------------------|--------------|
| Isophorone | < 10 | 10 | |
| 2-Methylphenol (o-Cresol) | < 10 | 10 | |
| 3&4-Methylphenol | < 20 | 20 | |
| 2-Nitroaniline | < 50 | 50 | |
| 3-Nitroaniline | < 50 | 50 | |
| 4-Nitroaniline | < 33 | 33 | |
| Nitrobenzene | < 1.2 | 1.2 | |
| 2-Nitrophenol | < 10 | 10 | |
| 4-Nitrophenol | < 50 | 50 | |
| N-Nitroso-di-n-propylamine | < 0.093 | 1.0 | 1 |
| N-Nitrosodiphenylamine | < 10 | 10 | |
| Pentachlorophenol | < 1.0 | 1 | |
| Phenol | < 10 | 10 | |
| 1,2,4-Trichlorobenzene | < 10 | 10 | |
| 2,4,5-Trichlorophenol | < 10 | 10 | |
| 2,4,6-Trichlorophenol | < 9 | 9 | |
| 2-Fluorophenol (surrogate) | 93% | | |
| Phenol-d6 (surrogate) | 91% | | |
| Nitrobenzene-d5 (surrogate) | 67% | | |
| 2-Fluorobiphenyl (surrogate) | 60% | | |
| 2,4,6-Tribromophenol (surrogate) | 93% | | |
| p-Terphenyl-d14 (surrogate) | 59% | | |
| Analysis Date/Time: | 06-28-23/19:41 | | |
| Analyst Initials: | gjd | | |
| Date Extracted: | 6/28/23 | | |
| Initial Sample Volume: | 1000 mL | | |
| Final Volume: | 1.0 mL | | |

PAH-SIM Analytical Batch: 062323PW2

| <u>PAH-SIM Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|--------------------------|------------------------------|-------------------------------|--------------|
| Acenaphthene | < 1.0 | 1.0 | |
| Acenaphthylene | < 1.0 | 1.0 | |
| Anthracene | < 0.10 | 0.10 | |
| Benzo(a)anthracene | < 0.10 | 0.10 | |
| Benzo(a)pyrene | < 0.10 | 0.10 | |
| Benzo(b)fluoranthene | < 0.10 | 0.10 | |
| Benzo(g,h,i)perylene | < 0.10 | 0.10 | |
| Benzo(k)fluoranthene | < 0.10 | 0.10 | |
| Chrysene | < 0.10 | 0.10 | |
| Dibenzo(a,h)anthracene | < 0.029 | 0.029 | |
| Fluoranthene | < 1.0 | 1.0 | |
| Fluorene | < 1.0 | 1.0 | |
| Indeno(1,2,3-cd)pyrene | < 0.022 | 0.022 | |
| 1-methylnaphthalene | < 1.0 | 1.0 | |
| 2-methylnaphthalene | < 1.0 | 1.0 | |
| Naphthalene | < 1.0 | 1.0 | |
| Phenanthrene | < 1.0 | 1.0 | |
| Pyrene | < 1.0 | 1.0 | |
| Analysis Date/Time: | 06-24-23/03:24 | | |
| Analyst Initials | gjd | | |



Analytical Report

ENVision Laboratories, Inc.
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Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 6010

Prep Method: EPA 3010A

Client Sample ID: SB-2-GW-2-12 **Sample Collection Date/Time:** 6/21/23 17:25

Envision Sample Number: 23-12288 **Sample Received Date/Time:** 6/22/23 9:25

Sample Matrix: water

| Compounds | Sample Results (ug/L) | Reporting Limit (ug/L) | Flags |
|--------------------|------------------------------|-------------------------------|--------------|
| Arsenic-Dissolved | < 10 | 10 | |
| Barium-Dissolved | < 100 | 100 | |
| Cadmium-Dissolved | < 5 | 5 | |
| Chromium-Dissolved | < 10 | 10 | |
| Lead-Dissolved | < 10 | 10 | |
| Selenium-Dissolved | < 10 | 10 | |
| Silver-Dissolved | < 50 | 50 | |

ICP Analysis Date/Time: 6-27-23/14:08

Analyst Initials: gjd

Date Digested: 6/27/23

Initial Sample Volume: 50 mL

Final Volume: 50 mL

Analytical Batch: 062723icp

Analytical & Prep Method: EPA 7470

| Compounds | Sample Results (ug/L) | Reporting Limit (ug/L) | Flags |
|-------------------|------------------------------|-------------------------------|--------------|
| Mercury-Dissolved | < 2 | 2 | |

Hg Analysis Date/Time: 6/28/23/11:37

Hg Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Volume: 50 mL

Final Volume: 50 mL

Analytical Batch: 062823hg



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL
Project ID: JX1600
Client Project Manager: TYLER ZSCHIEDRICH
ENVision Project Number: 2023-1274
Analytical Method: EPA 8260
Prep Method: EPA 5030B
Analytical Batch: 062523VW
Client Sample ID: SB-3-GW-7-17 **Sample Collection Date/Time:** 6/21/23 11:30
Envision Sample Number: 23-12289 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: water

| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------|------------------------------|-------------------------------|--------------|
| Acetone | < 100 | 100 | |
| Acrolein | < 1 | 1 | |
| Acrylonitrile | < 0.45 | 1 | 1 |
| Benzene | < 5 | 5 | |
| Bromobenzene | < 5 | 5 | |
| Bromochloromethane | < 5 | 5 | |
| Bromodichloromethane | < 5 | 5 | |
| Bromoform | < 5 | 5 | |
| Bromomethane | < 5 | 5 | |
| n-Butanol | < 50 | 50 | |
| 2-Butanone (MEK) | < 10 | 10 | |
| n-Butylbenzene | < 5 | 5 | |
| sec-Butylbenzene | < 5 | 5 | |
| tert-Butylbenzene | < 5 | 5 | |
| Carbon Disulfide | < 5 | 5 | |
| Carbon Tetrachloride | < 5 | 5 | |
| Chlorobenzene | < 5 | 5 | |
| Chloroethane | < 5 | 5 | |
| 2-Chloroethylvinylether | < 50 | 50 | |
| Chloroform | < 5 | 5 | |
| Chloromethane | < 5 | 5 | |
| 2-Chlorotoluene | < 5 | 5 | |
| 4-Chlorotoluene | < 5 | 5 | |
| 1,2-Dibromo-3-chloropropane | < 1 | 1 | |
| Dibromochloromethane | < 5 | 5 | |
| 1,2-Dibromoethane (EDB) | < 1 | 1 | |
| Dibromomethane | < 5 | 5 | |
| 1,2-Dichlorobenzene | < 5 | 5 | |
| 1,3-Dichlorobenzene | < 5 | 5 | |
| 1,4-Dichlorobenzene | < 5 | 5 | |
| trans-1,4-Dichloro-2-butene | < 1 | 1 | |
| Dichlorodifluoromethane | < 5 | 5 | |



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| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------------|------------------------------|-------------------------------|--------------|
| 1,1-Dichloroethane | < 5 | 5 | |
| 1,2-Dichloroethane | < 5 | 5 | |
| 1,1-Dichloroethene | < 5 | 5 | |
| cis-1,2-Dichloroethene | 12.2 | 5 | |
| trans-1,2-Dichloroethene | < 5 | 5 | |
| 1,2-Dichloropropane | < 5 | 5 | |
| 1,3-Dichloropropane | < 5 | 5 | |
| 2,2-Dichloropropane | < 5 | 5 | |
| 1,1-Dichloropropene | < 5 | 5 | |
| 1,3-Dichloropropene | < 4.1 | 4.1 | |
| Ethylbenzene | < 5 | 5 | |
| Ethyl methacrylate | < 100 | 100 | |
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| n-Hexane | < 10 | 10 | |
| 2-Hexanone | < 10 | 10 | |
| Iodomethane | < 10 | 10 | |
| Isopropylbenzene (Cumene) | < 5 | 5 | |
| p-Isopropyltoluene | < 5 | 5 | |
| Methylene chloride | < 5 | 5 | |
| 4-Methyl-2-pentanone (MIBK) | < 10 | 10 | |
| Methyl-tert-butyl-ether | < 5 | 5 | |
| n-Propylbenzene | < 5 | 5 | |
| Styrene | < 5 | 5 | |
| 1,1,1,2-Tetrachloroethane | < 5 | 5 | |
| 1,1,2,2-Tetrachloroethane | < 0.66 | 1 | 1 |
| Tetrachloroethene | < 5 | 5 | |
| Toluene | < 5 | 5 | |
| 1,2,3-Trichlorobenzene | < 5 | 5 | |
| 1,2,4-Trichlorobenzene | < 5 | 5 | |
| 1,1,1-Trichloroethane | < 5 | 5 | |
| 1,1,2-Trichloroethane | < 5 | 5 | |
| Trichloroethene | < 5 | 5 | |
| Trichlorofluoromethane | < 5 | 5 | |
| 1,2,3-Trichloropropane | < 1 | 1 | |
| 1,2,4-Trimethylbenzene | < 5 | 5 | |
| 1,3,5-Trimethylbenzene | < 5 | 5 | |
| Vinyl acetate | < 10 | 10 | |
| Vinyl chloride | < 2 | 2 | |
| Xylene, M&P | < 5 | 5 | |
| Xylene, Ortho | < 5 | 5 | |
| Xylene (Total) | < 10 | 10 | |
| Dibromofluoromethane (surrogate) | 90% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 93% | | |
| Toluene-d8 (surrogate) | 106% | | |
| 4-bromofluorobenzene (surrogate) | 112% | | |
| Analysis Date/Time: | 6-25-23/14:55 | | |
| Analyst Initials | tjg | | |



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL
Project ID: JX1600
Client Project Manager: TYLER ZSCHIEDRICH
ENVision Project Number: 2023-1274
Analytical Method: EPA 8270 BNA/PAH-SIM
Prep Method: EPA 3520C
BNA Analytical Batch: 062823BW

Client Sample ID: SB-3-GW-7-17 **Sample Collection Date/Time:** 6/21/23 11:30
Envision Sample Number: 23-12289 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: water

| <u>BNA Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------|------------------------------|-------------------------------|--------------|
| Aniline | < 10 | 10 | |
| Benzoic Acid | < 50 | 50 | |
| Benzyl Alcohol | < 20 | 20 | |
| 4-Bromophenylphenyl ether | < 10 | 10 | |
| Butylbenzylphthalate | < 10 | 10 | |
| Carbazole | < 20 | 20 | |
| 4-Chloro-3-methylphenol | < 20 | 20 | |
| 4-Chloroaniline | < 3.2 | 3.2 | |
| bis(2-Chloroethoxy)methane | < 10 | 10 | |
| bis(2-Chloroethyl)ether | < 0.12 | 1.0 | 1 |
| bis(2-Chloroisopropyl)ether | < 10 | 10 | |
| 2-Chloronaphthalene | < 10 | 10 | |
| 2-Chlorophenol | < 10 | 10 | |
| 4-Chlorophenylphenyl ether | < 10 | 10 | |
| Dibenzofuran | < 10 | 10 | |
| 1,2-Dichlorobenzene | < 10 | 10 | |
| 1,3-Dichlorobenzene | < 10 | 10 | |
| 1,4-Dichlorobenzene | < 10 | 10 | |
| 3,3-Dichlorobenzidine | < 1.1 | 1.1 | |
| 2,4-Dichlorophenol | < 10 | 10 | |
| Diethylphthalate | < 10 | 10 | |
| 2,4-Dimethylphenol | < 10 | 10 | |
| Dimethylphthalate | < 10 | 10 | |
| Di-n-butylphthalate | < 10 | 10 | |
| 4,6-Dinitro-2-methylphenol | < 1.2 | 1.2 | |
| 2,4-Dinitrophenol | < 30 | 50 | |
| 2,4-Dinitrotoluene | < 2.0 | 2 | |
| 2,6-Dinitrotoluene | < 10 | 10 | |
| Di-n-octylphthalate | < 10 | 10 | |
| bis(2-Ethylhexyl)phthalate | < 5 | 5 | |
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| Hexachlorobenzene | < 1.0 | 1 | |
| Hexachlorocyclopentadiene | < 25 | 25 | |
| Hexachloroethane | < 5.1 | 5.1 | |



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| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|----------------------------------|------------------------------|-------------------------------|--------------|
| Isophorone | < 10 | 10 | |
| 2-Methylphenol (o-Cresol) | < 10 | 10 | |
| 3&4-Methylphenol | < 20 | 20 | |
| 2-Nitroaniline | < 50 | 50 | |
| 3-Nitroaniline | < 50 | 50 | |
| 4-Nitroaniline | < 33 | 33 | |
| Nitrobenzene | < 1.2 | 1.2 | |
| 2-Nitrophenol | < 10 | 10 | |
| 4-Nitrophenol | < 50 | 50 | |
| N-Nitroso-di-n-propylamine | < 0.093 | 1.0 | 1 |
| N-Nitrosodiphenylamine | < 10 | 10 | |
| Pentachlorophenol | < 1.0 | 1 | |
| Phenol | < 10 | 10 | |
| 1,2,4-Trichlorobenzene | < 10 | 10 | |
| 2,4,5-Trichlorophenol | < 10 | 10 | |
| 2,4,6-Trichlorophenol | < 9 | 9 | |
| 2-Fluorophenol (surrogate) | 90% | | |
| Phenol-d6 (surrogate) | 112% | | |
| Nitrobenzene-d5 (surrogate) | 83% | | |
| 2-Fluorobiphenyl (surrogate) | 76% | | |
| 2,4,6-Tribromophenol (surrogate) | 100% | | |
| p-Terphenyl-d14 (surrogate) | 95% | | |
| Analysis Date/Time: | 06-28-23/20:06 | | |
| Analyst Initials: | gjd | | |
| Date Extracted: | 6/28/23 | | |
| Initial Sample Volume: | 1000 mL | | |
| Final Volume: | 1.0 mL | | |

PAH-SIM Analytical Batch: 062323PW2

| <u>PAH-SIM Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|--------------------------|------------------------------|-------------------------------|--------------|
| Acenaphthene | < 1.0 | 1.0 | |
| Acenaphthylene | < 1.0 | 1.0 | |
| Anthracene | < 0.10 | 0.10 | |
| Benzo(a)anthracene | < 0.10 | 0.10 | |
| Benzo(a)pyrene | < 0.10 | 0.10 | |
| Benzo(b)fluoranthene | < 0.10 | 0.10 | |
| Benzo(g,h,i)perylene | < 0.10 | 0.10 | |
| Benzo(k)fluoranthene | < 0.10 | 0.10 | |
| Chrysene | < 0.10 | 0.10 | |
| Dibenzo(a,h)anthracene | < 0.029 | 0.029 | |
| Fluoranthene | < 1.0 | 1.0 | |
| Fluorene | < 1.0 | 1.0 | |
| Indeno(1,2,3-cd)pyrene | < 0.022 | 0.022 | |
| 1-methylnaphthalene | < 1.0 | 1.0 | |
| 2-methylnaphthalene | < 1.0 | 1.0 | |
| Naphthalene | < 1.0 | 1.0 | |
| Phenanthrene | < 1.0 | 1.0 | |
| Pyrene | < 1.0 | 1.0 | |
| Analysis Date/Time: | 06-24-23/03:46 | | |
| Analyst Initials | gjd | | |



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 6010

Prep Method: EPA 3010A

Client Sample ID: SB-3-GW-7-17

Sample Collection Date/Time: 6/21/23 11:30

Envision Sample Number: 23-12289

Sample Received Date/Time: 6/22/23 9:25

Sample Matrix: water

| Compounds | Sample Results (ug/L) | Reporting Limit (ug/L) | Flags |
|--------------------|------------------------------|-------------------------------|--------------|
| Arsenic-Dissolved | < 10 | 10 | |
| Barium-Dissolved | < 100 | 100 | |
| Cadmium-Dissolved | < 5 | 5 | |
| Chromium-Dissolved | < 10 | 10 | |
| Lead-Dissolved | < 10 | 10 | |
| Selenium-Dissolved | < 10 | 10 | |
| Silver-Dissolved | < 50 | 50 | |

ICP Analysis Date/Time: 6-27-23/14:10

Analyst Initials: gjd

Date Digested: 6/27/23

Initial Sample Volume: 50 mL

Final Volume: 50 mL

Analytical Batch: 062723icp

Analytical & Prep Method: EPA 7470

| Compounds | Sample Results (ug/L) | Reporting Limit (ug/L) | Flags |
|-------------------|------------------------------|-------------------------------|--------------|
| Mercury-Dissolved | < 2 | 2 | |

Hg Analysis Date/Time: 6/28/23/11:38

Hg Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Volume: 50 mL

Final Volume: 50 mL

Analytical Batch: 062823hg



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL
Project ID: JX1600
Client Project Manager: TYLER ZSCHIEDRICH
ENVision Project Number: 2023-1274
Analytical Method: EPA 8260
Prep Method: EPA 5030B
Analytical Batch: 062523VW
Client Sample ID: SB-4-GW-6-16 **Sample Collection Date/Time:** 6/21/23 12:51
Envision Sample Number: 23-12290 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: water

| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------|------------------------------|-------------------------------|--------------|
| Acetone | < 100 | 100 | |
| Acrolein | < 1 | 1 | |
| Acrylonitrile | < 0.45 | 1 | 1 |
| Benzene | < 5 | 5 | |
| Bromobenzene | < 5 | 5 | |
| Bromochloromethane | < 5 | 5 | |
| Bromodichloromethane | < 5 | 5 | |
| Bromoform | < 5 | 5 | |
| Bromomethane | < 5 | 5 | |
| n-Butanol | < 50 | 50 | |
| 2-Butanone (MEK) | < 10 | 10 | |
| n-Butylbenzene | < 5 | 5 | |
| sec-Butylbenzene | < 5 | 5 | |
| tert-Butylbenzene | < 5 | 5 | |
| Carbon Disulfide | < 5 | 5 | |
| Carbon Tetrachloride | < 5 | 5 | |
| Chlorobenzene | < 5 | 5 | |
| Chloroethane | < 5 | 5 | |
| 2-Chloroethylvinylether | < 50 | 50 | |
| Chloroform | < 5 | 5 | |
| Chloromethane | < 5 | 5 | |
| 2-Chlorotoluene | < 5 | 5 | |
| 4-Chlorotoluene | < 5 | 5 | |
| 1,2-Dibromo-3-chloropropane | < 1 | 1 | |
| Dibromochloromethane | < 5 | 5 | |
| 1,2-Dibromoethane (EDB) | < 1 | 1 | |
| Dibromomethane | < 5 | 5 | |
| 1,2-Dichlorobenzene | < 5 | 5 | |
| 1,3-Dichlorobenzene | < 5 | 5 | |
| 1,4-Dichlorobenzene | < 5 | 5 | |
| trans-1,4-Dichloro-2-butene | < 1 | 1 | |
| Dichlorodifluoromethane | < 5 | 5 | |



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| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------------|------------------------------|-------------------------------|--------------|
| 1,1-Dichloroethane | < 5 | 5 | |
| 1,2-Dichloroethane | < 5 | 5 | |
| 1,1-Dichloroethene | < 5 | 5 | |
| cis-1,2-Dichloroethene | 11.7 | 5 | |
| trans-1,2-Dichloroethene | < 5 | 5 | |
| 1,2-Dichloropropane | < 5 | 5 | |
| 1,3-Dichloropropane | < 5 | 5 | |
| 2,2-Dichloropropane | < 5 | 5 | |
| 1,1-Dichloropropene | < 5 | 5 | |
| 1,3-Dichloropropene | < 4.1 | 4.1 | |
| Ethylbenzene | < 5 | 5 | |
| Ethyl methacrylate | < 100 | 100 | |
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| n-Hexane | < 10 | 10 | |
| 2-Hexanone | < 10 | 10 | |
| Iodomethane | < 10 | 10 | |
| Isopropylbenzene (Cumene) | < 5 | 5 | |
| p-Isopropyltoluene | < 5 | 5 | |
| Methylene chloride | < 5 | 5 | |
| 4-Methyl-2-pentanone (MIBK) | < 10 | 10 | |
| Methyl-tert-butyl-ether | < 5 | 5 | |
| n-Propylbenzene | < 5 | 5 | |
| Styrene | < 5 | 5 | |
| 1,1,1,2-Tetrachloroethane | < 5 | 5 | |
| 1,1,2,2-Tetrachloroethane | < 0.66 | 1 | 1 |
| Tetrachloroethene | 15.1 | 5 | |
| Toluene | < 5 | 5 | |
| 1,2,3-Trichlorobenzene | < 5 | 5 | |
| 1,2,4-Trichlorobenzene | < 5 | 5 | |
| 1,1,1-Trichloroethane | < 5 | 5 | |
| 1,1,2-Trichloroethane | < 5 | 5 | |
| Trichloroethene | < 5 | 5 | |
| Trichlorofluoromethane | < 5 | 5 | |
| 1,2,3-Trichloropropane | < 1 | 1 | |
| 1,2,4-Trimethylbenzene | < 5 | 5 | |
| 1,3,5-Trimethylbenzene | < 5 | 5 | |
| Vinyl acetate | < 10 | 10 | |
| Vinyl chloride | < 2 | 2 | |
| Xylene, M&P | < 5 | 5 | |
| Xylene, Ortho | < 5 | 5 | |
| Xylene (Total) | < 10 | 10 | |
| Dibromofluoromethane (surrogate) | 96% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 96% | | |
| Toluene-d8 (surrogate) | 109% | | |
| 4-bromofluorobenzene (surrogate) | 102% | | |
| Analysis Date/Time: | 6-25-23/15:27 | | |
| Analyst Initials | tjg | | |



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL
Project ID: JX1600
Client Project Manager: TYLER ZSCHIEDRICH
ENVision Project Number: 2023-1274
Analytical Method: EPA 8270 BNA/PAH-SIM
Prep Method: EPA 3520C
BNA Analytical Batch: 062823BW

