

Report of Analysis

| Client: | CDM Smith | | | | Date Collected: | 05/13/25 | | |
|----------------------|---------------------|----------------|-----------|------------------|--------------------|---------------|-----------|----------------|
| Project: | South River WM | Replacement | | | Date Received: | 05/13/25 | | |
| Client Sample ID: | COMP-1 | | | | SDG No.: | Q2032 | | |
| Lab Sample ID: | Q2032-09 | | | | Matrix: | TCLP | | |
| | | | | | | | | |
| Analytical Method | | | | | % Solid: | 0 | Decanted: | |
| Sample Wt/Vol: | 100 Units: | mL | | | Final Vol: | 10000 | uL | |
| Soil Aliquot Vol: | | uL | | | Test: | TCLP Pestic | cide | |
| Extraction Type: | | | | | Injection Volume : | | | |
| GPC Factor : | 1.0 | PH : | | | - | | | |
| | | 111. | | | | | | |
| Prep Method : | SW3541B | | | | | | | |
| File ID/Qc Batch: | Dilution: Prep Date | | Date | | Date Analyzed | Prep Batch ID | | |
| PD088624.D | 1 | 05/19/25 09:20 | | 05/19/25 22:16 | PB168066 | | | |
| CAS Number | Parameter | Conc. | Qualifier | MDL | | LOQ / C | RQL | Units |
| TARGETS | | | | | | | | |
| 58-89-9 | gamma-BHC (Lindane) | 0.037 | U | 0.037 | | | 0.50 | ug/L |
| 76-44-8 | Heptachlor | 0.027 | U | 0.027 | | | 0.50 | ug/L |
| 1024-57-3 | Heptachlor epoxide | 0.096 | U | 0.096 | | | 0.50 | ug/L |
| 72-20-8 | Endrin | 0.032 | U | 0.032 | | | 0.50 | ug/L |
| 72-43-5 | Methoxychlor | 0.11 | U | 0.11 | | | 0.50 | ug/L |
| | | 1.70 | U | 1.70 | | | 10.0 | ug/L |
| 8001-35-2 | Toxaphene | | | | | | 5.00 | ug/L |
| 8001-35-2 57-74-9 | Chlordane | 0.88 | U | 0.88 | | | 5.00 | ug/L |
| | • | | U | 0.88 | | | 5.00 | ug/L |
| 57-74-9 | • | | U | 0.88 43 - 140 | | | 5.00 | ug/L SPK: 2 |

Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates > 25% difference for detected

concentrations between the two GC columns

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

S = Indicates estimated value where valid five-point calibration

was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit