

## Report of Analysis

|                    |                           |                 |                      |
|--------------------|---------------------------|-----------------|----------------------|
| Client:            | Kleinfelder               | Date Collected: | 11/13/24             |
| Project:           | Webster School            | Date Received:  | 11/14/24             |
| Client Sample ID:  | COMP-2                    | SDG No.:        | P4852                |
| Lab Sample ID:     | P4852-02                  | Matrix:         | SOIL                 |
| Analytical Method: | SW8260                    | % Solid:        | 81.4                 |
| Sample Wt/Vol:     | 5.18      Units:    g     | Final Vol:      | 5000              uL |
| Soil Aliquot Vol:  | uL                        | Test:           | VOCMS Group1         |
| GC Column:         | RTX-VMS      ID :    0.18 | Level :         | LOW                  |
| Prep Method :      |                           |                 |                      |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VD080069.D        | 1         |           | 11/18/24 15:18 | VD111824      |

| CAS Number                | Parameter              | Conc. | Qualifier | MDL      | LOQ / CRQL | Units(Dry Weight) |
|---------------------------|------------------------|-------|-----------|----------|------------|-------------------|
| <b>TARGETS</b>            |                        |       |           |          |            |                   |
| 156-59-2                  | cis-1,2-Dichloroethene | 0.72  | U         | 0.72     | 5.90       | ug/Kg             |
| 71-55-6                   | 1,1,1-Trichloroethane  | 0.92  | U         | 0.92     | 5.90       | ug/Kg             |
| 71-43-2                   | Benzene                | 0.85  | U         | 0.85     | 5.90       | ug/Kg             |
| 79-01-6                   | Trichloroethene        | 0.89  | U         | 0.89     | 5.90       | ug/Kg             |
| 108-88-3                  | Toluene                | 0.79  | U         | 0.79     | 5.90       | ug/Kg             |
| 100-41-4                  | Ethyl Benzene          | 0.74  | U         | 0.74     | 5.90       | ug/Kg             |
| 1330-20-7                 | Total Xylenes          | 2.43  | U         | 2.43     | 17.8       | ug/Kg             |
| 98-82-8                   | Isopropylbenzene       | 0.79  | U         | 0.79     | 5.90       | ug/Kg             |
| <b>SURROGATES</b>         |                        |       |           |          |            |                   |
| 17060-07-0                | 1,2-Dichloroethane-d4  | 140   | *         | 50 - 163 | 286%       | SPK: 50           |
| 1868-53-7                 | Dibromofluoromethane   | 87.9  | *         | 54 - 147 | 176%       | SPK: 50           |
| 2037-26-5                 | Toluene-d8             | 41.6  |           | 58 - 134 | 83%        | SPK: 50           |
| 460-00-4                  | 4-Bromofluorobenzene   | 47.8  |           | 29 - 146 | 96%        | SPK: 50           |
| <b>INTERNAL STANDARDS</b> |                        |       |           |          |            |                   |
| 363-72-4                  | Pentafluorobenzene     | 10600 | 7.881     |          |            |                   |
| 540-36-3                  | 1,4-Difluorobenzene    | 22700 | 8.781     |          |            |                   |
| 3114-55-4                 | Chlorobenzene-d5       | 22500 | 11.581    |          |            |                   |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4 | 10100 | 13.516    |          |            |                   |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products