Client Sample ID: SB-4-GW-6-16 **Sample Collection Date/Time:** 6/21/23 12:51
Envision Sample Number: 23-12290 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: water

| <u>BNA Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------|------------------------------|-------------------------------|--------------|
| Aniline | < 10 | 10 | |
| Benzoic Acid | < 50 | 50 | |
| Benzyl Alcohol | < 20 | 20 | |
| 4-Bromophenylphenyl ether | < 10 | 10 | |
| Butylbenzylphthalate | < 10 | 10 | |
| Carbazole | < 20 | 20 | |
| 4-Chloro-3-methylphenol | < 20 | 20 | |
| 4-Chloroaniline | < 3.2 | 3.2 | |
| bis(2-Chloroethoxy)methane | < 10 | 10 | |
| bis(2-Chloroethyl)ether | < 0.12 | 1.0 | 1 |
| bis(2-Chloroisopropyl)ether | < 10 | 10 | |
| 2-Chloronaphthalene | < 10 | 10 | |
| 2-Chlorophenol | < 10 | 10 | |
| 4-Chlorophenylphenyl ether | < 10 | 10 | |
| Dibenzofuran | < 10 | 10 | |
| 1,2-Dichlorobenzene | < 10 | 10 | |
| 1,3-Dichlorobenzene | < 10 | 10 | |
| 1,4-Dichlorobenzene | < 10 | 10 | |
| 3,3-Dichlorobenzidine | < 1.1 | 1.1 | |
| 2,4-Dichlorophenol | < 10 | 10 | |
| Diethylphthalate | < 10 | 10 | |
| 2,4-Dimethylphenol | < 10 | 10 | |
| Dimethylphthalate | < 10 | 10 | |
| Di-n-butylphthalate | < 10 | 10 | |
| 4,6-Dinitro-2-methylphenol | < 1.2 | 1.2 | |
| 2,4-Dinitrophenol | < 30 | 50 | |
| 2,4-Dinitrotoluene | < 2.0 | 2 | |
| 2,6-Dinitrotoluene | < 10 | 10 | |
| Di-n-octylphthalate | < 10 | 10 | |
| bis(2-Ethylhexyl)phthalate | < 5 | 5 | |
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| Hexachlorobenzene | < 1.0 | 1 | |
| Hexachlorocyclopentadiene | < 25 | 25 | |
| Hexachloroethane | < 5.1 | 5.1 | |



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| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|----------------------------------|------------------------------|-------------------------------|--------------|
| Isophorone | < 10 | 10 | |
| 2-Methylphenol (o-Cresol) | < 10 | 10 | |
| 3&4-Methylphenol | < 20 | 20 | |
| 2-Nitroaniline | < 50 | 50 | |
| 3-Nitroaniline | < 50 | 50 | |
| 4-Nitroaniline | < 33 | 33 | |
| Nitrobenzene | < 1.2 | 1.2 | |
| 2-Nitrophenol | < 10 | 10 | |
| 4-Nitrophenol | < 50 | 50 | |
| N-Nitroso-di-n-propylamine | < 0.093 | 1.0 | 1 |
| N-Nitrosodiphenylamine | < 10 | 10 | |
| Pentachlorophenol | < 1.0 | 1 | |
| Phenol | < 10 | 10 | |
| 1,2,4-Trichlorobenzene | < 10 | 10 | |
| 2,4,5-Trichlorophenol | < 10 | 10 | |
| 2,4,6-Trichlorophenol | < 9 | 9 | |
| 2-Fluorophenol (surrogate) | 80% | | |
| Phenol-d6 (surrogate) | 92% | | |
| Nitrobenzene-d5 (surrogate) | 66% | | |
| 2-Fluorobiphenyl (surrogate) | 63% | | |
| 2,4,6-Tribromophenol (surrogate) | 80% | | |
| p-Terphenyl-d14 (surrogate) | 76% | | |
| Analysis Date/Time: | 06-28-23/20:32 | | |
| Analyst Initials: | gjd | | |
| Date Extracted: | 6/28/23 | | |
| Initial Sample Volume: | 1000 mL | | |
| Final Volume: | 1.0 mL | | |

PAH-SIM Analytical Batch: 062323PW2

| <u>PAH-SIM Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|--------------------------|------------------------------|-------------------------------|--------------|
| Acenaphthene | < 1.0 | 1.0 | |
| Acenaphthylene | < 1.0 | 1.0 | |
| Anthracene | < 0.10 | 0.10 | |
| Benzo(a)anthracene | < 0.10 | 0.10 | |
| Benzo(a)pyrene | < 0.10 | 0.10 | |
| Benzo(b)fluoranthene | < 0.10 | 0.10 | |
| Benzo(g,h,i)perylene | < 0.10 | 0.10 | |
| Benzo(k)fluoranthene | < 0.10 | 0.10 | |
| Chrysene | < 0.10 | 0.10 | |
| Dibenzo(a,h)anthracene | < 0.029 | 0.029 | |
| Fluoranthene | < 1.0 | 1.0 | |
| Fluorene | < 1.0 | 1.0 | |
| Indeno(1,2,3-cd)pyrene | < 0.022 | 0.022 | |
| 1-methylnaphthalene | < 1.0 | 1.0 | |
| 2-methylnaphthalene | < 1.0 | 1.0 | |
| Naphthalene | < 1.0 | 1.0 | |
| Phenanthrene | < 1.0 | 1.0 | |
| Pyrene | < 1.0 | 1.0 | |
| Analysis Date/Time: | 06-24-23/04:08 | | |
| Analyst Initials | gjd | | |



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 6010

Prep Method: EPA 3010A

Client Sample ID: SB-4-GW-6-16

Sample Collection Date/Time: 6/21/23 12:51

Envision Sample Number: 23-12290

Sample Received Date/Time: 6/22/23 9:25

Sample Matrix: water

| Compounds | Sample Results (ug/L) | Reporting Limit (ug/L) | Flags |
|--------------------|------------------------------|-------------------------------|--------------|
| Arsenic-Dissolved | < 10 | 10 | |
| Barium-Dissolved | < 100 | 100 | |
| Cadmium-Dissolved | < 5 | 5 | |
| Chromium-Dissolved | < 10 | 10 | |
| Lead-Dissolved | < 10 | 10 | |
| Selenium-Dissolved | < 10 | 10 | |
| Silver-Dissolved | < 50 | 50 | |

ICP Analysis Date/Time: 6-27-23/14:12

Analyst Initials: gjd

Date Digested: 6/27/23

Initial Sample Volume: 50 mL

Final Volume: 50 mL

Analytical Batch: 062723icp

Analytical & Prep Method: EPA 7470

| Compounds | Sample Results (ug/L) | Reporting Limit (ug/L) | Flags |
|-------------------|------------------------------|-------------------------------|--------------|
| Mercury-Dissolved | < 2 | 2 | |

Hg Analysis Date/Time: 6/28/23/11:35hg

Hg Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Volume: 50 mL

Final Volume: 50 mL

Analytical Batch: 062823hg



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL
Project ID: JX1600
Client Project Manager: TYLER ZSCHIEDRICH
ENVision Project Number: 2023-1274
Analytical Method: EPA 8260
Prep Method: EPA 5030B
Analytical Batch: 062523VW
Client Sample ID: SB-5-GW-6-16 **Sample Collection Date/Time:** 6/21/23 13:50
Envision Sample Number: 23-12291 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: water

| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------|------------------------------|-------------------------------|--------------|
| Acetone | < 100 | 100 | |
| Acrolein | < 1 | 1 | |
| Acrylonitrile | < 0.45 | 1 | 1 |
| Benzene | < 5 | 5 | |
| Bromobenzene | < 5 | 5 | |
| Bromochloromethane | < 5 | 5 | |
| Bromodichloromethane | < 5 | 5 | |
| Bromoform | < 5 | 5 | |
| Bromomethane | < 5 | 5 | |
| n-Butanol | < 50 | 50 | |
| 2-Butanone (MEK) | < 10 | 10 | |
| n-Butylbenzene | < 5 | 5 | |
| sec-Butylbenzene | < 5 | 5 | |
| tert-Butylbenzene | < 5 | 5 | |
| Carbon Disulfide | < 5 | 5 | |
| Carbon Tetrachloride | < 5 | 5 | |
| Chlorobenzene | < 5 | 5 | |
| Chloroethane | < 5 | 5 | |
| 2-Chloroethylvinylether | < 50 | 50 | |
| Chloroform | < 5 | 5 | |
| Chloromethane | < 5 | 5 | |
| 2-Chlorotoluene | < 5 | 5 | |
| 4-Chlorotoluene | < 5 | 5 | |
| 1,2-Dibromo-3-chloropropane | < 1 | 1 | |
| Dibromochloromethane | < 5 | 5 | |
| 1,2-Dibromoethane (EDB) | < 1 | 1 | |
| Dibromomethane | < 5 | 5 | |
| 1,2-Dichlorobenzene | < 5 | 5 | |
| 1,3-Dichlorobenzene | < 5 | 5 | |
| 1,4-Dichlorobenzene | < 5 | 5 | |
| trans-1,4-Dichloro-2-butene | < 1 | 1 | |
| Dichlorodifluoromethane | < 5 | 5 | |



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| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------------|------------------------------|-------------------------------|--------------|
| 1,1-Dichloroethane | < 5 | 5 | |
| 1,2-Dichloroethane | < 5 | 5 | |
| 1,1-Dichloroethene | < 5 | 5 | |
| cis-1,2-Dichloroethene | < 5 | 5 | |
| trans-1,2-Dichloroethene | < 5 | 5 | |
| 1,2-Dichloropropane | < 5 | 5 | |
| 1,3-Dichloropropane | < 5 | 5 | |
| 2,2-Dichloropropane | < 5 | 5 | |
| 1,1-Dichloropropene | < 5 | 5 | |
| 1,3-Dichloropropene | < 4.1 | 4.1 | |
| Ethylbenzene | < 5 | 5 | |
| Ethyl methacrylate | < 100 | 100 | |
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| n-Hexane | < 10 | 10 | |
| 2-Hexanone | < 10 | 10 | |
| Iodomethane | < 10 | 10 | |
| Isopropylbenzene (Cumene) | < 5 | 5 | |
| p-Isopropyltoluene | < 5 | 5 | |
| Methylene chloride | < 5 | 5 | |
| 4-Methyl-2-pentanone (MIBK) | < 10 | 10 | |
| Methyl-tert-butyl-ether | < 5 | 5 | |
| n-Propylbenzene | < 5 | 5 | |
| Styrene | < 5 | 5 | |
| 1,1,1,2-Tetrachloroethane | < 5 | 5 | |
| 1,1,2,2-Tetrachloroethane | < 0.66 | 1 | 1 |
| Tetrachloroethene | < 5 | 5 | |
| Toluene | < 5 | 5 | |
| 1,2,3-Trichlorobenzene | < 5 | 5 | |
| 1,2,4-Trichlorobenzene | < 5 | 5 | |
| 1,1,1-Trichloroethane | < 5 | 5 | |
| 1,1,2-Trichloroethane | < 5 | 5 | |
| Trichloroethene | < 5 | 5 | |
| Trichlorofluoromethane | < 5 | 5 | |
| 1,2,3-Trichloropropane | < 1 | 1 | |
| 1,2,4-Trimethylbenzene | < 5 | 5 | |
| 1,3,5-Trimethylbenzene | < 5 | 5 | |
| Vinyl acetate | < 10 | 10 | |
| Vinyl chloride | < 2 | 2 | |
| Xylene, M&P | < 5 | 5 | |
| Xylene, Ortho | < 5 | 5 | |
| Xylene (Total) | < 10 | 10 | |
| Dibromofluoromethane (surrogate) | 93% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 95% | | |
| Toluene-d8 (surrogate) | 103% | | |
| 4-bromofluorobenzene (surrogate) | 100% | | |
| Analysis Date/Time: | 6-25-23/15:43 | | |
| Analyst Initials | tjg | | |



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL
Project ID: JX1600
Client Project Manager: TYLER ZSCHIEDRICH
ENVision Project Number: 2023-1274
Analytical Method: EPA 8270 BNA/PAH-SIM
Prep Method: EPA 3520C
BNA Analytical Batch: 062823BW

Client Sample ID: SB-5-GW-6-16 **Sample Collection Date/Time:** 6/21/23 13:50
Envision Sample Number: 23-12291 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: water

| <u>BNA Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------|------------------------------|-------------------------------|--------------|
| Aniline | < 10 | 10 | |
| Benzoic Acid | < 50 | 50 | |
| Benzyl Alcohol | < 20 | 20 | |
| 4-Bromophenylphenyl ether | < 10 | 10 | |
| Butylbenzylphthalate | < 10 | 10 | |
| Carbazole | < 20 | 20 | |
| 4-Chloro-3-methylphenol | < 20 | 20 | |
| 4-Chloroaniline | < 3.2 | 3.2 | |
| bis(2-Chloroethoxy)methane | < 10 | 10 | |
| bis(2-Chloroethyl)ether | < 0.12 | 1.0 | 1 |
| bis(2-Chloroisopropyl)ether | < 10 | 10 | |
| 2-Chloronaphthalene | < 10 | 10 | |
| 2-Chlorophenol | < 10 | 10 | |
| 4-Chlorophenylphenyl ether | < 10 | 10 | |
| Dibenzofuran | < 10 | 10 | |
| 1,2-Dichlorobenzene | < 10 | 10 | |
| 1,3-Dichlorobenzene | < 10 | 10 | |
| 1,4-Dichlorobenzene | < 10 | 10 | |
| 3,3-Dichlorobenzidine | < 1.1 | 1.1 | |
| 2,4-Dichlorophenol | < 10 | 10 | |
| Diethylphthalate | < 10 | 10 | |
| 2,4-Dimethylphenol | < 10 | 10 | |
| Dimethylphthalate | < 10 | 10 | |
| Di-n-butylphthalate | < 10 | 10 | |
| 4,6-Dinitro-2-methylphenol | < 1.2 | 1.2 | |
| 2,4-Dinitrophenol | < 30 | 50 | |
| 2,4-Dinitrotoluene | < 2.0 | 2 | |
| 2,6-Dinitrotoluene | < 10 | 10 | |
| Di-n-octylphthalate | < 10 | 10 | |
| bis(2-Ethylhexyl)phthalate | < 5 | 5 | |
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| Hexachlorobenzene | < 1.0 | 1 | |
| Hexachlorocyclopentadiene | < 25 | 25 | |
| Hexachloroethane | < 5.1 | 5.1 | |



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| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|----------------------------------|------------------------------|-------------------------------|--------------|
| Isophorone | < 10 | 10 | |
| 2-Methylphenol (o-Cresol) | < 10 | 10 | |
| 3&4-Methylphenol | < 20 | 20 | |
| 2-Nitroaniline | < 50 | 50 | |
| 3-Nitroaniline | < 50 | 50 | |
| 4-Nitroaniline | < 33 | 33 | |
| Nitrobenzene | < 1.2 | 1.2 | |
| 2-Nitrophenol | < 10 | 10 | |
| 4-Nitrophenol | < 50 | 50 | |
| N-Nitroso-di-n-propylamine | < 0.093 | 1.0 | 1 |
| N-Nitrosodiphenylamine | < 10 | 10 | |
| Pentachlorophenol | < 1.0 | 1 | |
| Phenol | < 10 | 10 | |
| 1,2,4-Trichlorobenzene | < 10 | 10 | |
| 2,4,5-Trichlorophenol | < 10 | 10 | |
| 2,4,6-Trichlorophenol | < 9 | 9 | |
| 2-Fluorophenol (surrogate) | 89% | | |
| Phenol-d6 (surrogate) | 79% | | |
| Nitrobenzene-d5 (surrogate) | 75% | | |
| 2-Fluorobiphenyl (surrogate) | 70% | | |
| 2,4,6-Tribromophenol (surrogate) | 81% | | |
| p-Terphenyl-d14 (surrogate) | 64% | | |
| Analysis Date/Time: | 06-28-23/20:58 | | |
| Analyst Initials: | gjd | | |
| Date Extracted: | 6/28/23 | | |
| Initial Sample Volume: | 1000 mL | | |
| Final Volume: | 1.0 mL | | |

PAH-SIM Analytical Batch: 062323PW2

| <u>PAH-SIM Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|--------------------------|------------------------------|-------------------------------|--------------|
| Acenaphthene | < 1.0 | 1.0 | |
| Acenaphthylene | < 1.0 | 1.0 | |
| Anthracene | < 0.10 | 0.10 | |
| Benzo(a)anthracene | < 0.10 | 0.10 | |
| Benzo(a)pyrene | < 0.10 | 0.10 | |
| Benzo(b)fluoranthene | < 0.10 | 0.10 | |
| Benzo(g,h,i)perylene | < 0.10 | 0.10 | |
| Benzo(k)fluoranthene | < 0.10 | 0.10 | |
| Chrysene | < 0.10 | 0.10 | |
| Dibenzo(a,h)anthracene | < 0.029 | 0.029 | |
| Fluoranthene | < 1.0 | 1.0 | |
| Fluorene | < 1.0 | 1.0 | |
| Indeno(1,2,3-cd)pyrene | < 0.022 | 0.022 | |
| 1-methylnaphthalene | < 1.0 | 1.0 | |
| 2-methylnaphthalene | < 1.0 | 1.0 | |
| Naphthalene | < 1.0 | 1.0 | |
| Phenanthrene | < 1.0 | 1.0 | |
| Pyrene | < 1.0 | 1.0 | |
| Analysis Date/Time: | 06-24-23/04:30 | | |
| Analyst Initials | gjd | | |



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 6010

Prep Method: EPA 3010A

Client Sample ID: SB-5-GW-6-16

Sample Collection Date/Time: 6/21/23 13:50

Envision Sample Number: 23-12291

Sample Received Date/Time: 6/22/23 9:25

Sample Matrix: water

| Compounds | Sample Results (ug/L) | Reporting Limit (ug/L) | Flags |
|--------------------|------------------------------|-------------------------------|--------------|
| Arsenic-Dissolved | < 10 | 10 | |
| Barium-Dissolved | < 100 | 100 | |
| Cadmium-Dissolved | < 5 | 5 | |
| Chromium-Dissolved | < 10 | 10 | |
| Lead-Dissolved | < 10 | 10 | |
| Selenium-Dissolved | < 10 | 10 | |
| Silver-Dissolved | < 50 | 50 | |

ICP Analysis Date/Time: 6-27-23/14:14

Analyst Initials: gjd

Date Digested: 6/27/23

Initial Sample Volume: 50 mL

Final Volume: 50 mL

Analytical Batch: 062723icp

Analytical & Prep Method: EPA 7470

| Compounds | Sample Results (ug/L) | Reporting Limit (ug/L) | Flags |
|--------------------------|------------------------------|-------------------------------|--------------|
| Mercury-Dissolved | < 2 | 2 | |
| Hg Analysis Date/Time: | 6/28/23/11:45 | | |
| Hg Analyst Initials: | gjd | | |
| Date Digested: | 6/27/2023 | | |
| Initial Sample Volume: | 50 mL | | |
| Final Volume: | 50 mL | | |
| Analytical Batch: | 062823hg | | |



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL
Project ID: JX1600
Client Project Manager: TYLER ZSCHIEDRICH
ENVision Project Number: 2023-1274
Analytical Method: EPA 8260
Prep Method: EPA 5030B
Analytical Batch: 062523VW
Client Sample ID: SB-6-GW-6-16 **Sample Collection Date/Time:** 6/21/23 15:06
Envision Sample Number: 23-12292 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: water

| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------|------------------------------|-------------------------------|--------------|
| Acetone | < 100 | 100 | |
| Acrolein | < 1 | 1 | |
| Acrylonitrile | < 0.45 | 1 | 1 |
| Benzene | < 5 | 5 | |
| Bromobenzene | < 5 | 5 | |
| Bromochloromethane | < 5 | 5 | |
| Bromodichloromethane | < 5 | 5 | |
| Bromoform | < 5 | 5 | |
| Bromomethane | < 5 | 5 | |
| n-Butanol | < 50 | 50 | |
| 2-Butanone (MEK) | < 10 | 10 | |
| n-Butylbenzene | < 5 | 5 | |
| sec-Butylbenzene | < 5 | 5 | |
| tert-Butylbenzene | < 5 | 5 | |
| Carbon Disulfide | < 5 | 5 | |
| Carbon Tetrachloride | < 5 | 5 | |
| Chlorobenzene | < 5 | 5 | |
| Chloroethane | < 5 | 5 | |
| 2-Chloroethylvinylether | < 50 | 50 | |
| Chloroform | < 5 | 5 | |
| Chloromethane | < 5 | 5 | |
| 2-Chlorotoluene | < 5 | 5 | |
| 4-Chlorotoluene | < 5 | 5 | |
| 1,2-Dibromo-3-chloropropane | < 1 | 1 | |
| Dibromochloromethane | < 5 | 5 | |
| 1,2-Dibromoethane (EDB) | < 1 | 1 | |
| Dibromomethane | < 5 | 5 | |
| 1,2-Dichlorobenzene | < 5 | 5 | |
| 1,3-Dichlorobenzene | < 5 | 5 | |
| 1,4-Dichlorobenzene | < 5 | 5 | |
| trans-1,4-Dichloro-2-butene | < 1 | 1 | |
| Dichlorodifluoromethane | < 5 | 5 | |



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| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------------|------------------------------|-------------------------------|--------------|
| 1,1-Dichloroethane | < 5 | 5 | |
| 1,2-Dichloroethane | < 5 | 5 | |
| 1,1-Dichloroethene | < 5 | 5 | |
| cis-1,2-Dichloroethene | < 5 | 5 | |
| trans-1,2-Dichloroethene | < 5 | 5 | |
| 1,2-Dichloropropane | < 5 | 5 | |
| 1,3-Dichloropropane | < 5 | 5 | |
| 2,2-Dichloropropane | < 5 | 5 | |
| 1,1-Dichloropropene | < 5 | 5 | |
| 1,3-Dichloropropene | < 4.1 | 4.1 | |
| Ethylbenzene | < 5 | 5 | |
| Ethyl methacrylate | < 100 | 100 | |
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| n-Hexane | < 10 | 10 | |
| 2-Hexanone | < 10 | 10 | |
| Iodomethane | < 10 | 10 | |
| Isopropylbenzene (Cumene) | < 5 | 5 | |
| p-Isopropyltoluene | < 5 | 5 | |
| Methylene chloride | < 5 | 5 | |
| 4-Methyl-2-pentanone (MIBK) | < 10 | 10 | |
| Methyl-tert-butyl-ether | < 5 | 5 | |
| n-Propylbenzene | < 5 | 5 | |
| Styrene | < 5 | 5 | |
| 1,1,1,2-Tetrachloroethane | < 5 | 5 | |
| 1,1,2,2-Tetrachloroethane | < 0.66 | 1 | 1 |
| Tetrachloroethene | < 5 | 5 | |
| Toluene | < 5 | 5 | |
| 1,2,3-Trichlorobenzene | < 5 | 5 | |
| 1,2,4-Trichlorobenzene | < 5 | 5 | |
| 1,1,1-Trichloroethane | < 5 | 5 | |
| 1,1,2-Trichloroethane | < 5 | 5 | |
| Trichloroethene | < 5 | 5 | |
| Trichlorofluoromethane | < 5 | 5 | |
| 1,2,3-Trichloropropane | < 1 | 1 | |
| 1,2,4-Trimethylbenzene | < 5 | 5 | |
| 1,3,5-Trimethylbenzene | < 5 | 5 | |
| Vinyl acetate | < 10 | 10 | |
| Vinyl chloride | < 2 | 2 | |
| Xylene, M&P | < 5 | 5 | |
| Xylene, Ortho | < 5 | 5 | |
| Xylene (Total) | < 10 | 10 | |
| Dibromofluoromethane (surrogate) | 93% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 97% | | |
| Toluene-d8 (surrogate) | 106% | | |
| 4-bromofluorobenzene (surrogate) | 105% | | |
| Analysis Date/Time: | 6-25-23/15:59 | | |
| Analyst Initials | tjg | | |



