

Report of Analysis

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|--------------------|------------------------|----------|--------------------|----------|-----------|
| Client: | Nobis Group | | Date Collected: | 06/05/25 | |
| Project: | Raymark Superfund Site | | Date Received: | 06/05/25 | |
| Client Sample ID: | BU-03-060525MSD | | SDG No.: | Q2259 | |
| Lab Sample ID: | Q2246-01MSD | | Matrix: | SOIL | |
| Analytical Method: | 8082A | | % Solid: | 99 | Decanted: |
| Sample Wt/Vol: | 30.05 | Units: g | Final Vol: | 10000 | uL |
| Soil Aliquot Vol: | | uL | Test: | PCB | |
| Extraction Type: | | | Injection Volume : | | |
| GPC Factor : | 1.0 | PH : | | | |
| Prep Method : | SW3541B | | | | |

| | | | | |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed | Prep Batch ID |
| PP072747.D | 1 | 06/09/25 09:10 | 06/09/25 14:18 | PB168349 |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOD | LOQ / CRQL | Units(Dry Weight) |
|-------------------|----------------------|-------|-----------|----------|------|------------|-------------------|
| TARGETS | | | | | | | |
| 12674-11-2 | Aroclor-1016 | 141 | | 4.00 | 8.40 | 17.1 | ug/kg |
| 11104-28-2 | Aroclor-1221 | 13.1 | U | 4.10 | 13.1 | 17.1 | ug/kg |
| 11141-16-5 | Aroclor-1232 | 8.40 | U | 3.80 | 8.40 | 17.1 | ug/kg |
| 53469-21-9 | Aroclor-1242 | 8.40 | U | 4.00 | 8.40 | 17.1 | ug/kg |
| 12672-29-6 | Aroclor-1248 | 13.1 | U | 6.00 | 13.1 | 17.1 | ug/kg |
| 11097-69-1 | Aroclor-1254 | 8.40 | U | 3.20 | 8.40 | 17.1 | ug/kg |
| 37324-23-5 | Aroclor-1262 | 13.1 | U | 5.10 | 13.1 | 17.1 | ug/kg |
| 11100-14-4 | Aroclor-1268 | 8.40 | U | 3.60 | 8.40 | 17.1 | ug/kg |
| 11096-82-5 | Aroclor-1260 | 136 | | 3.30 | 8.40 | 17.1 | ug/kg |
| SURROGATES | | | | | | | |
| 877-09-8 | Tetrachloro-m-xylene | 21.9 | | 44 - 130 | | 110% | SPK: 20 |
| 2051-24-3 | Decachlorobiphenyl | 23.0 | | 60 - 125 | | 115% | SPK: 20 |

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