Analytical Report

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Client Name: AUGUST MACK ENVIRONMENTAL
Project ID: JX1600
Client Project Manager: TYLER ZSCHIEDRICH
ENVision Project Number: 2023-1274
Analytical Method: EPA 8270 BNA/PAH-SIM
Prep Method: EPA 3520C
BNA Analytical Batch: 062823BW

Client Sample ID: SB-6-GW-6-16 **Sample Collection Date/Time:** 6/21/23 15:06
Envision Sample Number: 23-12292 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: water

| <u>BNA Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|-----------------------------|------------------------------|-------------------------------|--------------|
| Aniline | < 10 | 10 | |
| Benzoic Acid | < 50 | 50 | |
| Benzyl Alcohol | < 20 | 20 | |
| 4-Bromophenylphenyl ether | < 10 | 10 | |
| Butylbenzylphthalate | < 10 | 10 | |
| Carbazole | < 20 | 20 | |
| 4-Chloro-3-methylphenol | < 20 | 20 | |
| 4-Chloroaniline | < 3.2 | 3.2 | |
| bis(2-Chloroethoxy)methane | < 10 | 10 | |
| bis(2-Chloroethyl)ether | < 0.12 | 1.0 | 1 |
| bis(2-Chloroisopropyl)ether | < 10 | 10 | |
| 2-Chloronaphthalene | < 10 | 10 | |
| 2-Chlorophenol | < 10 | 10 | |
| 4-Chlorophenylphenyl ether | < 10 | 10 | |
| Dibenzofuran | < 10 | 10 | |
| 1,2-Dichlorobenzene | < 10 | 10 | |
| 1,3-Dichlorobenzene | < 10 | 10 | |
| 1,4-Dichlorobenzene | < 10 | 10 | |
| 3,3-Dichlorobenzidine | < 1.1 | 1.1 | |
| 2,4-Dichlorophenol | < 10 | 10 | |
| Diethylphthalate | < 10 | 10 | |
| 2,4-Dimethylphenol | < 10 | 10 | |
| Dimethylphthalate | < 10 | 10 | |
| Di-n-butylphthalate | < 10 | 10 | |
| 4,6-Dinitro-2-methylphenol | < 1.2 | 1.2 | |
| 2,4-Dinitrophenol | < 30 | 50 | |
| 2,4-Dinitrotoluene | < 2.0 | 2 | |
| 2,6-Dinitrotoluene | < 10 | 10 | |
| Di-n-octylphthalate | < 10 | 10 | |
| bis(2-Ethylhexyl)phthalate | < 5 | 5 | |
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| Hexachlorobenzene | < 1.0 | 1 | |
| Hexachlorocyclopentadiene | < 25 | 25 | |
| Hexachloroethane | < 5.1 | 5.1 | |



Analytical Report

ENVision Laboratories, Inc.
 1439 Sadlier Circle West Drive
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8270 Continued...

| <u>Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|----------------------------------|------------------------------|-------------------------------|--------------|
| Isophorone | < 10 | 10 | |
| 2-Methylphenol (o-Cresol) | < 10 | 10 | |
| 3&4-Methylphenol | < 20 | 20 | |
| 2-Nitroaniline | < 50 | 50 | |
| 3-Nitroaniline | < 50 | 50 | |
| 4-Nitroaniline | < 33 | 33 | |
| Nitrobenzene | < 1.2 | 1.2 | |
| 2-Nitrophenol | < 10 | 10 | |
| 4-Nitrophenol | < 50 | 50 | |
| N-Nitroso-di-n-propylamine | < 0.093 | 1.0 | 1 |
| N-Nitrosodiphenylamine | < 10 | 10 | |
| Pentachlorophenol | < 1.0 | 1 | |
| Phenol | < 10 | 10 | |
| 1,2,4-Trichlorobenzene | < 10 | 10 | |
| 2,4,5-Trichlorophenol | < 10 | 10 | |
| 2,4,6-Trichlorophenol | < 9 | 9 | |
| 2-Fluorophenol (surrogate) | 65% | | |
| Phenol-d6 (surrogate) | 71% | | |
| Nitrobenzene-d5 (surrogate) | 71% | | |
| 2-Fluorobiphenyl (surrogate) | 65% | | |
| 2,4,6-Tribromophenol (surrogate) | 56% | | |
| p-Terphenyl-d14 (surrogate) | 51% | | |
| Analysis Date/Time: | 06-28-23/21:24 | | |
| Analyst Initials: | gjd | | |
| Date Extracted: | 6/28/23 | | |
| Initial Sample Volume: | 1000 mL | | |
| Final Volume: | 1.0 mL | | |

PAH-SIM Analytical Batch: 062323PW2

| <u>PAH-SIM Compounds</u> | <u>Sample Results (ug/L)</u> | <u>Reporting Limit (ug/L)</u> | <u>Flags</u> |
|--------------------------|------------------------------|-------------------------------|--------------|
| Acenaphthene | < 1.0 | 1.0 | |
| Acenaphthylene | < 1.0 | 1.0 | |
| Anthracene | < 0.10 | 0.10 | |
| Benzo(a)anthracene | < 0.10 | 0.10 | |
| Benzo(a)pyrene | < 0.10 | 0.10 | |
| Benzo(b)fluoranthene | < 0.10 | 0.10 | |
| Benzo(g,h,i)perylene | < 0.10 | 0.10 | |
| Benzo(k)fluoranthene | < 0.10 | 0.10 | |
| Chrysene | < 0.10 | 0.10 | |
| Dibenzo(a,h)anthracene | < 0.029 | 0.029 | |
| Fluoranthene | < 1.0 | 1.0 | |
| Fluorene | < 1.0 | 1.0 | |
| Indeno(1,2,3-cd)pyrene | < 0.022 | 0.022 | |
| 1-methylnaphthalene | < 1.0 | 1.0 | |
| 2-methylnaphthalene | < 1.0 | 1.0 | |
| Naphthalene | < 1.0 | 1.0 | |
| Phenanthrene | < 1.0 | 1.0 | |
| Pyrene | < 1.0 | 1.0 | |
| Analysis Date/Time: | 06-24-23/04:52 | | |
| Analyst Initials | gjd | | |



Analytical Report

ENVision Laboratories, Inc.
1439 Sadlier Circle West Drive
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Tel: 317.351.8632
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Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 6010

Prep Method: EPA 3010A

Client Sample ID: SB-6-GW-6-16 **Sample Collection Date/Time:** 6/21/23 15:06

Envision Sample Number: 23-12292 **Sample Received Date/Time:** 6/22/23 9:25

Sample Matrix: water

| Compounds | Sample Results (ug/L) | Reporting Limit (ug/L) | Flags |
|--------------------|------------------------------|-------------------------------|--------------|
| Arsenic-Dissolved | < 10 | 10 | |
| Barium-Dissolved | < 100 | 100 | |
| Cadmium-Dissolved | < 5 | 5 | |
| Chromium-Dissolved | < 10 | 10 | |
| Lead-Dissolved | < 10 | 10 | |
| Selenium-Dissolved | < 10 | 10 | |
| Silver-Dissolved | < 50 | 50 | |

ICP Analysis Date/Time: 6-27-23/14:17

Analyst Initials: gjd

Date Digested: 6/27/23

Initial Sample Volume: 50 mL

Final Volume: 50 mL

Analytical Batch: 062723icp

Analytical & Prep Method: EPA 7470

| Compounds | Sample Results (ug/L) | Reporting Limit (ug/L) | Flags |
|-------------------|------------------------------|-------------------------------|--------------|
| Mercury-Dissolved | < 2 | 2 | |

Hg Analysis Date/Time: 6/28/23/11:47

Hg Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Volume: 50 mL

Final Volume: 50 mL

Analytical Batch: 062823hg



Analytical Report

ENVision Laboratories, Inc.
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Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 8260

Prep Method: EPA 5035A

Analytical Batch: 062323VS

Client Sample ID: SB-7-1-1.5

Sample Collection Date/Time: 6/21/23

15:25

Envision Sample Number: 23-12293

Sample Received Date/Time: 6/22/23

9:25

Sample Matrix: soil

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------|------------------------|--------------------|-------|
| Acetone | < 0.122 | 0.122 | |
| Acrolein | < 0.00021 | 0.001 | 1 |
| Acrylonitrile | < 0.002 | 0.002 | |
| Benzene | < 0.006 | 0.006 | |
| Bromobenzene | < 0.006 | 0.006 | |
| Bromochloromethane | < 0.006 | 0.006 | |
| Bromodichloromethane | < 0.006 | 0.006 | |
| Bromoform | < 0.006 | 0.006 | |
| Bromomethane | < 0.006 | 0.006 | |
| n-Butanol | < 0.061 | 0.061 | |
| 2-Butanone (MEK) | < 0.012 | 0.012 | |
| n-Butylbenzene | 0.198 | 0.006 | |
| sec-Butylbenzene | 0.161 | 0.006 | |
| tert-Butylbenzene | < 0.006 | 0.006 | |
| Carbon Disulfide | < 0.006 | 0.006 | |
| Carbon Tetrachloride | < 0.006 | 0.006 | |
| Chlorobenzene | < 0.006 | 0.006 | |
| Chloroethane | < 0.006 | 0.006 | |
| 2-Chloroethylvinylether | < 0.061 | 0.061 | |
| Chloroform | < 0.006 | 0.006 | |
| Chloromethane | < 0.006 | 0.006 | |
| 2-Chlorotoluene | < 0.006 | 0.006 | |
| 4-Chlorotoluene | < 0.006 | 0.006 | |
| 1,2-Dibromo-3-chloropropane | < 0.0021 | 0.0021 | |
| Dibromochloromethane | < 0.006 | 0.006 | |
| 1,2-Dibromoethane (EDB) | < 0.00034 | 0.001 | 1 |
| Dibromomethane | < 0.006 | 0.006 | |
| 1,2-Dichlorobenzene | < 0.006 | 0.006 | |
| 1,3-Dichlorobenzene | < 0.006 | 0.006 | |
| 1,4-Dichlorobenzene | < 0.006 | 0.006 | |
| trans-1,4-Dichloro-2-butene | < 0.006 | 0.006 | |
| Dichlorodifluoromethane | < 0.006 | 0.006 | |
| 1,1-Dichloroethane | 0.0111 | 0.006 | |
| 1,2-Dichloroethane | < 0.006 | 0.006 | |
| 1,1-Dichloroethene | < 0.006 | 0.006 | |

**8260 continued...**

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------------|------------------------|--------------------|-------|
| cis-1,2-Dichloroethene | 0.357 | 0.305 | 2 |
| trans-1,2-Dichloroethene | < 0.006 | 0.006 | |
| 1,2-Dichloropropane | < 0.006 | 0.006 | |
| 1,3-Dichloropropane | < 0.006 | 0.006 | |
| 2,2-Dichloropropane | < 0.006 | 0.006 | |
| 1,1-Dichloropropene | < 0.006 | 0.006 | |
| 1,3-Dichloropropene | < 0.006 | 0.006 | |
| Ethylbenzene | 0.0285 | 0.006 | |
| Ethyl methacrylate | < 0.122 | 0.122 | |
| Hexachloro-1,3-butadiene | < 0.006 | 0.006 | |
| n-Hexane | < 0.012 | 0.012 | |
| 2-Hexanone | < 0.012 | 0.012 | |
| Iodomethane | < 0.012 | 0.012 | |
| Isopropylbenzene (Cumene) | 0.0482 | 0.006 | |
| p-Isopropyltoluene | 0.168 | 0.006 | |
| Methylene chloride | < 0.024 | 0.024 | |
| 4-Methyl-2-pentanone (MIBK) | < 0.012 | 0.012 | |
| Methyl-tert-butyl-ether | < 0.006 | 0.006 | |
| n-Propylbenzene | 0.132 | 0.006 | |
| Styrene | < 0.006 | 0.006 | |
| 1,1,1,2-Tetrachloroethane | < 0.006 | 0.006 | |
| 1,1,2,2-Tetrachloroethane | < 0.006 | 0.006 | |
| Tetrachloroethene | 0.817 | 0.305 | 2 |
| Toluene | < 0.006 | 0.006 | |
| 1,2,3-Trichlorobenzene | < 0.006 | 0.006 | |
| 1,2,4-Trichlorobenzene | < 0.006 | 0.006 | |
| 1,1,1-Trichloroethane | < 0.006 | 0.006 | |
| 1,1,2-Trichloroethane | < 0.006 | 0.006 | |
| Trichloroethene | 1.50 | 0.305 | 2 |
| Trichlorofluoromethane | < 0.006 | 0.006 | |
| 1,2,3-Trichloropropane | < 0.006 | 0.006 | |
| 1,2,4-Trimethylbenzene | 1.67 | 0.305 | 2 |
| 1,3,5-Trimethylbenzene | 0.580 | 0.305 | 2 |
| Vinyl acetate | < 0.012 | 0.012 | |
| Vinyl chloride | < 0.002 | 0.002 | |
| Xylene, M&P | 0.0815 | 0.006 | |
| Xylene, Ortho | 0.116 | 0.006 | |
| Xylene, Total | 0.197 | 0.012 | |
| Dibromofluoromethane (surrogate) | 111% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 90% | | |
| Toluene-d8 (surrogate) | 104% | | |
| 4-bromofluorobenzene (surrogate) | 112% | | |
| Analysis Date: | 6/23/23 | | |
| Analysis Time: | 21:33 | | |
| Analyst Initials | tjg | | |

Percent Solids: 82%

All results reported on dry weight basis.



Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 8270 PAH

Prep Method: EPA 3550C

Analytical Batch: 062623BS

Client Sample ID: SB-7-1-1.5 **Sample Collection Date/Time:** 6/21/23 15:25
Envision Sample Number: 23-12293 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: soil

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------|------------------------|--------------------|-------|
| Acenaphthene | < 0.41 | 0.41 | |
| Acenaphthylene | < 0.41 | 0.41 | |
| Aniline | < 0.41 | 0.41 | |
| Anthracene | < 0.41 | 0.41 | |
| Benzo(a)anthracene | < 0.41 | 0.41 | |
| Benzo(a)pyrene | < 0.081 | 0.081 | |
| Benzo(b)fluoranthene | < 0.41 | 0.41 | |
| Benzo(g,h,i)perylene | < 0.41 | 0.41 | |
| Benzo(k)fluoranthene | < 0.41 | 0.41 | |
| Benzoic Acid | < 2.03 | 2.03 | |
| Benzyl Alcohol | < 0.81 | 0.81 | |
| 4-Bromophenylphenyl ether | < 0.41 | 0.41 | |
| Butylbenzylphthalate | < 0.41 | 0.41 | |
| Carbazole | < 0.81 | 0.81 | |
| 4-Chloro-3-methylphenol | < 0.81 | 0.81 | |
| 4-Chloroaniline | < 0.033 | 0.040 | 1 |
| bis(2-Chloroethoxy)methane | < 0.081 | 0.081 | |
| bis(2-Chloroethyl)ether | < 0.081 | 0.081 | |
| bis(2-Chloroisopropyl)ether | < 0.41 | 0.41 | |
| 2-Chloronaphthalene | < 0.41 | 0.41 | |
| 2-Chlorophenol | < 0.41 | 0.41 | |
| 4-Chlorophenylphenyl ether | < 0.41 | 0.41 | |
| Chrysene | < 0.41 | 0.41 | |
| Dibenzo(a,h)anthracene | < 0.081 | 0.081 | |
| Dibenzofuran | < 0.41 | 0.41 | |
| 1,2-Dichlorobenzene | < 0.41 | 0.41 | |
| 1,3-Dichlorobenzene | < 0.41 | 0.41 | |
| 1,4-Dichlorobenzene | < 0.41 | 0.41 | |
| 3,3-Dichlorobenzidine | < 0.17 | 0.81 | |
| 2,4-Dichlorophenol | < 0.41 | 0.41 | |
| Diethylphthalate | < 0.41 | 0.41 | |
| 2,4-Dimethylphenol | < 0.41 | 0.41 | |
| Dimethylphthalate | < 0.41 | 0.41 | |
| Di-n-butylphthalate | < 0.41 | 0.41 | |

**8270 continued...**

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|----------------------------------|------------------------|--------------------|-------|
| 4,6-Dinitro-2-methylphenol | < 0.050 | 0.050 | |
| 2,4-Dinitrophenol | < 0.081 | 0.081 | |
| 2,4-Dinitrotoluene | < 0.066 | 0.066 | |
| 2,6-Dinitrotoluene | < 0.41 | 0.41 | |
| Di-n-octylphthalate | < 0.41 | 0.41 | |
| bis(2-Ethylhexyl)phthalate | < 0.41 | 0.41 | |
| Fluoranthene | < 0.41 | 0.41 | |
| Fluorene | < 0.41 | 0.41 | |
| Hexachloro-1,3-butadiene | < 0.081 | 0.081 | |
| Hexachlorobenzene | < 0.081 | 0.081 | |
| Hexachlorocyclopentadiene | < 0.41 | 0.41 | |
| Hexachloroethane | < 0.081 | 0.081 | |
| Indeno(1,2,3-cd)pyrene | < 0.41 | 0.41 | |
| Isophorone | < 0.41 | 0.41 | |
| 2-Methylphenol (o-Cresol) | < 0.41 | 0.41 | |
| 3&4-Methylphenol | < 0.81 | 0.81 | |
| 1-Methylnaphthalene | 0.519 | 0.41 | |
| 2-Methylnaphthalene | 0.684 | 0.41 | |
| Naphthalene | 1.00 | 0.081 | |
| 2-Nitroaniline | < 1.63 | 1.63 | |
| 3-Nitroaniline | < 2.03 | 2.03 | |
| 4-Nitroaniline | < 0.081 | 0.081 | |
| Nitrobenzene | < 0.04 | 0.04 | |
| 2-Nitrophenol | < 0.41 | 0.41 | |
| 4-Nitrophenol | < 2.03 | 2.03 | |
| N-Nitroso-di-n-propylamine | < 0.081 | 0.081 | |
| N-Nitrosodiphenylamine | < 0.41 | 0.41 | |
| Pentachlorophenol | < 0.081 | 0.081 | |
| Phenanthrene | < 0.41 | 0.41 | |
| Phenol | < 0.41 | 0.41 | |
| Pyrene | < 0.41 | 0.41 | |
| 1,2,4-Trichlorobenzene | < 0.41 | 0.41 | |
| 2,4,5-Trichlorophenol | < 0.41 | 0.41 | |
| 2,4,6-Trichlorophenol | < 0.41 | 0.41 | |
| 2-Fluorophenol (surrogate) | 35% | | |
| Phenol-d6 (surrogate) | 29% | | |
| Nitrobenzene-d5 (surrogate) | 55% | | |
| 2-Fluorobiphenyl (surrogate) | 50% | | |
| 2,4,6-Tribromophenol (surrogate) | 45% | | |
| p-Terphenyl-d14 (surrogate) | 46% | | |
| Analysis Date: | 6/27/2023 | | |
| Analysis Time: | 1:44 | | |
| Analyst Initials: | JAK | | |
| Date Extracted: | 6/26/23 | | |
| Initial Sample Weight (g): | 30 | | |
| Final Volume (mL): | 1 | | |
| Percent Solids | 82% | | |



Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 6010B

Prep Method: EPA 3050B

Client Sample ID: SB-7-1-1.5 **Sample Collection Date/Time:** 6/21/23 15:25

Envision Sample Number: 23-12293 **Sample Received Date/Time:** 6/22/23 9:25

Sample Matrix: soil

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|------------------|-------------------------------|---------------------------|--------------|
| Arsenic | < 2 | 2 | |
| Barium | 58 | 2 | |
| Cadmium | < 2 | 2 | |
| Chromium | 17 | 2 | |
| Lead | 95 | 2 | |
| Selenium | < 2 | 2 | |
| Silver | < 2 | 2 | |

Analysis Date:

Analysis Time: 6-27-23/14:19

Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Weight: 1.0 g

Final Volume: 50 mL

Analytical Batch: 062723icp

Analytical Method: EPA 7471A

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|------------------|-------------------------------|---------------------------|--------------|
| Mercury | < 1 | 1 | |

Hg Analysis Date: 6/28/2023

Hg Analysis Time: 11:48

Hg Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Weight: 0.6 g

Final Volume: 50 mL

Analytical Batch: 062823hg

Percent Solids 82%

All results reported on dry weight basis.



Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

| | | | | |
|--------------------------------|------------|-------------------------------------|---------|-------|
| Client Sample ID: | SB-7-1-1.5 | Sample Collection Date/Time: | 6/21/23 | 15:25 |
| Envision Sample Number: | 23-12293 | Sample Received Date/Time: | 6/22/23 | 9:25 |
| Sample Matrix: | soil | | | |

| Analyte | Sample Results | Flags | Method |
|------------------|-----------------------|--------------|---------------|
| Percent Moisture | 18.0% | | EPA 1684 |
| Percent Solids | 82.0% | | EPA 1684 |
| Analysis Date: | 6/23/23 | | |
| Analyst Initials | NR | | |



Analytical Report

ENVision Laboratories, Inc.
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Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 8260

Prep Method: EPA 5035A

Analytical Batch: 062323VS

| | | | | |
|--------------------------------|----------|-------------------------------------|---------|-------|
| Client Sample ID: | SB-1-2-4 | Sample Collection Date/Time: | 6/21/23 | 15:55 |
| Envision Sample Number: | 23-12294 | Sample Received Date/Time: | 6/22/23 | 9:25 |
| Sample Matrix: | soil | | | |

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|------------------|-------------------------------|---------------------------|--------------|
|------------------|-------------------------------|---------------------------|--------------|

| | | | |
|-----------------------------|-----------|--------|---|
| Acetone | < 0.118 | 0.118 | |
| Acrolein | < 0.00020 | 0.001 | 1 |
| Acrylonitrile | < 0.002 | 0.002 | |
| Benzene | < 0.006 | 0.006 | |
| Bromobenzene | < 0.006 | 0.006 | |
| Bromochloromethane | < 0.006 | 0.006 | |
| Bromodichloromethane | < 0.006 | 0.006 | |
| Bromoform | < 0.006 | 0.006 | |
| Bromomethane | < 0.006 | 0.006 | |
| n-Butanol | < 0.059 | 0.059 | |
| 2-Butanone (MEK) | < 0.012 | 0.012 | |
| n-Butylbenzene | < 0.006 | 0.006 | |
| sec-Butylbenzene | < 0.006 | 0.006 | |
| tert-Butylbenzene | < 0.006 | 0.006 | |
| Carbon Disulfide | < 0.006 | 0.006 | |
| Carbon Tetrachloride | < 0.006 | 0.006 | |
| Chlorobenzene | < 0.006 | 0.006 | |
| Chloroethane | < 0.006 | 0.006 | |
| 2-Chloroethylvinylether | < 0.059 | 0.059 | |
| Chloroform | < 0.006 | 0.006 | |
| Chloromethane | < 0.006 | 0.006 | |
| 2-Chlorotoluene | < 0.006 | 0.006 | |
| 4-Chlorotoluene | < 0.006 | 0.006 | |
| 1,2-Dibromo-3-chloropropane | < 0.0020 | 0.0020 | |
| Dibromochloromethane | < 0.006 | 0.006 | |
| 1,2-Dibromoethane (EDB) | < 0.00033 | 0.001 | 1 |
| Dibromomethane | < 0.006 | 0.006 | |
| 1,2-Dichlorobenzene | < 0.006 | 0.006 | |
| 1,3-Dichlorobenzene | < 0.006 | 0.006 | |
| 1,4-Dichlorobenzene | < 0.006 | 0.006 | |
| trans-1,4-Dichloro-2-butene | < 0.006 | 0.006 | |
| Dichlorodifluoromethane | < 0.006 | 0.006 | |
| 1,1-Dichloroethane | < 0.006 | 0.006 | |
| 1,2-Dichloroethane | < 0.006 | 0.006 | |
| 1,1-Dichloroethene | < 0.006 | 0.006 | |

**8260 continued...**

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------------|------------------------|--------------------|-------|
| cis-1,2-Dichloroethene | < 0.006 | 0.006 | |
| trans-1,2-Dichloroethene | < 0.006 | 0.006 | |
| 1,2-Dichloropropane | < 0.006 | 0.006 | |
| 1,3-Dichloropropane | < 0.006 | 0.006 | |
| 2,2-Dichloropropane | < 0.006 | 0.006 | |
| 1,1-Dichloropropene | < 0.006 | 0.006 | |
| 1,3-Dichloropropene | < 0.006 | 0.006 | |
| Ethylbenzene | < 0.006 | 0.006 | |
| Ethyl methacrylate | < 0.118 | 0.118 | |
| Hexachloro-1,3-butadiene | < 0.006 | 0.006 | |
| n-Hexane | < 0.012 | 0.012 | |
| 2-Hexanone | < 0.012 | 0.012 | |
| Iodomethane | < 0.012 | 0.012 | |
| Isopropylbenzene (Cumene) | < 0.006 | 0.006 | |
| p-Isopropyltoluene | < 0.006 | 0.006 | |
| Methylene chloride | < 0.024 | 0.024 | |
| 4-Methyl-2-pentanone (MIBK) | < 0.012 | 0.012 | |
| Methyl-tert-butyl-ether | < 0.006 | 0.006 | |
| n-Propylbenzene | < 0.006 | 0.006 | |
| Styrene | < 0.006 | 0.006 | |
| 1,1,1,2-Tetrachloroethane | < 0.006 | 0.006 | |
| 1,1,2,2-Tetrachloroethane | < 0.006 | 0.006 | |
| Tetrachloroethene | 0.00599 | 0.006 | |
| Toluene | < 0.006 | 0.006 | |
| 1,2,3-Trichlorobenzene | < 0.006 | 0.006 | |
| 1,2,4-Trichlorobenzene | < 0.006 | 0.006 | |
| 1,1,1-Trichloroethane | < 0.006 | 0.006 | |
| 1,1,2-Trichloroethane | < 0.006 | 0.006 | |
| Trichloroethene | < 0.006 | 0.006 | |
| Trichlorofluoromethane | < 0.006 | 0.006 | |
| 1,2,3-Trichloropropane | < 0.006 | 0.006 | |
| 1,2,4-Trimethylbenzene | < 0.006 | 0.006 | |
| 1,3,5-Trimethylbenzene | < 0.006 | 0.006 | |
| Vinyl acetate | < 0.012 | 0.012 | |
| Vinyl chloride | < 0.002 | 0.002 | |
| Xylene, M&P | < 0.006 | 0.006 | |
| Xylene, Ortho | < 0.006 | 0.006 | |
| Xylene, Total | < 0.012 | 0.012 | |
| Dibromofluoromethane (surrogate) | 100% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 87% | | |
| Toluene-d8 (surrogate) | 94% | | |
| 4-bromofluorobenzene (surrogate) | 86% | | |
| Analysis Date: | 6/23/23 | | |
| Analysis Time: | 21:49 | | |
| Analyst Initials | tjg | | |
| Percent Solids: | 85% | | |

All results reported on dry weight basis.



Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 8270 PAH

Prep Method: EPA 3550C

Analytical Batch: 062623BS

Client Sample ID: SB-1-2-4 **Sample Collection Date/Time:** 6/21/23 15:55
Envision Sample Number: 23-12294 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: soil

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------|------------------------|--------------------|-------|
| Acenaphthene | < 0.39 | 0.39 | |
| Acenaphthylene | < 0.39 | 0.39 | |
| Aniline | < 0.39 | 0.39 | |
| Anthracene | < 0.39 | 0.39 | |
| Benzo(a)anthracene | 0.988 | 0.39 | |
| Benzo(a)pyrene | 1.01 | 0.078 | |
| Benzo(b)fluoranthene | 1.06 | 0.39 | |
| Benzo(g,h,i)perylene | 0.947 | 0.39 | |
| Benzo(k)fluoranthene | 0.411 | 0.39 | |
| Benzoic Acid | < 1.96 | 1.96 | |
| Benzyl Alcohol | < 0.78 | 0.78 | |
| 4-Bromophenylphenyl ether | < 0.39 | 0.39 | |
| Butylbenzylphthalate | < 0.39 | 0.39 | |
| Carbazole | < 0.78 | 0.78 | |
| 4-Chloro-3-methylphenol | < 0.78 | 0.78 | |
| 4-Chloroaniline | < 0.032 | 0.039 | 1 |
| bis(2-Chloroethoxy)methane | < 0.078 | 0.078 | |
| bis(2-Chloroethyl)ether | < 0.078 | 0.078 | |
| bis(2-Chloroisopropyl)ether | < 0.39 | 0.39 | |
| 2-Chloronaphthalene | < 0.39 | 0.39 | |
| 2-Chlorophenol | < 0.39 | 0.39 | |
| 4-Chlorophenylphenyl ether | < 0.39 | 0.39 | |
| Chrysene | 1.03 | 0.39 | |
| Dibenzo(a,h)anthracene | < 0.078 | 0.078 | |
| Dibenzofuran | < 0.39 | 0.39 | |
| 1,2-Dichlorobenzene | < 0.39 | 0.39 | |
| 1,3-Dichlorobenzene | < 0.39 | 0.39 | |
| 1,4-Dichlorobenzene | < 0.39 | 0.39 | |
| 3,3-Dichlorobenzidine | < 0.16 | 0.78 | |
| 2,4-Dichlorophenol | < 0.39 | 0.39 | |
| Diethylphthalate | < 0.39 | 0.39 | |
| 2,4-Dimethylphenol | < 0.39 | 0.39 | |
| Dimethylphthalate | < 0.39 | 0.39 | |
| Di-n-butylphthalate | < 0.39 | 0.39 | |

**8270 continued...**

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|----------------------------------|------------------------|--------------------|-------|
| 4,6-Dinitro-2-methylphenol | < 0.048 | 0.048 | |
| 2,4-Dinitrophenol | < 0.078 | 0.078 | |
| 2,4-Dinitrotoluene | < 0.064 | 0.064 | |
| 2,6-Dinitrotoluene | < 0.39 | 0.39 | |
| Di-n-octylphthalate | < 0.39 | 0.39 | |
| bis(2-Ethylhexyl)phthalate | < 0.39 | 0.39 | |
| Fluoranthene | 1.80 | 0.39 | |
| Fluorene | < 0.39 | 0.39 | |
| Hexachloro-1,3-butadiene | < 0.078 | 0.078 | |
| Hexachlorobenzene | < 0.078 | 0.078 | |
| Hexachlorocyclopentadiene | < 0.39 | 0.39 | |
| Hexachloroethane | < 0.078 | 0.078 | |
| Indeno(1,2,3-cd)pyrene | 0.885 | 0.39 | |
| Isophorone | < 0.39 | 0.39 | |
| 2-Methylphenol (o-Cresol) | < 0.39 | 0.39 | |
| 3&4-Methylphenol | < 0.78 | 0.78 | |
| 1-Methylnaphthalene | < 0.39 | 0.39 | |
| 2-Methylnaphthalene | < 0.39 | 0.39 | |
| Naphthalene | < 0.078 | 0.078 | |
| 2-Nitroaniline | < 1.57 | 1.57 | |
| 3-Nitroaniline | < 1.96 | 1.96 | |
| 4-Nitroaniline | < 0.078 | 0.078 | |
| Nitrobenzene | < 0.04 | 0.04 | |
| 2-Nitrophenol | < 0.39 | 0.39 | |
| 4-Nitrophenol | < 1.96 | 1.96 | |
| N-Nitroso-di-n-propylamine | < 0.078 | 0.078 | |
| N-Nitrosodiphenylamine | < 0.39 | 0.39 | |
| Pentachlorophenol | < 0.078 | 0.078 | |
| Phenanthrene | 0.791 | 0.39 | |
| Phenol | < 0.39 | 0.39 | |
| Pyrene | 1.54 | 0.39 | |
| 1,2,4-Trichlorobenzene | < 0.39 | 0.39 | |
| 2,4,5-Trichlorophenol | < 0.39 | 0.39 | |
| 2,4,6-Trichlorophenol | < 0.39 | 0.39 | |
| 2-Fluorophenol (surrogate) | 35% | | |
| Phenol-d6 (surrogate) | 31% | | |
| Nitrobenzene-d5 (surrogate) | 26% | | |
| 2-Fluorobiphenyl (surrogate) | 31% | | |
| 2,4,6-Tribromophenol (surrogate) | 40% | | |
| p-Terphenyl-d14 (surrogate) | 27% | | |
| Analysis Date: | 6/27/2023 | | |
| Analysis Time: | 2:10 | | |
| Analyst Initials: | JAK | | |
| Date Extracted: | 6/26/23 | | |
| Initial Sample Weight (g): | 30 | | |
| Final Volume (mL): | 1 | | |
| Percent Solids | 85% | | |

All results reported on dry weight basis.



Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 6010B

Prep Method: EPA 3050B

| | | | | |
|--------------------------------|----------|-------------------------------------|---------|-------|
| Client Sample ID: | SB-1-2-4 | Sample Collection Date/Time: | 6/21/23 | 15:55 |
| Envision Sample Number: | 23-12294 | Sample Received Date/Time: | 6/22/23 | 9:25 |
| Sample Matrix: | soil | | | |

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|------------------|-------------------------------|---------------------------|--------------|
| Arsenic | < 2 | 2 | |
| Barium | 31 | 2 | |
| Cadmium | < 2 | 2 | |
| Chromium | 8.8 | 2 | |
| Lead | 11 | 2 | |
| Selenium | < 2 | 2 | |
| Silver | < 2 | 2 | |

Analysis Date:

Analysis Time: 6-27-23/14:21

Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Weight: 1.0 g

Final Volume: 50 mL

Analytical Batch: 062723icp

Analytical Method: EPA 7471A

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|------------------|-------------------------------|---------------------------|--------------|
| Mercury | < 1 | 1 | |

Hg Analysis Date: 6/28/2023

Hg Analysis Time: 11:50

Hg Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Weight: 0.6 g

Final Volume: 50 mL

Analytical Batch: 062823hg

Percent Solids 85%

All results reported on dry weight basis.



Analytical Report

ENVision Laboratories, Inc.
1439 Sadlier Circle West Drive
Indianapolis, IN 46239
Tel: 317.351.8632
Fax: 317.351.8639
www.envisionlaboratories.com

Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

| | | | | |
|--------------------------------|----------|-------------------------------------|---------|-------|
| Client Sample ID: | SB-1-2-4 | Sample Collection Date/Time: | 6/21/23 | 15:55 |
| Envision Sample Number: | 23-12294 | Sample Received Date/Time: | 6/22/23 | 9:25 |
| Sample Matrix: | soil | | | |

| Analyte | Sample Results | Flags | Method |
|------------------|-----------------------|--------------|---------------|
| Percent Moisture | 15.0% | | EPA 1684 |
| Percent Solids | 85.0% | | EPA 1684 |
| Analysis Date: | 6/23/23 | | |
| Analyst Initials | NR | | |



Analytical Report

ENVision Laboratories, Inc.
1439 Sadlier Circle West Drive
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Tel: 317.351.8632
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Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 8260

Prep Method: EPA 5035A

Analytical Batch: 062323VS

| | | | | |
|--------------------------------|----------|-------------------------------------|---------|-------|
| Client Sample ID: | SB-2-2-4 | Sample Collection Date/Time: | 6/21/23 | 17:05 |
| Envision Sample Number: | 23-12295 | Sample Received Date/Time: | 6/22/23 | 9:25 |
| Sample Matrix: | soil | | | |

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|------------------|-------------------------------|---------------------------|--------------|
|------------------|-------------------------------|---------------------------|--------------|

| | | | |
|-----------------------------|-----------|--------|---|
| Acetone | < 0.104 | 0.104 | |
| Acrolein | < 0.00018 | 0.001 | 1 |
| Acrylonitrile | < 0.002 | 0.002 | |
| Benzene | < 0.005 | 0.005 | |
| Bromobenzene | < 0.005 | 0.005 | |
| Bromochloromethane | < 0.005 | 0.005 | |
| Bromodichloromethane | < 0.005 | 0.005 | |
| Bromoform | < 0.005 | 0.005 | |
| Bromomethane | < 0.005 | 0.005 | |
| n-Butanol | < 0.052 | 0.052 | |
| 2-Butanone (MEK) | < 0.010 | 0.010 | |
| n-Butylbenzene | < 0.005 | 0.005 | |
| sec-Butylbenzene | < 0.005 | 0.005 | |
| tert-Butylbenzene | < 0.005 | 0.005 | |
| Carbon Disulfide | < 0.005 | 0.005 | |
| Carbon Tetrachloride | < 0.005 | 0.005 | |
| Chlorobenzene | < 0.005 | 0.005 | |
| Chloroethane | < 0.005 | 0.005 | |
| 2-Chloroethylvinylether | < 0.052 | 0.052 | |
| Chloroform | < 0.005 | 0.005 | |
| Chloromethane | < 0.005 | 0.005 | |
| 2-Chlorotoluene | < 0.005 | 0.005 | |
| 4-Chlorotoluene | < 0.005 | 0.005 | |
| 1,2-Dibromo-3-chloropropane | < 0.0018 | 0.0018 | |
| Dibromochloromethane | < 0.005 | 0.005 | |
| 1,2-Dibromoethane (EDB) | < 0.00029 | 0.001 | 1 |
| Dibromomethane | < 0.005 | 0.005 | |
| 1,2-Dichlorobenzene | < 0.005 | 0.005 | |
| 1,3-Dichlorobenzene | < 0.005 | 0.005 | |
| 1,4-Dichlorobenzene | < 0.005 | 0.005 | |
| trans-1,4-Dichloro-2-butene | < 0.005 | 0.005 | |
| Dichlorodifluoromethane | < 0.005 | 0.005 | |
| 1,1-Dichloroethane | < 0.005 | 0.005 | |
| 1,2-Dichloroethane | < 0.005 | 0.005 | |
| 1,1-Dichloroethene | < 0.005 | 0.005 | |

**8260 continued...**

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------------|------------------------|--------------------|-------|
| cis-1,2-Dichloroethene | < 0.005 | 0.005 | |
| trans-1,2-Dichloroethene | < 0.005 | 0.005 | |
| 1,2-Dichloropropane | < 0.005 | 0.005 | |
| 1,3-Dichloropropane | < 0.005 | 0.005 | |
| 2,2-Dichloropropane | < 0.005 | 0.005 | |
| 1,1-Dichloropropene | < 0.005 | 0.005 | |
| 1,3-Dichloropropene | < 0.005 | 0.005 | |
| Ethylbenzene | < 0.005 | 0.005 | |
| Ethyl methacrylate | < 0.104 | 0.104 | |
| Hexachloro-1,3-butadiene | < 0.005 | 0.005 | |
| n-Hexane | < 0.010 | 0.010 | |
| 2-Hexanone | < 0.010 | 0.010 | |
| Iodomethane | < 0.010 | 0.010 | |
| Isopropylbenzene (Cumene) | < 0.005 | 0.005 | |
| p-Isopropyltoluene | < 0.005 | 0.005 | |
| Methylene chloride | < 0.021 | 0.021 | |
| 4-Methyl-2-pentanone (MIBK) | < 0.010 | 0.010 | |
| Methyl-tert-butyl-ether | < 0.005 | 0.005 | |
| n-Propylbenzene | < 0.005 | 0.005 | |
| Styrene | < 0.005 | 0.005 | |
| 1,1,1,2-Tetrachloroethane | < 0.005 | 0.005 | |
| 1,1,2,2-Tetrachloroethane | < 0.005 | 0.005 | |
| Tetrachloroethene | 0.0188 | 0.005 | |
| Toluene | < 0.005 | 0.005 | |
| 1,2,3-Trichlorobenzene | < 0.005 | 0.005 | |
| 1,2,4-Trichlorobenzene | < 0.005 | 0.005 | |
| 1,1,1-Trichloroethane | < 0.005 | 0.005 | |
| 1,1,2-Trichloroethane | < 0.005 | 0.005 | |
| Trichloroethene | < 0.005 | 0.005 | |
| Trichlorofluoromethane | < 0.005 | 0.005 | |
| 1,2,3-Trichloropropane | < 0.005 | 0.005 | |
| 1,2,4-Trimethylbenzene | < 0.005 | 0.005 | |
| 1,3,5-Trimethylbenzene | < 0.005 | 0.005 | |
| Vinyl acetate | < 0.010 | 0.010 | |
| Vinyl chloride | < 0.002 | 0.002 | |
| Xylene, M&P | < 0.005 | 0.005 | |
| Xylene, Ortho | < 0.005 | 0.005 | |
| Xylene, Total | < 0.010 | 0.010 | |
| Dibromofluoromethane (surrogate) | 99% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 101% | | |
| Toluene-d8 (surrogate) | 92% | | |
| 4-bromofluorobenzene (surrogate) | 109% | | |
| Analysis Date: | 6/23/23 | | |
| Analysis Time: | 22:05 | | |
| Analyst Initials | tjg | | |
| Percent Solids: | 96% | | |

All results reported on dry weight basis.



Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 8270 PAH

Prep Method: EPA 3550C

Analytical Batch: 062623BS

Client Sample ID: SB-2-2-4 **Sample Collection Date/Time:** 6/21/23 17:05
Envision Sample Number: 23-12295 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: soil

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------|------------------------|--------------------|-------|
| Acenaphthene | < 0.35 | 0.35 | |
| Acenaphthylene | < 0.35 | 0.35 | |
| Aniline | < 0.35 | 0.35 | |
| Anthracene | < 0.35 | 0.35 | |
| Benzo(a)anthracene | < 0.35 | 0.35 | |
| Benzo(a)pyrene | < 0.069 | 0.069 | |
| Benzo(b)fluoranthene | < 0.35 | 0.35 | |
| Benzo(g,h,i)perylene | < 0.35 | 0.35 | |
| Benzo(k)fluoranthene | < 0.35 | 0.35 | |
| Benzoic Acid | < 1.74 | 1.74 | |
| Benzyl Alcohol | < 0.69 | 0.69 | |
| 4-Bromophenylphenyl ether | < 0.35 | 0.35 | |
| Butylbenzylphthalate | < 0.35 | 0.35 | |
| Carbazole | < 0.69 | 0.69 | |
| 4-Chloro-3-methylphenol | < 0.69 | 0.69 | |
| 4-Chloroaniline | < 0.028 | 0.034 | 1 |
| bis(2-Chloroethoxy)methane | < 0.069 | 0.069 | |
| bis(2-Chloroethyl)ether | < 0.069 | 0.069 | |
| bis(2-Chloroisopropyl)ether | < 0.35 | 0.35 | |
| 2-Chloronaphthalene | < 0.35 | 0.35 | |
| 2-Chlorophenol | < 0.35 | 0.35 | |
| 4-Chlorophenylphenyl ether | < 0.35 | 0.35 | |
| Chrysene | < 0.35 | 0.35 | |
| Dibenzo(a,h)anthracene | < 0.069 | 0.069 | |
| Dibenzofuran | < 0.35 | 0.35 | |
| 1,2-Dichlorobenzene | < 0.35 | 0.35 | |
| 1,3-Dichlorobenzene | < 0.35 | 0.35 | |
| 1,4-Dichlorobenzene | < 0.35 | 0.35 | |
| 3,3-Dichlorobenzidine | < 0.15 | 0.69 | |
| 2,4-Dichlorophenol | < 0.35 | 0.35 | |
| Diethylphthalate | < 0.35 | 0.35 | |
| 2,4-Dimethylphenol | < 0.35 | 0.35 | |
| Dimethylphthalate | < 0.35 | 0.35 | |
| Di-n-butylphthalate | < 0.35 | 0.35 | |

**8270 continued...**

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|----------------------------------|------------------------|--------------------|-------|
| 4,6-Dinitro-2-methylphenol | < 0.043 | 0.043 | |
| 2,4-Dinitrophenol | < 0.069 | 0.069 | |
| 2,4-Dinitrotoluene | < 0.056 | 0.056 | |
| 2,6-Dinitrotoluene | < 0.35 | 0.35 | |
| Di-n-octylphthalate | < 0.35 | 0.35 | |
| bis(2-Ethylhexyl)phthalate | < 0.35 | 0.35 | |
| Fluoranthene | < 0.35 | 0.35 | |
| Fluorene | < 0.35 | 0.35 | |
| Hexachloro-1,3-butadiene | < 0.069 | 0.069 | |
| Hexachlorobenzene | < 0.069 | 0.069 | |
| Hexachlorocyclopentadiene | < 0.35 | 0.35 | |
| Hexachloroethane | < 0.069 | 0.069 | |
| Indeno(1,2,3-cd)pyrene | < 0.35 | 0.35 | |
| Isophorone | < 0.35 | 0.35 | |
| 2-Methylphenol (o-Cresol) | < 0.35 | 0.35 | |
| 3&4-Methylphenol | < 0.69 | 0.69 | |
| 1-Methylnaphthalene | < 0.35 | 0.35 | |
| 2-Methylnaphthalene | < 0.35 | 0.35 | |
| Naphthalene | < 0.069 | 0.069 | |
| 2-Nitroaniline | < 1.39 | 1.39 | |
| 3-Nitroaniline | < 1.74 | 1.74 | |
| 4-Nitroaniline | < 0.069 | 0.069 | |
| Nitrobenzene | < 0.03 | 0.03 | |
| 2-Nitrophenol | < 0.35 | 0.35 | |
| 4-Nitrophenol | < 1.74 | 1.74 | |
| N-Nitroso-di-n-propylamine | < 0.069 | 0.069 | |
| N-Nitrosodiphenylamine | < 0.35 | 0.35 | |
| Pentachlorophenol | < 0.069 | 0.069 | |
| Phenanthrone | < 0.35 | 0.35 | |
| Phenol | < 0.35 | 0.35 | |
| Pyrene | < 0.35 | 0.35 | |
| 1,2,4-Trichlorobenzene | < 0.35 | 0.35 | |
| 2,4,5-Trichlorophenol | < 0.35 | 0.35 | |
| 2,4,6-Trichlorophenol | < 0.35 | 0.35 | |
| 2-Fluorophenol (surrogate) | 36% | | |
| Phenol-d6 (surrogate) | 37% | | |
| Nitrobenzene-d5 (surrogate) | 27% | | |
| 2-Fluorobiphenyl (surrogate) | 34% | | |
| 2,4,6-Tribromophenol (surrogate) | 36% | | |
| p-Terphenyl-d14 (surrogate) | 31% | | |
| Analysis Date: | 6/27/2023 | | |
| Analysis Time: | 2:36 | | |
| Analyst Initials: | JAK | | |
| Date Extracted: | 6/26/23 | | |
| Initial Sample Weight (g): | 30 | | |
| Final Volume (mL): | 1 | | |
| Percent Solids | 96% | | |



Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 6010B

Prep Method: EPA 3050B

| | | | | |
|--------------------------------|----------|-------------------------------------|---------|-------|
| Client Sample ID: | SB-2-2-4 | Sample Collection Date/Time: | 6/21/23 | 17:05 |
| Envision Sample Number: | 23-12295 | Sample Received Date/Time: | 6/22/23 | 9:25 |
| Sample Matrix: | soil | | | |

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|------------------|-------------------------------|---------------------------|--------------|
| Arsenic | < 2 | 2 | |
| Barium | 4.7 | 2 | |
| Cadmium | < 2 | 2 | |
| Chromium | 7.3 | 2 | |
| Lead | < 2 | 2 | |
| Selenium | < 2 | 2 | |
| Silver | < 2 | 2 | |

Analysis Date:

Analysis Time: 6-27-23/14:23

Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Weight: 1.0 g

Final Volume: 50 mL

Analytical Batch: 062723icp

Analytical Method: EPA 7471A

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|------------------|-------------------------------|---------------------------|--------------|
| Mercury | < 1 | 1 | |

Hg Analysis Date: 6/28/2023

Hg Analysis Time: 11:52

Hg Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Weight: 0.6 g

Final Volume: 50 mL

Analytical Batch: 062823hg

Percent Solids 96%

All results reported on dry weight basis.



Analytical Report

ENVision Laboratories, Inc.
1439 Sadlier Circle West Drive
Indianapolis, IN 46239
Tel: 317.351.8632
Fax: 317.351.8639
www.envisionlaboratories.com

Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

| | | | | |
|--------------------------------|----------|-------------------------------------|---------|-------|
| Client Sample ID: | SB-2-2-4 | Sample Collection Date/Time: | 6/21/23 | 17:05 |
| Envision Sample Number: | 23-12295 | Sample Received Date/Time: | 6/22/23 | 9:25 |
| Sample Matrix: | soil | | | |

| Analyte | Sample Results | Flags | Method |
|------------------|-----------------------|--------------|---------------|
| Percent Moisture | 4.0% | | EPA 1684 |
| Percent Solids | 96.0% | | EPA 1684 |
| Analysis Date: | 6/23/23 | | |
| Analyst Initials | NR | | |



Analytical Report

ENVision Laboratories, Inc.
1439 Sadlier Circle West Drive
Indianapolis, IN 46239
Tel: 317.351.8632
Fax: 317.351.8639
www.envisionlaboratories.com

Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 8260

Prep Method: EPA 5035A

Analytical Batch: 062323VS

| | | | | |
|--------------------------------|----------|-------------------------------------|---------|-------|
| Client Sample ID: | SB-3-2-4 | Sample Collection Date/Time: | 6/21/23 | 11:05 |
| Envision Sample Number: | 23-12296 | Sample Received Date/Time: | 6/22/23 | 9:25 |
| Sample Matrix: | soil | | | |

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------|------------------------|--------------------|-------|
| Acetone | < 0.119 | 0.119 | |
| Acrolein | < 0.00020 | 0.001 | 1 |
| Acrylonitrile | < 0.002 | 0.002 | |
| Benzene | < 0.006 | 0.006 | |
| Bromobenzene | < 0.006 | 0.006 | |
| Bromochloromethane | < 0.006 | 0.006 | |
| Bromodichloromethane | < 0.006 | 0.006 | |
| Bromoform | < 0.006 | 0.006 | |
| Bromomethane | < 0.006 | 0.006 | |
| n-Butanol | < 0.060 | 0.060 | |
| 2-Butanone (MEK) | < 0.012 | 0.012 | |
| n-Butylbenzene | < 0.006 | 0.006 | |
| sec-Butylbenzene | < 0.006 | 0.006 | |
| tert-Butylbenzene | < 0.006 | 0.006 | |
| Carbon Disulfide | < 0.006 | 0.006 | |
| Carbon Tetrachloride | < 0.006 | 0.006 | |
| Chlorobenzene | < 0.006 | 0.006 | |
| Chloroethane | < 0.006 | 0.006 | |
| 2-Chloroethylvinylether | < 0.060 | 0.060 | |
| Chloroform | < 0.006 | 0.006 | |
| Chloromethane | < 0.006 | 0.006 | |
| 2-Chlorotoluene | < 0.006 | 0.006 | |
| 4-Chlorotoluene | < 0.006 | 0.006 | |
| 1,2-Dibromo-3-chloropropane | < 0.0020 | 0.0020 | |
| Dibromochloromethane | < 0.006 | 0.006 | |
| 1,2-Dibromoethane (EDB) | < 0.00033 | 0.001 | 1 |
| Dibromomethane | < 0.006 | 0.006 | |
| 1,2-Dichlorobenzene | < 0.006 | 0.006 | |
| 1,3-Dichlorobenzene | < 0.006 | 0.006 | |
| 1,4-Dichlorobenzene | < 0.006 | 0.006 | |
| trans-1,4-Dichloro-2-butene | < 0.006 | 0.006 | |
| Dichlorodifluoromethane | < 0.006 | 0.006 | |
| 1,1-Dichloroethane | < 0.006 | 0.006 | |
| 1,2-Dichloroethane | < 0.006 | 0.006 | |
| 1,1-Dichloroethene | < 0.006 | 0.006 | |

**8260 continued...**

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------------|------------------------|--------------------|-------|
| cis-1,2-Dichloroethene | < 0.006 | 0.006 | |
| trans-1,2-Dichloroethene | < 0.006 | 0.006 | |
| 1,2-Dichloropropane | < 0.006 | 0.006 | |
| 1,3-Dichloropropane | < 0.006 | 0.006 | |
| 2,2-Dichloropropane | < 0.006 | 0.006 | |
| 1,1-Dichloropropene | < 0.006 | 0.006 | |
| 1,3-Dichloropropene | < 0.006 | 0.006 | |
| Ethylbenzene | < 0.006 | 0.006 | |
| Ethyl methacrylate | < 0.119 | 0.119 | |
| Hexachloro-1,3-butadiene | < 0.006 | 0.006 | |
| n-Hexane | < 0.012 | 0.012 | |
| 2-Hexanone | < 0.012 | 0.012 | |
| Iodomethane | < 0.012 | 0.012 | |
| Isopropylbenzene (Cumene) | < 0.006 | 0.006 | |
| p-Isopropyltoluene | < 0.006 | 0.006 | |
| Methylene chloride | < 0.024 | 0.024 | |
| 4-Methyl-2-pentanone (MIBK) | < 0.012 | 0.012 | |
| Methyl-tert-butyl-ether | < 0.006 | 0.006 | |
| n-Propylbenzene | < 0.006 | 0.006 | |
| Styrene | < 0.006 | 0.006 | |
| 1,1,1,2-Tetrachloroethane | < 0.006 | 0.006 | |
| 1,1,2,2-Tetrachloroethane | < 0.006 | 0.006 | |
| Tetrachloroethene | < 0.006 | 0.006 | |
| Toluene | < 0.006 | 0.006 | |
| 1,2,3-Trichlorobenzene | < 0.006 | 0.006 | |
| 1,2,4-Trichlorobenzene | < 0.006 | 0.006 | |
| 1,1,1-Trichloroethane | < 0.006 | 0.006 | |
| 1,1,2-Trichloroethane | < 0.006 | 0.006 | |
| Trichloroethene | < 0.006 | 0.006 | |
| Trichlorofluoromethane | < 0.006 | 0.006 | |
| 1,2,3-Trichloropropane | < 0.006 | 0.006 | |
| 1,2,4-Trimethylbenzene | < 0.006 | 0.006 | |
| 1,3,5-Trimethylbenzene | < 0.006 | 0.006 | |
| Vinyl acetate | < 0.012 | 0.012 | |
| Vinyl chloride | < 0.002 | 0.002 | |
| Xylene, M&P | < 0.006 | 0.006 | |
| Xylene, Ortho | < 0.006 | 0.006 | |
| Xylene, Total | < 0.012 | 0.012 | |
| Dibromofluoromethane (surrogate) | 96% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 96% | | |
| Toluene-d8 (surrogate) | 95% | | |
| 4-bromofluorobenzene (surrogate) | 101% | | |
| Analysis Date: | 6/23/23 | | |
| Analysis Time: | 22:21 | | |
| Analyst Initials | tjg | | |
| Percent Solids: | 84% | | |

All results reported on dry weight basis.



Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 8270 PAH

Prep Method: EPA 3550C

Analytical Batch: 062623BS

Client Sample ID: SB-3-2-4 **Sample Collection Date/Time:** 6/21/23 11:05
Envision Sample Number: 23-12296 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: soil

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------|------------------------|--------------------|-------|
| Acenaphthene | < 0.40 | 0.40 | |
| Acenaphthylene | < 0.40 | 0.40 | |
| Aniline | < 0.40 | 0.40 | |
| Anthracene | < 0.40 | 0.40 | |
| Benzo(a)anthracene | < 0.40 | 0.40 | |
| Benzo(a)pyrene | < 0.079 | 0.079 | |
| Benzo(b)fluoranthene | < 0.40 | 0.40 | |
| Benzo(g,h,i)perylene | < 0.40 | 0.40 | |
| Benzo(k)fluoranthene | < 0.40 | 0.40 | |
| Benzoic Acid | < 1.98 | 1.98 | |
| Benzyl Alcohol | < 0.79 | 0.79 | |
| 4-Bromophenylphenyl ether | < 0.40 | 0.40 | |
| Butylbenzylphthalate | < 0.40 | 0.40 | |
| Carbazole | < 0.79 | 0.79 | |
| 4-Chloro-3-methylphenol | < 0.79 | 0.79 | |
| 4-Chloroaniline | < 0.032 | 0.039 | 1 |
| bis(2-Chloroethoxy)methane | < 0.079 | 0.079 | |
| bis(2-Chloroethyl)ether | < 0.079 | 0.079 | |
| bis(2-Chloroisopropyl)ether | < 0.40 | 0.40 | |
| 2-Chloronaphthalene | < 0.40 | 0.40 | |
| 2-Chlorophenol | < 0.40 | 0.40 | |
| 4-Chlorophenylphenyl ether | < 0.40 | 0.40 | |
| Chrysene | < 0.40 | 0.40 | |
| Dibenzo(a,h)anthracene | < 0.079 | 0.079 | |
| Dibenzofuran | < 0.40 | 0.40 | |
| 1,2-Dichlorobenzene | < 0.40 | 0.40 | |
| 1,3-Dichlorobenzene | < 0.40 | 0.40 | |
| 1,4-Dichlorobenzene | < 0.40 | 0.40 | |
| 3,3-Dichlorobenzidine | < 0.17 | 0.79 | |
| 2,4-Dichlorophenol | < 0.40 | 0.40 | |
| Diethylphthalate | < 0.40 | 0.40 | |
| 2,4-Dimethylphenol | < 0.40 | 0.40 | |
| Dimethylphthalate | < 0.40 | 0.40 | |
| Di-n-butylphthalate | < 0.40 | 0.40 | |

**8270 continued...**

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|----------------------------------|------------------------|--------------------|-------|
| 4,6-Dinitro-2-methylphenol | < 0.049 | 0.049 | |
| 2,4-Dinitrophenol | < 0.079 | 0.079 | |
| 2,4-Dinitrotoluene | < 0.064 | 0.064 | |
| 2,6-Dinitrotoluene | < 0.40 | 0.40 | |
| Di-n-octylphthalate | < 0.40 | 0.40 | |
| bis(2-Ethylhexyl)phthalate | < 0.40 | 0.40 | |
| Fluoranthene | 0.762 | 0.40 | |
| Fluorene | < 0.40 | 0.40 | |
| Hexachloro-1,3-butadiene | < 0.079 | 0.079 | |
| Hexachlorobenzene | < 0.079 | 0.079 | |
| Hexachlorocyclopentadiene | < 0.40 | 0.40 | |
| Hexachloroethane | < 0.079 | 0.079 | |
| Indeno(1,2,3-cd)pyrene | < 0.40 | 0.40 | |
| Isophorone | < 0.40 | 0.40 | |
| 2-Methylphenol (o-Cresol) | < 0.40 | 0.40 | |
| 3&4-Methylphenol | < 0.79 | 0.79 | |
| 1-Methylnaphthalene | < 0.40 | 0.40 | |
| 2-Methylnaphthalene | < 0.40 | 0.40 | |
| Naphthalene | < 0.079 | 0.079 | |
| 2-Nitroaniline | < 1.59 | 1.59 | |
| 3-Nitroaniline | < 1.98 | 1.98 | |
| 4-Nitroaniline | < 0.079 | 0.079 | |
| Nitrobenzene | < 0.04 | 0.04 | |
| 2-Nitrophenol | < 0.40 | 0.40 | |
| 4-Nitrophenol | < 1.98 | 1.98 | |
| N-Nitroso-di-n-propylamine | < 0.079 | 0.079 | |
| N-Nitrosodiphenylamine | < 0.40 | 0.40 | |
| Pentachlorophenol | < 0.079 | 0.079 | |
| Phenanthrene | 0.668 | 0.40 | |
| Phenol | < 0.40 | 0.40 | |
| Pyrene | 0.678 | 0.40 | |
| 1,2,4-Trichlorobenzene | < 0.40 | 0.40 | |
| 2,4,5-Trichlorophenol | < 0.40 | 0.40 | |
| 2,4,6-Trichlorophenol | < 0.40 | 0.40 | |
| 2-Fluorophenol (surrogate) | 34% | | |
| Phenol-d6 (surrogate) | 32% | | |
| Nitrobenzene-d5 (surrogate) | 21% | | |
| 2-Fluorobiphenyl (surrogate) | 25% | | |
| 2,4,6-Tribromophenol (surrogate) | 31% | | |
| p-Terphenyl-d14 (surrogate) | 24% | | |
| Analysis Date: | 6/27/2023 | | |
| Analysis Time: | 3:02 | | |
| Analyst Initials: | JAK | | |
| Date Extracted: | 6/26/23 | | |
| Initial Sample Weight (g): | 30 | | |
| Final Volume (mL): | 1 | | |
| Percent Solids | 84% | | |



Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 6010B

Prep Method: EPA 3050B

Client Sample ID: SB-3-2-4 **Sample Collection Date/Time:** 6/21/23 11:05

Envision Sample Number: 23-12296 **Sample Received Date/Time:** 6/22/23 9:25

Sample Matrix: soil

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|------------------|-------------------------------|---------------------------|--------------|
| Arsenic | < 2 | 2 | |
| Barium | 17 | 2 | |
| Cadmium | < 2 | 2 | |
| Chromium | 4.0 | 2 | |
| Lead | 8.9 | 2 | |
| Selenium | < 2 | 2 | |
| Silver | < 2 | 2 | |

Analysis Date:

Analysis Time: 6-27-23/14:26

Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Weight: 1.0 g

Final Volume: 50 mL

Analytical Batch: 062723icp

Analytical Method: EPA 7471A

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|------------------|-------------------------------|---------------------------|--------------|
| Mercury | < 1 | 1 | |

Hg Analysis Date: 6/28/2023

Hg Analysis Time: 11:53

Hg Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Weight: 0.6 g

Final Volume: 50 mL

Analytical Batch: 062823hg

Percent Solids 84%

All results reported on dry weight basis.



Analytical Report

ENVision Laboratories, Inc.
1439 Sadlier Circle West Drive
Indianapolis, IN 46239
Tel: 317.351.8632
Fax: 317.351.8639
www.envisionlaboratories.com

Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

| | | | | |
|--------------------------------|----------|-------------------------------------|---------|-------|
| Client Sample ID: | SB-3-2-4 | Sample Collection Date/Time: | 6/21/23 | 11:05 |
| Envision Sample Number: | 23-12296 | Sample Received Date/Time: | 6/22/23 | 9:25 |
| Sample Matrix: | soil | | | |

| Analyte | Sample Results | Flags | Method |
|------------------|-----------------------|--------------|---------------|
| Percent Moisture | 16.0% | | EPA 1684 |
| Percent Solids | 84.0% | | EPA 1684 |
| Analysis Date: | 6/23/23 | | |
| Analyst Initials | NR | | |



Analytical Report

ENVision Laboratories, Inc.
1439 Sadlier Circle West Drive
Indianapolis, IN 46239
Tel: 317.351.8632
Fax: 317.351.8639
www.envisionlaboratories.com

Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 8260

Prep Method: EPA 5035A

Analytical Batch: 062323VS

| | | | | |
|--------------------------------|----------|-------------------------------------|---------|-------|
| Client Sample ID: | SB-6-2-4 | Sample Collection Date/Time: | 6/21/23 | 14:30 |
| Envision Sample Number: | 23-12297 | Sample Received Date/Time: | 6/22/23 | 9:25 |
| Sample Matrix: | soil | | | |

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------|------------------------|--------------------|-------|
| Acetone | < 0.110 | 0.110 | |
| Acrolein | < 0.00019 | 0.001 | 1 |
| Acrylonitrile | < 0.002 | 0.002 | |
| Benzene | < 0.005 | 0.005 | |
| Bromobenzene | < 0.005 | 0.005 | |
| Bromochloromethane | < 0.005 | 0.005 | |
| Bromodichloromethane | < 0.005 | 0.005 | |
| Bromoform | < 0.005 | 0.005 | |
| Bromomethane | < 0.005 | 0.005 | |
| n-Butanol | < 0.055 | 0.055 | |
| 2-Butanone (MEK) | < 0.011 | 0.011 | |
| n-Butylbenzene | < 0.005 | 0.005 | |
| sec-Butylbenzene | 0.0137 | 0.005 | |
| tert-Butylbenzene | < 0.005 | 0.005 | |
| Carbon Disulfide | < 0.005 | 0.005 | |
| Carbon Tetrachloride | < 0.005 | 0.005 | |
| Chlorobenzene | < 0.005 | 0.005 | |
| Chloroethane | < 0.005 | 0.005 | |
| 2-Chloroethylvinylether | < 0.055 | 0.055 | |
| Chloroform | < 0.005 | 0.005 | |
| Chloromethane | < 0.005 | 0.005 | |
| 2-Chlorotoluene | < 0.005 | 0.005 | |
| 4-Chlorotoluene | < 0.005 | 0.005 | |
| 1,2-Dibromo-3-chloropropane | < 0.0019 | 0.0019 | |
| Dibromochloromethane | < 0.005 | 0.005 | |
| 1,2-Dibromoethane (EDB) | < 0.00031 | 0.001 | 1 |
| Dibromomethane | < 0.005 | 0.005 | |
| 1,2-Dichlorobenzene | < 0.005 | 0.005 | |
| 1,3-Dichlorobenzene | < 0.005 | 0.005 | |
| 1,4-Dichlorobenzene | < 0.005 | 0.005 | |
| trans-1,4-Dichloro-2-butene | < 0.005 | 0.005 | |
| Dichlorodifluoromethane | < 0.005 | 0.005 | |
| 1,1-Dichloroethane | < 0.005 | 0.005 | |
| 1,2-Dichloroethane | < 0.005 | 0.005 | |
| 1,1-Dichloroethene | < 0.005 | 0.005 | |

**8260 continued...**

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------------|------------------------|--------------------|-------|
| cis-1,2-Dichloroethene | < 0.005 | 0.005 | |
| trans-1,2-Dichloroethene | < 0.005 | 0.005 | |
| 1,2-Dichloropropane | < 0.005 | 0.005 | |
| 1,3-Dichloropropane | < 0.005 | 0.005 | |
| 2,2-Dichloropropane | < 0.005 | 0.005 | |
| 1,1-Dichloropropene | < 0.005 | 0.005 | |
| 1,3-Dichloropropene | < 0.005 | 0.005 | |
| Ethylbenzene | < 0.005 | 0.005 | |
| Ethyl methacrylate | < 0.110 | 0.110 | |
| Hexachloro-1,3-butadiene | < 0.005 | 0.005 | |
| n-Hexane | < 0.011 | 0.011 | |
| 2-Hexanone | < 0.011 | 0.011 | |
| Iodomethane | < 0.011 | 0.011 | |
| Isopropylbenzene (Cumene) | < 0.005 | 0.005 | |
| p-Isopropyltoluene | 0.0872 | 0.005 | |
| Methylene chloride | < 0.022 | 0.022 | |
| 4-Methyl-2-pentanone (MIBK) | < 0.011 | 0.011 | |
| Methyl-tert-butyl-ether | < 0.005 | 0.005 | |
| n-Propylbenzene | 0.00720 | 0.005 | |
| Styrene | < 0.005 | 0.005 | |
| 1,1,1,2-Tetrachloroethane | < 0.005 | 0.005 | |
| 1,1,2,2-Tetrachloroethane | < 0.005 | 0.005 | |
| Tetrachloroethene | < 0.005 | 0.005 | |
| Toluene | < 0.005 | 0.005 | |
| 1,2,3-Trichlorobenzene | < 0.005 | 0.005 | |
| 1,2,4-Trichlorobenzene | < 0.005 | 0.005 | |
| 1,1,1-Trichloroethane | < 0.005 | 0.005 | |
| 1,1,2-Trichloroethane | < 0.005 | 0.005 | |
| Trichloroethene | < 0.005 | 0.005 | |
| Trichlorofluoromethane | < 0.005 | 0.005 | |
| 1,2,3-Trichloropropane | < 0.005 | 0.005 | |
| 1,2,4-Trimethylbenzene | 0.211 | 0.005 | |
| 1,3,5-Trimethylbenzene | 0.622 | 0.275 | 2 |
| Vinyl acetate | < 0.011 | 0.011 | |
| Vinyl chloride | < 0.002 | 0.002 | |
| Xylene, M&P | 0.0113 | 0.005 | |
| Xylene, Ortho | 0.131 | 0.005 | |
| Xylene, Total | 0.142 | 0.011 | |
| Dibromofluoromethane (surrogate) | 112% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 107% | | |
| Toluene-d8 (surrogate) | 108% | | |
| 4-bromofluorobenzene (surrogate) | 108% | | |
| Analysis Date: | 6/23/23 | | |
| Analysis Time: | 22:37 | | |
| Analyst Initials | tjg | | |
| Percent Solids: | 91% | | |

All results reported on dry weight basis.



Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 8270 PAH

Prep Method: EPA 3550C

Analytical Batch: 062623BS

Client Sample ID: SB-6-2-4 **Sample Collection Date/Time:** 6/21/23 14:30
Envision Sample Number: 23-12297 **Sample Received Date/Time:** 6/22/23 9:25
Sample Matrix: soil

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|-----------------------------|------------------------|--------------------|-------|
| Acenaphthene | < 0.37 | 0.37 | |
| Acenaphthylene | < 0.37 | 0.37 | |
| Aniline | < 0.37 | 0.37 | |
| Anthracene | < 0.37 | 0.37 | |
| Benzo(a)anthracene | < 0.37 | 0.37 | |
| Benzo(a)pyrene | < 0.073 | 0.073 | |
| Benzo(b)fluoranthene | < 0.37 | 0.37 | |
| Benzo(g,h,i)perylene | < 0.37 | 0.37 | |
| Benzo(k)fluoranthene | < 0.37 | 0.37 | |
| Benzoic Acid | < 1.83 | 1.83 | |
| Benzyl Alcohol | < 0.73 | 0.73 | |
| 4-Bromophenylphenyl ether | < 0.37 | 0.37 | |
| Butylbenzylphthalate | < 0.37 | 0.37 | |
| Carbazole | < 0.73 | 0.73 | |
| 4-Chloro-3-methylphenol | < 0.73 | 0.73 | |
| 4-Chloroaniline | < 0.030 | 0.036 | 1 |
| bis(2-Chloroethoxy)methane | < 0.073 | 0.073 | |
| bis(2-Chloroethyl)ether | < 0.073 | 0.073 | |
| bis(2-Chloroisopropyl)ether | < 0.37 | 0.37 | |
| 2-Chloronaphthalene | < 0.37 | 0.37 | |
| 2-Chlorophenol | < 0.37 | 0.37 | |
| 4-Chlorophenylphenyl ether | < 0.37 | 0.37 | |
| Chrysene | < 0.37 | 0.37 | |
| Dibenzo(a,h)anthracene | < 0.073 | 0.073 | |
| Dibenzofuran | < 0.37 | 0.37 | |
| 1,2-Dichlorobenzene | < 0.37 | 0.37 | |
| 1,3-Dichlorobenzene | < 0.37 | 0.37 | |
| 1,4-Dichlorobenzene | < 0.37 | 0.37 | |
| 3,3-Dichlorobenzidine | < 0.15 | 0.73 | |
| 2,4-Dichlorophenol | < 0.37 | 0.37 | |
| Diethylphthalate | < 0.37 | 0.37 | |
| 2,4-Dimethylphenol | < 0.37 | 0.37 | |
| Dimethylphthalate | < 0.37 | 0.37 | |
| Di-n-butylphthalate | < 0.37 | 0.37 | |

**8270 continued...**

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|----------------------------------|------------------------|--------------------|-------|
| 4,6-Dinitro-2-methylphenol | < 0.045 | 0.045 | |
| 2,4-Dinitrophenol | < 0.073 | 0.073 | |
| 2,4-Dinitrotoluene | < 0.059 | 0.059 | |
| 2,6-Dinitrotoluene | < 0.37 | 0.37 | |
| Di-n-octylphthalate | < 0.37 | 0.37 | |
| bis(2-Ethylhexyl)phthalate | < 0.37 | 0.37 | |
| Fluoranthene | < 0.37 | 0.37 | |
| Fluorene | < 0.37 | 0.37 | |
| Hexachloro-1,3-butadiene | < 0.073 | 0.073 | |
| Hexachlorobenzene | < 0.073 | 0.073 | |
| Hexachlorocyclopentadiene | < 0.37 | 0.37 | |
| Hexachloroethane | < 0.073 | 0.073 | |
| Indeno(1,2,3-cd)pyrene | < 0.37 | 0.37 | |
| Isophorone | < 0.37 | 0.37 | |
| 2-Methylphenol (o-Cresol) | < 0.37 | 0.37 | |
| 3&4-Methylphenol | < 0.73 | 0.73 | |
| 1-Methylnaphthalene | < 0.37 | 0.37 | |
| 2-Methylnaphthalene | < 0.37 | 0.37 | |
| Naphthalene | 0.893 | 0.073 | |
| 2-Nitroaniline | < 1.47 | 1.47 | |
| 3-Nitroaniline | < 1.83 | 1.83 | |
| 4-Nitroaniline | < 0.073 | 0.073 | |
| Nitrobenzene | < 0.04 | 0.04 | |
| 2-Nitrophenol | < 0.37 | 0.37 | |
| 4-Nitrophenol | < 1.83 | 1.83 | |
| N-Nitroso-di-n-propylamine | < 0.073 | 0.073 | |
| N-Nitrosodiphenylamine | < 0.37 | 0.37 | |
| Pentachlorophenol | < 0.073 | 0.073 | |
| Phenanthrone | < 0.37 | 0.37 | |
| Phenol | < 0.37 | 0.37 | |
| Pyrene | < 0.37 | 0.37 | |
| 1,2,4-Trichlorobenzene | < 0.37 | 0.37 | |
| 2,4,5-Trichlorophenol | < 0.37 | 0.37 | |
| 2,4,6-Trichlorophenol | < 0.37 | 0.37 | |
| 2-Fluorophenol (surrogate) | 28% | | |
| Phenol-d6 (surrogate) | 31% | | |
| Nitrobenzene-d5 (surrogate) | 53% | | |
| 2-Fluorobiphenyl (surrogate) | 50% | | |
| 2,4,6-Tribromophenol (surrogate) | 45% | | |
| p-Terphenyl-d14 (surrogate) | 42% | | |
| Analysis Date: | 6/27/2023 | | |
| Analysis Time: | 3:28 | | |
| Analyst Initials: | JAK | | |
| Date Extracted: | 6/26/23 | | |
| Initial Sample Weight (g): | 30 | | |
| Final Volume (mL): | 1 | | |
| Percent Solids | 91% | | |



Client Name: AUGUST MACK ENVIRONMENTAL

Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

Analytical Method: EPA 6010B

Prep Method: EPA 3050B

Client Sample ID: SB-6-2-4 **Sample Collection Date/Time:** 6/21/23 14:30

Envision Sample Number: 23-12297 **Sample Received Date/Time:** 6/22/23 9:25

Sample Matrix: soil

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|------------------|-------------------------------|---------------------------|--------------|
| Arsenic | < 2 | 2 | |
| Barium | 196 | 2 | |
| Cadmium | < 2 | 2 | |
| Chromium | 43 | 2 | |
| Lead | 36 | 2 | |
| Selenium | < 2 | 2 | |
| Silver | < 2 | 2 | |

Analysis Date:

Analysis Time: 6-27-23/14:28

Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Weight: 1.0 g

Final Volume: 50 mL

Analytical Batch: 062723icp

Analytical Method: EPA 7471A

| Compounds | Sample Results (mg/kg) | Rep. Limit (mg/kg) | Flags |
|------------------|-------------------------------|---------------------------|--------------|
| Mercury | < 1 | 1 | |

Hg Analysis Date: 6/28/2023

Hg Analysis Time: 11:55

Hg Analyst Initials: gjd

Date Digested: 6/27/2023

Initial Sample Weight: 0.6 g

Final Volume: 50 mL

Analytical Batch: 062823hg

Percent Solids 91%

All results reported on dry weight basis.



Analytical Report

ENVision Laboratories, Inc.
1439 Sadlier Circle West Drive
Indianapolis, IN 46239
Tel: 317.351.8632
Fax: 317.351.8639
www.envisionlaboratories.com

Client Name: AUGUST MACK ENVIRONMENTAL

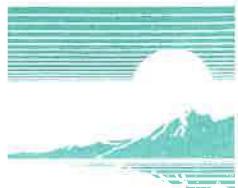
Project ID: JX1600

Client Project Manager: TYLER ZSCHIEDRICH

ENVision Project Number: 2023-1274

| | | | | |
|--------------------------------|----------|-------------------------------------|---------|-------|
| Client Sample ID: | SB-6-2-4 | Sample Collection Date/Time: | 6/21/23 | 14:30 |
| Envision Sample Number: | 23-12297 | Sample Received Date/Time: | 6/22/23 | 9:25 |
| Sample Matrix: | soil | | | |

| Analyte | Sample Results | Flags | Method |
|------------------|-----------------------|--------------|---------------|
| Percent Moisture | 9.0% | | EPA 1684 |
| Percent Solids | 91.0% | | EPA 1684 |
| Analysis Date: | 6/23/23 | | |
| Analyst Initials | NR | | |



**First
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IL ELAP / NELAC Accreditation # 100292

July 07, 2023

Ms. Cheryl Crum
ENVISION LABORATORIES, INC.
1439 Sandlier Cir. W. Drive
Indianapolis, IN 46239

Project ID: 2023-1274
First Environmental File ID: 23-5329
Date Received: June 23, 2023

Dear Ms. Cheryl Crum:

The above referenced project was analyzed as directed on the enclosed chain of custody record.

All Quality Control criteria as outlined in the methods and current IL ELAP/NELAP have been met unless otherwise noted. QA/QC documentation and raw data will remain on file for future reference. Our accreditation number is 100292 and our current certificate is number:

1002922023-10: effective 03/07/2023 through 02/28/2024.

I thank you for the opportunity to be of service to you and look forward to working with you again in the future. Should you have any questions regarding any of the enclosed analytical data or need additional information, please contact me at (630) 778-1200.

Sincerely,

Joy Geraci
Project Manager



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IL ELAP / NELAC Accreditation # 100292

Case Narrative

ENVISION LABORATORIES, INC.

Project ID: **2023-1274**

Lab File ID: **23-5329**

Date Received: **June 23, 2023**

All quality control criteria, as outlined in the methods, have been met except as noted below or on the following analytical report.

The results in this report apply to the samples in the following table:

| Laboratory Sample ID | Client Sample Identifier | Date/Time Collected | |
|----------------------|--------------------------|---------------------|-------|
| 23-5329-001 | 23-12293/SB-7-1-1.5 | 6/21/2023 | 15:25 |
| 23-5329-002 | 23-12294/SB-1-2-4 | 6/21/2023 | 15:55 |
| 23-5329-003 | 23-12295/SB-2-2-4 | 6/21/2023 | 17:05 |
| 23-5329-004 | 23-12296/SB-3-2-4 | 6/21/2023 | 11:05 |
| 23-5329-005 | 23-12297/SB-6-2-4 | 6/21/2023 | 14:30 |

Sample Batch Comments:

Sample acceptance criteria were met.



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Case Narrative

ENVISION LABORATORIES, INC.

Lab File ID: **23-5329**

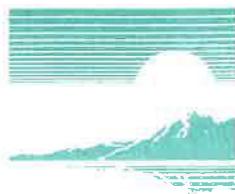
Project ID: **2023-1274**

Date Received: **June 23, 2023**

All quality control criteria, as outlined in the methods, have been met except as noted below or on the following analytical report.

The following is a definition of flags that may be used in this report:

| Flag | Description | Flag | Description |
|------|----------------------------------------------------------------------------------------------------------------------|------|----------------------------------------------------------------------------------------------------------|
| A | Method holding time is 15 minutes from collection. Lab analysis was performed as soon as possible. | | |
| B | Analyte was found in the method blank. | L | LCS recovery outside control limits. |
| < | Analyte not detected at or above the reporting limit. | M | MS recovery outside control limits; LCS acceptable. |
| C | Sample received in an improper container for this test. | P | Chemical preservation pH adjusted in lab. |
| D | Surrogates diluted out; recovery not available. | Q | Result was determined by a GC/MS database search. |
| E | Estimated result; concentration exceeds calibration range. | S | Analysis was subcontracted to another laboratory. |
| G | Surrogate recovery outside control limits. | T | Result is less than three times the MDL value. |
| H | Analysis or extraction holding time exceeded. | W | Reporting limit elevated due to sample matrix. |
| I | ICVS % rec outside 95-105% but within 90-110% | | |
| J | Estimated result; concentration is less than routine RL but greater than MDL. | N | Analyte is not part of our NELAC accreditation or accreditation may not be available for this parameter. |
| RL | Routine Reporting Limit (Lowest amount that can be detected when routine weights/volumes are used without dilution.) | ND | Analyte was not detected using a library search routine; No calibration standard was analyzed. |



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Analytical Report

Client: ENVISION LABORATORIES, INC.

Date Collected: 06/21/23

Project ID: 2023-1274

Time Collected: 15:25

Sample ID: 23-12293/SB-7-1-1.5

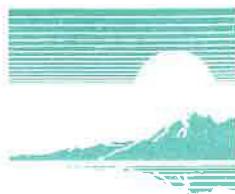
Date Received: 06/23/23

Sample No: 23-5329-001

Date Reported: 07/07/23

Results are reported on a dry weight basis.

| Analyte | Method | Result | R.L. | Units | Flags |
|-----------------------------------------|-----------------------------|-----------|-------|-------|------------------|
| Solids, Total | Method: 2540G 2011 | | | | |
| Analysis Date: | 06/29/23 | | | | |
| Total Solids | | 90.11 | | % | |
| Polychlorinated biphenyls (PCBs) | Method: 8082 | | | | |
| Analysis Date: | 07/07/23 | | | | |
| Aroclor 1016 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1221 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1232 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1242 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1248 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1254 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1260 | | < 0.016 | 0.016 | mg/kg | |
| Chromium, Hexavalent | Method: 3060A/7196A | | | | |
| Analysis Date: | 06/29/23 | | | | |
| Chromium, Hexavalent | | < 2.5 | 2.5 | mg/kg | |
| <i>Sample QC Summary:</i> | <i>Surrogate Recovery</i> | | | | |
| | | | | | <i>%R Limits</i> |
| Method | Analyte | QC Result | | Low | High |
| 8082 | Decachlorobiphenyl (Surr) | %R: | 61.9 | | 28 - 136 |
| 8082 | Tetrachloro-m-xylene (Surr) | %R: | 81.7 | | 61 - 127 |



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Analytical Report

Client: ENVISION LABORATORIES, INC.

Date Collected: 06/21/23

Project ID: 2023-1274

Time Collected: 15:55

Sample ID: 23-12294/SB-1-2-4

Date Received: 06/23/23

Sample No: 23-5329-002

Date Reported: 07/07/23

Results are reported on a dry weight basis.

| Analyte | Method | Result | R.L. | Units | Flags |
|-----------------------------------------|-----------------------------|-----------|-------|-------|---------------------------------|
| Solids, Total | Method: 2540G 2011 | | | | |
| Analysis Date: | 06/29/23 | | | | |
| Total Solids | | 82.61 | | % | |
| Polychlorinated biphenyls (PCBs) | Method: 8082 | | | | Preparation Method 3540C |
| Analysis Date: | 07/07/23 | | | | Preparation Date: 06/26/23 |
| Aroclor 1016 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1221 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1232 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1242 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1248 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1254 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1260 | | < 0.016 | 0.016 | mg/kg | |
| Chromium, Hexavalent | Method: 3060A/7196A | | | | |
| Analysis Date: | 06/29/23 | | | | |
| Chromium, Hexavalent | | < 2.5 | 2.5 | mg/kg | |
| <i>Sample QC Summary:</i> | <i>Surrogate Recovery</i> | | | | <i>%R Limits</i> |
| Method | Analyte | QC Result | | Low | High |
| 8082 | Decachlorobiphenyl (Surr) | %R: | 56.2 | | 28 - 136 |
| 8082 | Tetrachloro-m-xylene (Surr) | %R: | 79.1 | | 61 - 127 |



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IL ELAP / NELAC Accreditation # 100292

Analytical Report

Client: ENVISION LABORATORIES, INC.

Date Collected: 06/21/23

Project ID: 2023-1274

Time Collected: 17:05

Sample ID: 23-12295/SB-2-2-4

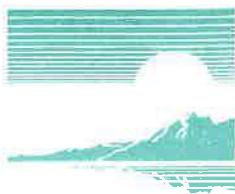
Date Received: 06/23/23

Sample No: 23-5329-003

Date Reported: 07/07/23

Results are reported on a dry weight basis.

| Analyte | Result | R.L. | Units | Flags |
|-----------------------------------------|-----------------------------|------------------|-------|------------------|
| Solids, Total | Method: 2540G 2011 | | | |
| Analysis Date: | 06/29/23 | | | |
| Total Solids | 95.34 | | % | |
| Polychlorinated biphenyls (PCBs) | Method: 8082 | | | |
| Analysis Date: | 07/07/23 | | | |
| Aroclor 1016 | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1221 | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1232 | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1242 | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1248 | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1254 | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1260 | < 0.016 | 0.016 | mg/kg | |
| Chromium, Hexavalent | Method: 3060A/7196A | | | |
| Analysis Date: | 06/29/23 | | | |
| Chromium, Hexavalent | < 2.5 | 2.5 | mg/kg | |
| <i>Sample QC Summary:</i> | <i>Surrogate Recovery</i> | | | |
| <i>Method</i> | <i>Analyte</i> | <i>QC Result</i> | | <i>%R Limits</i> |
| 8082 | Decachlorobiphenyl (Surr) | %R: | 74.6 | 28 - 136 |
| 8082 | Tetrachloro-m-xylene (Surr) | %R: | 95.5 | 61 - 127 |



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IL ELAP / NELAC Accreditation # 100292

Analytical Report

Client: ENVISION LABORATORIES, INC.

Date Collected: 06/21/23

Project ID: 2023-1274

Time Collected: 11:05

Sample ID: 23-12296/SB-3-2-4

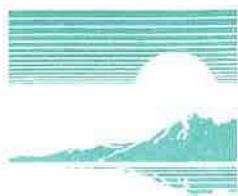
Date Received: 06/23/23

Sample No: 23-5329-004

Date Reported: 07/07/23

Results are reported on a dry weight basis.

| Analyte | Method | Result | R.L. | Units | Flags |
|-----------------------------------------|-----------------------------|-----------|-------|-------|---------------------------------|
| Solids, Total | Method: 2540G 2011 | | | | |
| Analysis Date: | 06/29/23 | | | | |
| Total Solids | | 93.93 | | % | |
| Polychlorinated biphenyls (PCBs) | Method: 8082 | | | | Preparation Method 3540C |
| Analysis Date: | 07/07/23 | | | | Preparation Date: 06/26/23 |
| Aroclor 1016 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1221 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1232 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1242 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1248 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1254 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1260 | | < 0.016 | 0.016 | mg/kg | |
| Chromium, Hexavalent | Method: 3060A/7196A | | | | |
| Analysis Date: | 06/29/23 | | | | |
| Chromium, Hexavalent | | < 2.5 | 2.5 | mg/kg | |
| <i>Sample QC Summary:</i> | <i>Surrogate Recovery</i> | | | | <i>%R Limits</i> |
| Method | Analyte | QC Result | Low | High | |
| 8082 | Decachlorobiphenyl (Surr) | 0 | | | 28 - 136 |
| 8082 | Tetrachloro-m-xylene (Surr) | 0 | | | 61 - 127 |



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Analytical Report

Client: ENVISION LABORATORIES, INC.

Date Collected: 06/21/23

Project ID: 2023-1274

Time Collected: 14:30

Sample ID: 23-12297/SB-6-2-4

Date Received: 06/23/23

Sample No: 23-5329-005

Date Reported: 07/07/23

Results are reported on a dry weight basis.

| Analyte | Method | Result | R.L. | Units | Flags |
|-----------------------------------------|-----------------------------|---------|------------------|-------|---------------------------------|
| Solids, Total | Method: 2540G 2011 | | | | |
| Analysis Date: | 06/29/23 | | | | |
| Total Solids | | 89.60 | | % | |
| Polychlorinated biphenyls (PCBs) | Method: 8082 | | | | Preparation Method 3540C |
| Analysis Date: | 07/07/23 | | | | Preparation Date: 06/26/23 |
| Aroclor 1016 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1221 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1232 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1242 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1248 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1254 | | < 0.016 | 0.016 | mg/kg | |
| Aroclor 1260 | | < 0.016 | 0.016 | mg/kg | |
| Chromium, Hexavalent | Method: 3060A/7196A | | | | |
| Analysis Date: | 06/29/23 | | | | |
| Chromium, Hexavalent | | < 2.5 | 2.5 | mg/kg | |
| <i>Sample QC Summary:</i> | <i>Surrogate Recovery</i> | | | | <i>%R Limits</i> |
| <i>Method</i> | <i>Analyte</i> | | <i>QC Result</i> | | <i>Low</i> <i>High</i> |
| 8082 | Decachlorobiphenyl (Surr) | | %R: | 35.3 | 28 - 136 |
| 8082 | Tetrachloro-m-xylene (Surr) | | %R: | 66.3 | 61 - 127 |



First Environmental Laboratories, Inc.

IL ELAP / NELAC Accreditation # 100292

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Quality Control Summary

Client: ENVISION LABORATORIES, INC.

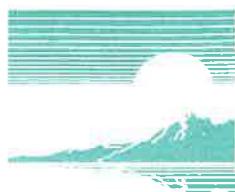
Lab File ID: 23-5329

Project ID: 2023-1274

| QC Lab# | Time QC Code | Parameter | Reported Result | Units | QC Result | %R Limits Low | %R Limits High | RPD Limit | |
|---------------|----------------------|---------------------------|-------------------------------------------------------------------------|-------|------------------|---------------|----------------|-----------|-----|
| Parameter: | Chromium, Hexavalent | Analytical Method: | 3060A/7196A | | Analytical WS #: | 233380 | Analysis Date: | 6/29/2023 | |
| 23-5329-003MS | MS | Chromium, Hex (Insoluble) | 1660 | mg/kg | %R: 98 | 75 | - | 125 | |
| | MS | Chromium, Hex (Soluble) | 27.6 | mg/kg | %R: 42 | * | 75 | - | 125 |
| | | | MS outside control limits. All other QCIs are within acceptance limits. | | | | | | |
| CCB818265 | CB | Chromium, Hexavalent | < 0.005 | mg/L | 0 | - | - | - | |
| CCB818266 | CB | Chromium, Hexavalent | < 0.005 | mg/L | 0 | - | - | - | |
| CCB818267 | CB | Chromium, Hexavalent | < 0.005 | mg/L | 0 | - | - | - | |
| CCVS818268 | CCVS | Chromium, Hexavalent | 0.122 | mg/L | %R: 97.6 | 90 | - | 110 | |
| CCVS818269 | CCVS | Chromium, Hexavalent | 0.127 | mg/L | %R: 101.6 | 90 | - | 110 | |
| CCVS818270 | CCVS | Chromium, Hexavalent | 0.127 | mg/L | %R: 101.6 | 90 | - | 110 | |
| LCS818271 | LCS | Chromium, Hex (Soluble) | 1.28 | mg/L | %R: 102.4 | 80 | - | 120 | |
| LCS818274 | LCS | Chromium, Hex (Insoluble) | 33.3 | mg/L | %R: 103.6 | 80 | - | 120 | |
| PB818273 | PB | Chromium, Hexavalent | < 0.05 | mg/L | 0 | - | - | - | |

* The QC indicator is outside control limits. %R = percent recovery; RPD = Relative percent difference
CB = Calibration Blank; CCVS = Continuing Calibration Verification Standard; MS = Matrix Spike;
MSD = Matrix Spike Duplicate; LCS = Laboratory Control Spike; SURR = Surrogate Spiking Compound;
PB = Procedure Blank; BLK = Method Blank; D = QCI diluted out.





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Quality Control Summary

Client: ENVISION LABORATORIES, INC.

Lab File ID: 23-5329

Project ID: 2023-1274

| QC Lab# | Time QC Code | Parameter | Reported Result | Units | QC Result | %R Limits Low | %R Limits High | RPD Limit |
|------------------|----------------------------------|--------------------|-----------------|-------|-------------------------|---------------|-------------------------|-----------|
| Parameter: | Polychlorinated biphenyls (PCBs) | Analytical Method: | 8082 | | Analytical WS #: 233635 | | Analysis Date: 7/7/2023 | |
| | | Prep Method: | 3540C | | Prep WS#: 39569 | | Prep Date: 6/26/2023 | |
| LCS74470 | LCS | Aroclor 1016 | | ug/kg | | 72 - 126 | | |
| | LCS | Aroclor 1260 | | ug/kg | | 56 - 121 | | |
| Method Blank7447 | BLK | Aroclor 1016 | | ug/kg | | - | | |
| | BLK | Aroclor 1221 | | ug/kg | | - | | |
| | BLK | Aroclor 1232 | | ug/kg | | - | | |
| | BLK | Aroclor 1242 | | ug/kg | | - | | |
| | BLK | Aroclor 1248 | | ug/kg | | - | | |
| | BLK | Aroclor 1254 | | ug/kg | | - | | |
| | BLK | Aroclor 1260 | | ug/kg | | - | | |
| | BLK | Aroclor 1268 | | ug/kg | | - | | |

* The QC indicator is outside control limits. %R = percent recovery; RPD = Relative percent difference
CB = Calibration Blank; CCVS = Continuing Calibration Verification Standard; MS = Matrix Spike;
MSD = Matrix Spike Duplicate; LCS = Laboratory Control Spike; SURR = Surrogate Spiking Compound;
PB = Procedure Blank; BLK = Method Blank; D = QCI diluted out.





CHAIN OF CUSTODY RECORD

ENVision Laboratories, Inc. [1439 Sadlier Circle West Drive, Indianapolis, IN 46239] Phone: 317-351-8632 Fax: 317-351-8639



ENVision Laboratories, Inc.
1439 Sadlier Circle West Drive
Indianapolis, IN 46239
Tel: 317.351.8632
Fax: 317.351.8639
www.envisionlaboratories.com

EPA 8260 Quality Control Data

ENVision Batch Number: 062323VS

| Method Blank (MB): | MB Results (ug/kg) | Rep Lim (ug/kg) | Flag |
|-----------------------------|---------------------------|------------------------|-------------|
| Acetone | < 100 | 100 | |
| Acrolein | < 0.17 | 1 | |
| Acrylonitrile | < 2 | 2 | |
| Benzene | < 5 | 5 | |
| Bromobenzene | < 5 | 5 | |
| Bromochloromethane | < 5 | 5 | |
| Bromodichloromethane | < 5 | 5 | |
| Bromoform | < 5 | 5 | |
| Bromomethane | < 5 | 5 | |
| n-Butanol | < 50 | 50 | |
| 2-Butanone (MEK) | < 10 | 10 | |
| n-Butylbenzene | < 5 | 5 | |
| sec-Butylbenzene | < 5 | 5 | |
| tert-Butylbenzene | < 5 | 5 | |
| Carbon Disulfide | < 5 | 5 | |
| Carbon Tetrachloride | < 5 | 5 | |
| Chlorobenzene | < 5 | 5 | |
| Chloroethane | < 5 | 5 | |
| 2-Chloroethylvinylether | < 50 | 50 | |
| Chloroform | < 5 | 5 | |
| Chloromethane | < 5 | 5 | |
| 2-Chlorotoluene | < 5 | 5 | |
| 4-Chlorotoluene | < 5 | 5 | |
| 1,2-Dibromo-3-chloropropane | < 1.7 | 1.7 | |
| Dibromochloromethane | < 5 | 5 | |
| 1,2-Dibromoethane (EDB) | < 0.28 | 1 | 1 |
| Dibromomethane | < 5 | 5 | |
| 1,2-Dichlorobenzene | < 5 | 5 | |
| 1,3-Dichlorobenzene | < 5 | 5 | |
| 1,4-Dichlorobenzene | < 5 | 5 | |
| trans-1,4-Dichloro-2-butene | < 5 | 5 | |
| Dichlorodifluoromethane | < 5 | 5 | |
| 1,1-Dichloroethane | < 5 | 5 | |
| 1,2-Dichloroethane | < 5 | 5 | |
| 1,1-Dichloroethene | < 5 | 5 | |
| cis-1,2-Dichloroethene | < 5 | 5 | |
| trans-1,2-Dichloroethene | < 5 | 5 | |
| 1,2-Dichloropropane | < 5 | 5 | |
| 1,3-Dichloropropane | < 5 | 5 | |
| 2,2-Dichloropropane | < 5 | 5 | |
| 1,1-Dichloropropene | < 5 | 5 | |
| 1,3-Dichloropropene | < 5 | 5 | |
| Ethylbenzene | < 5 | 5 | |
| Ethyl methacrylate | < 100 | 100 | |



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8260 QC Continued...

| <u>Method Blank (MB)</u> | <u>MB Results (ug/kg)</u> | <u>Rep Lim (ug/kg)</u> | <u>Flag</u> |
|-----------------------------------|---------------------------|------------------------|-------------|
| Hexachloro-1,3-butadiene | < 5 | 5 | |
| 2-Hexanone | < 10 | 10 | |
| n-Hexane | < 10 | 10 | |
| Iodomethane | < 10 | 10 | |
| Isopropylbenzene (Cumene) | < 5 | 5 | |
| p-Isopropyltoluene | < 5 | 5 | |
| Methylene chloride | < 20 | 20 | |
| 4-Methyl-2-pentanone (MIBK) | < 10 | 10 | |
| Methyl-tert-butyl-ether | < 5 | 5 | |
| 1-Methylnaphthalene | < 5 | 5 | |
| 2-Methylnaphthalene | < 5 | 5 | |
| Naphthalene | < 5 | 5 | |
| n-Propylbenzene | < 5 | 5 | |
| Styrene | < 5 | 5 | |
| 1,1,1,2-Tetrachloroethane | < 5 | 5 | |
| 1,1,2,2-Tetrachloroethane | < 5 | 5 | |
| Tetrachloroethene | < 5 | 5 | |
| Toluene | < 5 | 5 | |
| 1,2,3-Trichlorobenzene | < 5 | 5 | |
| 1,2,4-Trichlorobenzene | < 5 | 5 | |
| 1,1,1-Trichloroethane | < 5 | 5 | |
| 1,1,2-Trichloroethane | < 5 | 5 | |
| Trichloroethene | < 5 | 5 | |
| Trichlorofluoromethane | < 5 | 5 | |
| 1,2,3-Trichloropropane | < 5 | 5 | |
| 1,2,4-Trimethylbenzene | < 5 | 5 | |
| 1,3,5-Trimethylbenzene | < 5 | 5 | |
| Vinyl acetate | < 10 | 10 | |
| Vinyl chloride | < 2 | 2 | |
| Xylene, M&P | < 5 | 5 | |
| Xylene, Ortho | < 5 | 5 | |
| Xylenes, Total | < 10 | 10 | |
| Dibromofluoromethane (surrogate) | 93% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 92% | | |
| Toluene-d8 (surrogate) | 98% | | |
| 4-bromofluorobenzene (surrogate) | 92% | | |
| Analysis Date/Time: | 6-23-23/19:43 | | |
| Analyst Initials | tjg | | |



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8260 QC Continued...

| <u>LCS/LCSD:</u> | <u>LCS Results (ug/kg)</u> | <u>LCS/LCSD Conc. (ug/kg)</u> | <u>LCSD Result (ug/kg)</u> | <u>LCS Rec.</u> | <u>LCSD Rec.</u> | <u>% D</u> | <u>Flag</u> |
|-----------------------------------|----------------------------|-----------------------------------|--------------------------------|-----------------|----------------------|------------|-------------|
| Vinyl Chloride | 50.5 | 50 | 51.9 | 101% | 104% | 2.7 | |
| 1,1-Dichloroethene | 47.5 | 50 | 47.0 | 95% | 94% | 1.1 | |
| trans-1,2-Dichloroethene | 51.4 | 50 | 48.9 | 103% | 98% | 5.0 | |
| Methyl-tert-butyl ether | 47.0 | 50 | 47.5 | 94% | 95% | 1.1 | |
| 1,1-Dichloroethane | 47.1 | 50 | 47.1 | 94% | 94% | 0.0 | |
| cis-1,2-Dichloroethene | 44.4 | 50 | 49.6 | 89% | 99% | 11.1 | |
| Chloroform | 47.3 | 50 | 51.6 | 95% | 103% | 8.7 | |
| 1,1,1-Trichloroethane | 48.4 | 50 | 53.2 | 97% | 106% | 9.4 | |
| Benzene | 48.3 | 50 | 52.1 | 97% | 104% | 7.6 | |
| Trichloroethene | 52.1 | 50 | 56.3 | 104% | 113% | 7.7 | |
| Toluene | 53.3 | 50 | 51.5 | 107% | 103% | 3.4 | |
| 1,1,1,2-Tetrachloroethane | 51.2 | 50 | 55.6 | 102% | 111% | 8.2 | |
| Chlorobenzene | 46.5 | 50 | 50.9 | 93% | 102% | 9.0 | |
| Ethylbenzene | 44.8 | 50 | 49.0 | 90% | 98% | 9.0 | |
| o-Xylene | 46.9 | 50 | 51.9 | 94% | 104% | 10.1 | |
| n-Propylbenzene | 46.5 | 50 | 52.0 | 93% | 104% | 11.2 | |
| Dibromofluoromethane (surrogate) | 100% | | 98% | | | | |
| 1,2-Dichloroethane-d4 (surrogate) | 106% | | 94% | | | | |
| Toluene-d8 (surrogate) | 112% | | 110% | | | | |
| 4-bromofluorobenzene (surrogate) | 95% | | 95% | | | | |
| Analysis Date/Time: | 6-23-23/18:55 | | 6-23-23/19:11 | | | | |
| Analyst Initials | tjg | | tjg | | | | |



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EPA 8270 Quality Control Data

ENVision Batch Number: 062623BS

| <u>Method Blank (MB):</u> | <u>Method Blank Results (mg/kg)</u> | <u>Reporting Limit (mg/kg)</u> | <u>Flag</u> |
|-----------------------------|-------------------------------------|--------------------------------|-------------|
| Acenaphthene | < 0.33 | 0.33 | |
| Acenaphthylene | < 0.33 | 0.33 | |
| Aniline | < 0.33 | 0.33 | |
| Anthracene | < 0.33 | 0.33 | |
| Benzo(a)anthracene | < 0.33 | 0.33 | |
| Benzo(a)pyrene | < 0.33 | 0.33 | |
| Benzo(b)fluoranthene | < 0.33 | 0.33 | |
| Benzo(g,h,i)perylene | < 0.33 | 0.33 | |
| Benzo(k)fluoranthene | < 0.33 | 0.33 | |
| Benzoic Acid | < 1.6 | 1.6 | |
| Benzyl Alcohol | < 0.66 | 0.66 | |
| 4-Bromophenylphenyl ether | < 0.33 | 0.33 | |
| Butylbenzylphthalate | < 0.33 | 0.33 | |
| Carbazole | < 0.66 | 0.66 | |
| 4-Chloro-3-methylphenol | < 0.66 | 0.66 | |
| 4-Chloroaniline | < 0.66 | 0.66 | |
| bis(2-Chloroethoxy)methane | < 0.33 | 0.33 | |
| bis(2-Chloroethyl)ether | < 0.33 | 0.33 | |
| bis(2-Chloroisopropyl)ether | < 0.33 | 0.33 | |
| 2-Chloronaphthalene | < 0.33 | 0.33 | |
| 2-Chlorophenol | < 0.33 | 0.33 | |
| 4-Chlorophenylphenyl ether | < 0.33 | 0.33 | |
| Chrysene | < 0.33 | 0.33 | |
| Dibenzo(a,h)anthracene | < 0.33 | 0.33 | |
| Dibenzofuran | < 0.33 | 0.33 | |
| 1,2-Dichlorobenzene | < 0.33 | 0.33 | |
| 1,3-Dichlorobenzene | < 0.33 | 0.33 | |
| 1,4-Dichlorobenzene | < 0.33 | 0.33 | |
| 3,3-Dichlorobenzidine | < 0.66 | 0.66 | |
| 2,4-Dichlorophenol | < 0.33 | 0.33 | |
| Diethylphthalate | < 0.33 | 0.33 | |
| 2,4-Dimethylphenol | < 0.33 | 0.33 | |
| Dimethylphthalate | < 0.33 | 0.33 | |
| Di-n-butylphthalate | < 0.33 | 0.33 | |
| 4,6-Dinitro-2-methylphenol | < 1.6 | 1.6 | |
| 2,4-Dinitrophenol | < 1.6 | 1.6 | |
| 2,4-Dinitrotoluene | < 0.33 | 0.33 | |
| 2,6-Dinitrotoluene | < 0.33 | 0.33 | |
| Di-n-octylphthalate | < 0.33 | 0.33 | |
| bis(2-Ethylhexyl)phthalate | < 0.33 | 0.33 | |
| Fluoranthene | < 0.33 | 0.33 | |
| Fluorene | < 0.33 | 0.33 | |
| Hexachloro-1,3-butadiene | < 0.33 | 0.33 | |
| Hexachlorobenzene | < 0.33 | 0.33 | |



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8270 QC continued...

| <u>Method Blank (MB):</u> | <u>Method Blank Results (mg/kg)</u> | <u>Reporting Limit (mg/kg)</u> | <u>Flag</u> |
|----------------------------------|-----------------------------------------|------------------------------------|-------------|
| Hexachlorocyclopentadiene | < 0.33 | 0.33 | |
| Hexachloroethane | < 0.33 | 0.33 | |
| Indeno(1,2,3-cd)pyrene | < 0.33 | 0.33 | |
| Isophorone | < 0.33 | 0.33 | |
| 2-Methylnaphthalene | < 0.33 | 0.33 | |
| 2-Methylphenol (o-Cresol) | < 0.33 | 0.33 | |
| 3&4-Methylphenol | < 0.66 | 0.66 | |
| Naphthalene | < 0.33 | 0.33 | |
| 2-Nitroaniline | < 1.6 | 1.6 | |
| 3-Nitroaniline | < 1.6 | 1.6 | |
| 4-Nitroaniline | < 1.6 | 1.6 | |
| Nitrobenzene | < 0.33 | 0.33 | |
| 2-Nitrophenol | < 0.33 | 0.33 | |
| 4-Nitrophenol | < 1.6 | 1.6 | |
| N-Nitroso-di-n-propylamine | < 0.33 | 0.33 | |
| N-Nitrosodiphenylamine | < 0.33 | 0.33 | |
| Pentachlorophenol | < 1.6 | 1.6 | |
| Phenanthren | <0.3 | 0.3 | |
| Phenol | < 0.33 | 0.33 | |
| Pyrene | < 0.33 | 0.33 | |
| 1,2,4-Trichlorobenzene | < 0.33 | 0.33 | |
| 2,4,5-Trichlorophenol | < 0.33 | 0.33 | |
| 2,4,6-Trichlorophenol | < 0.33 | 0.33 | |
| 2-Fluorophenol (surrogate) | 31% | | |
| Phenol-d6 (surrogate) | 28% | | |
| Nitrobenzene-d5 (surrogate) | 43% | | |
| 2-Fluorobiphenyl (surrogate) | 34% | | |
| 2,4,6-Tribromophenol (surrogate) | 32% | | |
| p-Terphenyl-d14 (surrogate) | 32% | | |
| Analysis Date/Time: | 06-27-23/00:27 | | |
| Analyst Initials: | gjd | | |
| Date Extracted: | 6/26/2023 | | |
| Initial Sample Weight: | 30 g | | |
| Final Volume: | 1.0 mL | | |



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| <u>LCS/LCSD:</u> | <u>LCS Results</u> | <u>LCS Concentration</u> | <u>LCSD Results</u> | <u>LCS Recovery</u> | <u>LCSD Recovery</u> | <u>RPD</u> | <u>Flag</u> |
|----------------------------------|--------------------|--------------------------|---------------------|---------------------|----------------------|------------|-------------|
| Phenol | 20.51 | 50.0 | 22.3 | 41% | 45% | 8.5% | |
| 2-Chlorophenol | 20.95 | 50.0 | 23.6 | 42% | 47% | 11.8% | |
| 1,4-Dichlorobenzene | 20.49 | 50.0 | 20.0 | 41% | 40% | 2.3% | |
| N-Nitroso-di-n-propylamine | 20.67 | 50.0 | 21.9 | 41% | 44% | 5.9% | |
| 1,2,4-Trichlorobenzene | 20.50 | 50.0 | 22.0 | 41% | 44% | 7.1% | |
| 4-Chloro-3-methylphenol | 21.15 | 50.0 | 21.7 | 42% | 43% | 2.7% | |
| 2,4,5-Trichlorophenol | 20.83 | 50.0 | 21.3 | 42% | 43% | 2.0% | |
| 2-Nitroaniline | 21.87 | 50.0 | 22.4 | 44% | 45% | 2.3% | |
| 3-Nitroaniline | 22.07 | 50.0 | 20.8 | 44% | 42% | 6.0% | |
| Acenaphthene | 20.96 | 50.0 | 22.8 | 42% | 46% | 8.4% | |
| 4-Nitrophenol | 20.60 | 100.0 | 21.0 | 21% | 21% | 1.7% | |
| 2,4-Dinitrotoluene | 23.97 | 50.0 | 21.1 | 48% | 42% | 13.0% | |
| 4-Nitroaniline | 21.20 | 50.0 | 20.8 | 42% | 42% | 1.9% | |
| 4,6-Dinitro-2-methylphenol | 23.51 | 50.0 | 25.3 | 47% | 51% | 7.2% | |
| Pentachlorophenol | 21.44 | 50.0 | 22.7 | 43% | 45% | 5.7% | |
| Pyrene | 20.88 | 50.0 | 23.5 | 42% | 47% | 11.8% | |
| 2-Fluorophenol (surrogate) | 35% | | 42% | | | | |
| Phenol-d6 (surrogate) | 33% | | 42% | | | | |
| Nitrobenzene-d5 (surrogate) | 26% | | 29% | | | | |
| 2-Fluorobiphenyl (surrogate) | 31% | | 39% | | | | |
| 2,4,6-Tribromophenol (surrogate) | 35% | | 42% | | | | |
| p-Terphenyl-d14 (surrogate) | 27% | | 38% | | | | |
| Analysis Date/Time: | 06-27-23/00:53 | | 06-27-23/01:18 | | | | |
| Analyst Initials: | | gjd | | gjd | | | |
| Date Extracted: | 6/26/2023 | | 6/26/2023 | | | | |
| Initial Sample Weight: | 30 g | | 30 g | | | | |
| Final Volume: | 1.0 mL | | 1.0 mL | | | | |



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EPA 6010B/7471A Metals Quality Control Data

ENVision Batch Number: 062723icp/062823hg

| <u>Method Blank (MB):</u> | <u>MB Results (mg/kg)</u> | <u>Rep Lim (mg/kg)</u> | <u>Flag</u> |
|---------------------------|---------------------------|------------------------|-------------|
| Arsenic | < 2 | 2 | |
| Barium | < 2 | 2 | |
| Cadmium | < 2 | 2 | |
| Chromium | < 2 | 2 | |
| Lead | < 2 | 2 | |
| Mercury | < 1 | 1 | |
| Selenium | < 2 | 2 | |
| Silver | < 2 | 2 | |

Analysis Date/Time: 6-27-23/10:26icp/6/28/23/11:17hg

Analyst Initials: gjd

| <u>Laboratory Control Standard:</u> | <u>LCS Results(ppm)</u> | <u>LCS Conc(ppm)</u> | <u>% Rec</u> | <u>Flag</u> |
|-------------------------------------|-------------------------|----------------------|--------------|-------------|
| Arsenic | 0.54 | 0.50 | 108% | |
| Barium | 0.52 | 0.50 | 104% | |
| Cadmium | 0.53 | 0.50 | 106% | |
| Chromium | 0.49 | 0.50 | 98% | |
| Lead | 0.53 | 0.50 | 106% | |
| Mercury | 0.0056 | 0.005 | 112% | |
| Selenium | 0.52 | 0.50 | 104% | |
| Silver | 0.51 | 0.50 | 102% | |

Analysis Date/Time: 6-27-23/10:24icp/06/28/23/11:15hg

Analyst Initials: gjd



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EPA 8260 Quality Control Data

ENVision Batch Number: 062523VS

| <u>Method Blank (MB):</u> | <u>MB Results (ug/L)</u> | <u>Rep Lim (ug/L)</u> | <u>Flag</u> |
|-----------------------------|--------------------------|-----------------------|-------------|
| Acetone | < 100 | 100 | |
| Acrolein | < 1 | 1 | |
| Acrylonitrile | < 0.45 | 1 | 1 |
| Benzene | < 5 | 5 | |
| Bromobenzene | < 5 | 5 | |
| Bromochloromethane | < 5 | 5 | |
| Bromodichloromethane | < 5 | 5 | |
| Bromoform | < 5 | 5 | |
| Bromomethane | < 5 | 5 | |
| n-Butanol | < 50 | 50 | |
| 2-Butanone (MEK) | < 10 | 10 | |
| n-Butylbenzene | < 5 | 5 | |
| sec-Butylbenzene | < 5 | 5 | |
| tert-Butylbenzene | < 5 | 5 | |
| Carbon Disulfide | < 5 | 5 | |
| Carbon Tetrachloride | < 5 | 5 | |
| Chlorobenzene | < 5 | 5 | |
| Chloroethane | < 5 | 5 | |
| 2-Chloroethylvinylether | < 50 | 50 | |
| Chloroform | < 5 | 5 | |
| Chloromethane | < 5 | 5 | |
| 2-Chlorotoluene | < 5 | 5 | |
| 4-Chlorotoluene | < 5 | 5 | |
| 1,2-Dibromo-3-chloropropane | < 1 | 1 | |
| Dibromochloromethane | < 5 | 5 | |
| 1,2-Dibromoethane (EDB) | < 1 | 1 | |
| Dibromomethane | < 5 | 5 | |
| 1,2-Dichlorobenzene | < 5 | 5 | |
| 1,3-Dichlorobenzene | < 5 | 5 | |
| 1,4-Dichlorobenzene | < 5 | 5 | |
| trans-1,4-Dichloro-2-butene | < 1 | 1 | |
| Dichlorodifluoromethane | < 5 | 5 | |
| 1,1-Dichloroethane | < 5 | 5 | |
| 1,2-Dichloroethane | < 5 | 5 | |
| 1,1-Dichloroethene | < 5 | 5 | |
| cis-1,2-Dichloroethene | < 5 | 5 | |
| trans-1,2-Dichloroethene | < 5 | 5 | |
| 1,2-Dichloropropane | < 5 | 5 | |
| 1,3-Dichloropropane | < 5 | 5 | |
| 2,2-Dichloropropane | < 5 | 5 | |
| 1,1-Dichloropropene | < 5 | 5 | |
| 1,3-Dichloropropene | < 4.1 | 4.1 | |
| Ethylbenzene | < 5 | 5 | |
| Ethyl methacrylate | < 100 | 100 | |



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8260 QC Continued...

| Method Blank (MB): | MB Results (ug/L) | Rep Lim (ug/L) | Flag |
|-----------------------------------------------|--------------------------|-----------------------|-------------|
| Hexachloro-1,3-butadiene | < 2.6 | 2.6 | |
| 2-Hexanone | < 10 | 10 | |
| n-Hexane | < 10 | 10 | |
| Iodomethane | < 10 | 10 | |
| Isopropylbenzene (Cumene) | < 5 | 5 | |
| p-Isopropyltoluene | < 5 | 5 | |
| Methylene chloride | < 5 | 5 | |
| 4-Methyl-2-pentanone (MIBK) | < 10 | 10 | |
| Methyl-tert-butyl-ether | < 5 | 5 | |
| 1-Methylnaphthalene | < 5 | 5 | |
| 2-Methylnaphthalene | < 5 | 5 | |
| Naphthalene | < 1 | 1 | |
| n-Propylbenzene | < 5 | 5 | |
| Styrene | < 5 | 5 | |
| 1,1,1,2-Tetrachloroethane | < 5 | 5 | |
| 1,1,2,2-Tetrachloroethane | < 0.66 | 1 | 1 |
| Tetrachloroethene | < 5 | 5 | |
| Toluene | < 5 | 5 | |
| 1,2,3-Trichlorobenzene | < 5 | 5 | |
| 1,2,4-Trichlorobenzene | < 5 | 5 | |
| 1,1,1-Trichloroethane | < 5 | 5 | |
| 1,1,2-Trichloroethane | < 5 | 5 | |
| Trichloroethene | < 5 | 5 | |
| Trichlorofluoromethane | < 5 | 5 | |
| 1,2,3-Trichloropropane | < 1 | 1 | |
| 1,2,4-Trimethylbenzene | < 5 | 5 | |
| 1,3,5-Trimethylbenzene | < 5 | 5 | |
| Vinyl acetate | < 10 | 10 | |
| Vinyl chloride | < 2 | 2 | |
| Xylene, M&P | < 5 | 5 | |
| Xylene, Ortho | < 5 | 5 | |
| Xylene (total) | < 10 | 10 | |
| Dibromo ^f luoromethane (surrogate) | 95% | | |
| 1,2-Dichloroethane-d4 (surrogate) | 95% | | |
| Toluene-d8 (surrogate) | 107% | | |
| 4-bromofluorobenzene (surrogate) | 97% | | |
| Analysis Date/Time: | 6-25-23/06:06 | | |
| Analyst Initials | tjg | | |



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8260 QC Continued...

| <u>LCS/LCSD</u> | <u>LCS Results (ug/L)</u> | <u>LCS/LCSD Conc. (ug/L)</u> | <u>LCSD Result (ug/L)</u> | <u>LCS Rec.</u> | <u>LCSD Rec.</u> | <u>% D</u> | <u>Flag</u> |
|-----------------------------------|---------------------------|----------------------------------|-------------------------------|-----------------|------------------|------------|-------------|
| Vinyl Chloride | 47.9 | 50 | 52.9 | 96% | 106% | 9.9 | |
| 1,1-Dichloroethene | 50.5 | 50 | 46.6 | 101% | 93% | 8.0 | |
| trans-1,2-Dichloroethene | 48.6 | 50 | 46.4 | 97% | 93% | 4.6 | |
| Methyl-tert-butyl-ether | 49.7 | 50 | 47.6 | 99% | 95% | 4.3 | |
| 1,1-Dichloroethane | 46.5 | 50 | 45.3 | 93% | 91% | 2.6 | |
| cis-1,2-Dichloroethene | 50.6 | 50 | 48.8 | 101% | 98% | 3.6 | |
| Chloroform | 50.4 | 50 | 49.2 | 101% | 98% | 2.4 | |
| 1,1,1-Trichloroethane | 51.3 | 50 | 50.6 | 103% | 101% | 1.4 | |
| Benzene | 53.1 | 50 | 50.8 | 106% | 102% | 4.4 | |
| Trichloroethene | 54.7 | 50 | 53.3 | 109% | 107% | 2.6 | |
| Toluene | 51.3 | 50 | 56.6 | 103% | 113% | 9.8 | |
| 1,1,1,2-Tetrachloroethane | 58.4 | 50 | 51.5 | 117% | 103% | 12.6 | |
| Chlorobenzene | 52.9 | 50 | 46.1 | 106% | 92% | 13.7 | |
| Ethylbenzene | 50.9 | 50 | 44.3 | 102% | 89% | 13.9 | |
| o-Xylene | 53.8 | 50 | 46.7 | 108% | 93% | 14.1 | |
| n-Propylbenzene | 55.2 | 50 | 46.6 | 110% | 93% | 16.9 | |
| Dibromofluoromethane (surrogate) | 93% | | 102% | | | | |
| 1,2-Dichloroethane-d4 (surrogate) | 96% | | 98% | | | | |
| Toluene-d8 (surrogate) | 106% | | 116% | | | | |
| 4-bromofluorobenzene (surrogate) | 100% | | 94% | | | | |
| Analysis Date/Time: | 6-25-23/05:03 | | 6-25-23/05:35 | | | | |
| Analyst Initials | tjg | | tjg | | | | |



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EPA 8270 Quality Control Data

ENVision Batch Number: 062823BW

| <u>BNA Method Blank (MB):</u> | <u>Method Blank</u> <u>Result (ug/L)</u> | <u>Reporting Limit</u> <u>(ug/L)</u> | <u>Flag</u> |
|-------------------------------|---------------------------------------------|-----------------------------------------|-------------|
| Aniline | < 10 | 10 | |
| Benzoic Acid | < 50 | 50 | |
| Benzyl Alcohol | < 20 | 20 | |
| 4-Bromophenylphenyl ether | < 10 | 10 | |
| Butylbenzylphthalate | < 10 | 10 | |
| Carbazole | < 20 | 20 | |
| 4-Chloro-3-methylphenol | < 20 | 20 | |
| 4-Chloroaniline | < 20 | 20 | |
| bis(2-Chloroethoxy)methane | < 10 | 10 | |
| bis(2-Chloroethyl)ether | < 10 | 10 | |
| bis(2-Chloroisopropyl)ether | < 10 | 10 | |
| 2-Chloronaphthalene | < 10 | 10 | |
| 2-Chlorophenol | < 10 | 10 | |
| 4-Chlorophenylphenyl ether | < 10 | 10 | |
| Dibenzofuran | < 10 | 10 | |
| 1,2-Dichlorobenzene | < 10 | 10 | |
| 1,3-Dichlorobenzene | < 10 | 10 | |
| 1,4-Dichlorobenzene | < 10 | 10 | |
| 3,3-Dichlorobenzidine | < 20 | 20 | |
| 2,4-Dichlorophenol | < 10 | 10 | |
| Diethylphthalate | < 10 | 10 | |
| 2,4-Dimethylphenol | < 20 | 20 | |
| Dimethylphthalate | < 10 | 10 | |
| Di-n-butylphthalate | < 10 | 10 | |
| 4,6-Dinitro-2-methylphenol | < 50 | 50 | |
| 2,4-Dinitrophenol | < 50 | 50 | |
| 2,4-Dinitrotoluene | < 10 | 10 | |
| 2,6-Dinitrotoluene | < 10 | 10 | |
| Di-n-octylphthalate | < 10 | 10 | |
| bis(2-Ethylhexyl)phthalate | < 5 | 5 | |
| Hexachloro-1,3-butadiene | < 10 | 10 | |
| Hexachlorobenzene | < 5 | 5 | |
| Hexachlorocyclopentadiene | < 25 | 25 | |
| Hexachloroethane | < 10 | 10 | |
| Isophorone | < 10 | 10 | |
| 2-Methylphenol (o-Cresol) | < 10 | 10 | |
| 3&4-Methylphenol | < 20 | 20 | |
| 2-Nitroaniline | < 50 | 50 | |
| 3-Nitroaniline | < 50 | 50 | |
| 4-Nitroaniline | < 50 | 50 | |
| Nitrobenzene | < 10 | 10 | |
| 2-Nitrophenol | < 10 | 10 | |
| 4-Nitrophenol | < 50 | 50 | |
| N-Nitroso-di-n-propylamine | < 10 | 10 | |
| N-Nitrosodiphenylamine | < 10 | 10 | |



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8270 QC Continued...

| <u>BNA Method Blank (MB):</u> | <u>Method Blank</u> | <u>Reporting Limit</u> | <u>Flag</u> |
|----------------------------------|----------------------|------------------------|-------------|
| | <u>Result (ug/L)</u> | <u>(ug/L)</u> | |
| Pentachlorophenol | < 50 | 50 | |
| Phenol | < 10 | 10 | |
| 1,2,4-Trichlorobenzene | < 10 | 10 | |
| 2,4,5-Trichlorophenol | < 10 | 10 | |
| 2,4,6-Trichlorophenol | < 10 | 10 | |
| 2-Fluorophenol (surrogate) | 90% | | |
| Phenol-d6 (surrogate) | 49% | | |
| Nitrobenzene-d5 (surrogate) | 80% | | |
| 2-Fluorobiphenyl (surrogate) | 77% | | |
| 2,4,6-Tribromophenol (surrogate) | 93% | | |
| p-Terphenyl-d14 (surrogate) | 87% | | |
| Analysis Date/Time: | 06-28-23/17:57 | | |
| Analyst Initials: | JAK | | |
| Date Extracted: | 6/28/2023 | | |
| Initial Sample Volume: | 1000 mL | | |
| Final Volume: | 1.0 mL | | |

| <u>PAH-SIM Method Blank (MB):</u> | <u>Method Blank</u> | <u>Reporting Limit</u> | <u>Flag</u> |
|-----------------------------------|----------------------|------------------------|-------------|
| | <u>Result (ug/L)</u> | <u>(ug/L)</u> | |
| Acenaphthene | < 1.0 | 1.0 | |
| Acenaphthylene | < 1.0 | 1.0 | |
| Anthracene | < 0.10 | 0.10 | |
| Benzo(a)anthracene | < 0.10 | 0.10 | |
| Benzo(a)pyrene | < 0.10 | 0.10 | |
| Benzo(b)fluoranthene | < 0.10 | 0.10 | |
| Benzo(g,h,i)perylene | < 0.10 | 0.10 | |
| Benzo(k)fluoranthene | < 0.10 | 0.10 | |
| Chrysene | < 0.10 | 0.10 | |
| Dibenzo(a,h)anthracene | < 0.10 | 0.10 | |
| Fluoranthene | < 1.0 | 1.0 | |
| Fluorene | < 1.0 | 1.0 | |
| Indeno(1,2,3-cd)pyrene | < 0.022 | 0.022 | |
| 2-methylnaphthalene | < 1.0 | 1.0 | |
| Naphthalene | < 1.0 | 1.0 | |
| Phenanthrene | < 1.0 | 1.0 | |
| Pyrene | < 1.0 | 1.0 | |
| Analysis Date/Time: | 06-23-23/21:52 | | |
| Analyst Initials | JAK | | |



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8270 QC continued...

| <u>LCS/LCSD:</u> | <u>LCS Result (ug/L)</u> | <u>Conc. (ug/L)</u> | <u>LCSD Result (ug/L)</u> | <u>LCS Recovery</u> | <u>LCSD Recovery</u> | <u>RPD</u> | <u>Flag</u> |
|----------------------------------|--------------------------|---------------------|---------------------------|---------------------|----------------------|------------|-------------|
| Acenaphthene | 20.53 | 50.00 | 21.35 | 41.1% | 42.7% | 3.9% | |
| 4-Chloro-3-methylphenol | 24.27 | 50.00 | 23.89 | 48.5% | 47.8% | 1.6% | |
| 2-Chlorophenol | 25.96 | 50.00 | 26.70 | 51.9% | 53.4% | 2.8% | |
| 1,4-Dichlorobenzene | 21.88 | 50.00 | 21.21 | 43.8% | 42.4% | 3.1% | |
| 4,6-Dinitro-2-methylphenol | 20.89 | 50.00 | 21.20 | 41.8% | 42.4% | 1.5% | |
| 2,4-Dinitrotoluene | 23.80 | 50.00 | 25.25 | 47.6% | 50.5% | 5.9% | |
| 2-Nitroaniline | 25.61 | 50.00 | 25.11 | 51.2% | 50.2% | 2.0% | |
| 3-Nitroaniline | 20.23 | 50.00 | 22.54 | 40.5% | 45.1% | 10.8% | |
| 4-Nitroaniline | 24.85 | 50.00 | 26.06 | 49.7% | 52.1% | 4.8% | |
| 4-Nitrophenol | 27.30 | 100.00 | 27.80 | 27.3% | 27.8% | 1.8% | |
| N-Nitroso-di-n-propylamine | 21.38 | 50.00 | 21.80 | 42.8% | 43.6% | 1.9% | |
| Pentachlorophenol | 24.34 | 50.00 | 24.42 | 48.7% | 48.8% | 0.3% | |
| Phenol | 25.57 | 50.00 | 24.36 | 51.1% | 48.7% | 4.8% | |
| Pyrene | 22.57 | 50.00 | 21.06 | 45.1% | 42.1% | 6.9% | |
| 1,2,4-Trichlorobenzene | 23.02 | 50.00 | 22.94 | 46.0% | 45.9% | 0.3% | |
| 2,4,5-Trichlorophenol | 24.20 | 50.00 | 26.85 | 48.4% | 53.7% | 10.4% | |
| 2-Fluorophenol (surrogate) | 88% | | 80% | | | | |
| Phenol-d6 (surrogate) | 30% | | 89% | | | | |
| Nitrobenzene-d5 (surrogate) | 77% | | 76% | | | | |
| 2-Fluorobiphenyl (surrogate) | 71% | | 75% | | | | |
| 2,4,6-Tribromophenol (surrogate) | 93% | | 92% | | | | |
| p-Terphenyl-d14 (surrogate) | 93% | | 91% | | | | |
| Analysis Date/Time: | 06-28-23/18:23 | | 06-28-23/18:49 | | | | |
| Analyst Initials: | JAK | | JAK | | | | |
| Date Extracted: | 6/28/2023 | | 6/28/2023 | | | | |
| Initial Sample Volume: | 1000 mL | | 1000 mL | | | | |
| Final Volume: | 1.0 mL | | 1.0 mL | | | | |



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EPA 6010B/7470A Metals Quality Control Data

ENVision Batch Number: 062723icp/062823hg

| <u>Method Blank (MB):</u> | <u>MB Results (mg/L)</u> | <u>Rep Lim (mg/L)</u> | <u>Flag</u> |
|---------------------------|----------------------------------|-----------------------|-------------|
| Arsenic, dissolved | < 0.01 | 0.01 | |
| Barium, dissolved | < 0.1 | 0.1 | |
| Cadmium, dissolved | < 0.005 | 0.005 | |
| Chromium, dissolved | < 0.01 | 0.01 | |
| Lead, dissolved | < 0.01 | 0.01 | |
| Mercury, dissolved | < 0.002 | 0.002 | |
| Selenium, dissolved | < 0.01 | 0.01 | |
| Silver, dissolved | < 0.05 | 0.05 | |
| Analysis Date/Time: | 6-27-23/10:22icp/6/28/23/11:17hg | | |
| Analyst Initials: | gjd | | |

| <u>Laboratory Control Standard (LCS):</u> | <u>LCS Results(mg/L)</u> | <u>LCS Conc(mg/L)</u> | <u>% Rec</u> | <u>Flag</u> |
|-------------------------------------------|----------------------------------|-----------------------|--------------|-------------|
| Arsenic, dissolved | 0.48 | 0.50 | 96 | |
| Barium, dissolved | 0.49 | 0.50 | 98 | |
| Cadmium, dissolved | 0.49 | 0.50 | 98 | |
| Chromium, dissolved | 0.51 | 0.50 | 102 | |
| Lead, dissolved | 0.53 | 0.50 | 106 | |
| Mercury, dissolved | 0.0056 | 0.005 | 112 | |
| Selenium, dissolved | 0.48 | 0.50 | 96 | |
| Silver, dissolved | 0.48 | 0.50 | 96 | |
| Analysis Date/Time: | 6-27-23/10:20icp/6/28/23/11:15hg | | | |
| Analyst Initials: | gjd | | | |



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Flag Number

- | | <u>Comments</u> |
|---|----------------------------------------------------------------|
| 1 | Reported value is below the reporting limit but above the MDL. |
| 2 | Reported value is from a 50x dilution. TJG 6/29/23 |



CHAIN OF CUSTODY RECORD

ENVision Laboratories, Inc. | 1439 Sadlier Circle West Drive | Indianapolis, IN 46239 | Phone: (317) 351-8632 | Fax: (317) 351-8639

| Client: AME | | Invoice Address: Same | | REQUESTED PARAMETERS | | | | | | Sample Integrity: | | | | | | | |
|---------------------------------------------------------------------------|------------|--------------------------------------------------------------|-------------------|---------------------------------------------------------------|-----|------------------|--------------------------------|------|-------|-------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|--------------------|----------|--|--|--|--|--|
| Report 1302 N McRidian Address: St. Indianapolis, IN | | Project Name: JX1600 | | VOCs SVOCs + CR + G VOCs + PCP + PCB + I.M VOCs + CRAS + Diss | | | | | | Cooler Temp: 2 °C (Circle) Samples on Ice? Yes No Samples Intact? Yes No Custody Seal: Yes No ENVision provided bottles: Yes No VOC vials free of head-space: Yes No N/A pH checked? Yes No N/A Method 5035 collection used? Yes No 5035 samples received within 48 hr of Collection? Yes No | | | | | | | |
| Report To: Tyler Z | | Lab Contact: Cheryl Crum | | | | | | | | Please indicate number of containers per preservative below | | | | | | | |
| Phone: | | Sampled by: A HICKS | | | | | | | | | | | | | | | |
| Fax: | | P.O. Number: | | | | | | | | | | | | | | | |
| Desired TAT: (Please Circle One) 1-day 2-day 3-day Std (5-7 bus. days) | | QA/QC Required: (circle if applicable) Level III Level IV | | | | | | | | | | | | | | | |
| Sample ID | Coll. Date | Coll. Time | Comp (C) Grab (G) | Matrix | HCl | HNO ₃ | H ₂ SO ₄ | NaOH | Other | None | ENVision Sample ID | | | | | | |
| SB-1-GW-6-16 | 6/21 | 1630 | G | WT | X | X | X | | | | X | 23-12287 | | | | | |
| SB-2-GW-2-12 | | 1725 | | WT | X | X | X | | | | X | 12288 | | | | | |
| SB-3-GW-7-17 | | 1130 | | WT | X | X | X | | X | | | 12289 | | | | | |
| SB-4-GW-6-16 | | 1251 | | WT | X | X | X | | X | | | 12290 | | | | | |
| SB-5-GW-6-14 | | 1350 | | WT | X | X | X | | X | | | 12291 | | | | | |
| SB-6-GW-6-16 | | 1506 | | WT | X | X | X | | | | X | 12292 | | | | | |
| SB-7-1-1.5 | | 1525 | | SL | X | X | X | | | | X | 12293 | | | | | |
| SB-1-2-4 | | 1555 | | SL | X | X | X | | | | X | 12294 | | | | | |
| SB-2-2-4 | | 1705 | | SL | X | X | X | | | | X | 12295 | | | | | |
| SB-3-2-4 | | 1105 | | SL | X | X | X | | | | X | 12296 | | | | | |
| SB-6-2-4 | V | 1430 | V | SL | X | X | X | | | | X | 12297 | | | | | |

Comments:

| Relinquished by: | Date | Time | Received by: | Date | Time |
|------------------|---------|------|--------------|---------|------|
| Jane Rutherford | 6-22-23 | 9:25 | 4th Danforth | 6-22-23 | 9:25 |

5035 CHECK-IN SHEET

Client Name: AUGUST MACK ENV.

ENVision project#: 2023-1274

Cooler Temp: 2°C

Method 5035A used: YES X NO

ENVision provided tared vials w/stir bars & Terra Core T-handles: YES X NO

5035A samples were received within 48 hrs of collection: YES X NO

5035A samples were frozen within 48 hrs of collection by lab: YES X NO

If NO, did client freeze samples? YES NO

5035ATable A.1 Reference:

Sample is extruded into an empty sealed vial and cooled to 4° ± 2°C for no more than 48 hours then frozen to < -7°C upon laboratory receipt.

Methanol was added to a vial from each sample for Medium-Level dilution within 48 hrs of collection: YES X NO

5035ATable A.1 Reference:

Sample is extruded into an empty sealed vial and cooled to 4° ± 2°C for no more than 48 hours then preserved with methanol upon laboratory receipt.

Performed by/Date: LISA DAULTON 06-22-23