

## Report of Analysis

|                    |                        |                 |               |
|--------------------|------------------------|-----------------|---------------|
| Client:            | Nobis Group            | Date Collected: | 05/07/25      |
| Project:           | Raymark Superfund Site | Date Received:  | 05/08/25      |
| Client Sample ID:  | OU4-TS-28-050725       | SDG No.:        | Q1984         |
| Lab Sample ID:     | Q1984-15               | Matrix:         | SOIL          |
| Analytical Method: | 8270E                  | % Solid:        | 64.6          |
| Sample Wt/Vol:     | 30.03 Units: g         | Final Vol:      | 1000 uL       |
| Soil Aliquot Vol:  | uL                     | Test:           | SVOCMS Group3 |
| Extraction Type :  | Decanted : N           | Level :         | LOW           |
| Injection Volume : | GPC Factor : 1.0       | GPC Cleanup :   | N PH :        |
| Prep Method :      | SW3541                 |                 |               |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP024646.D        | 1         | 05/13/25 09:05 | 05/15/25 20:45 | PB167973      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOD | LOQ / CRQL | Units(Dry Weight) |
|------------|-----------|-------|-----------|-----|-----|------------|-------------------|
|------------|-----------|-------|-----------|-----|-----|------------|-------------------|

### TARGETS

|          |                        |      |   |       |      |      |       |
|----------|------------------------|------|---|-------|------|------|-------|
| 91-20-3  | Naphthalene            | 0.20 | U | 0.035 | 0.20 | 0.26 | mg/Kg |
| 91-57-6  | 2-Methylnaphthalene    | 0.20 | U | 0.040 | 0.20 | 0.26 | mg/Kg |
| 208-96-8 | Acenaphthylene         | 0.20 | U | 0.045 | 0.20 | 0.26 | mg/Kg |
| 83-32-9  | Acenaphthene           | 0.20 | U | 0.033 | 0.20 | 0.26 | mg/Kg |
| 86-73-7  | Fluorene               | 0.20 | U | 0.039 | 0.20 | 0.26 | mg/Kg |
| 85-01-8  | Phenanthrene           | 0.20 | U | 0.032 | 0.20 | 0.26 | mg/Kg |
| 120-12-7 | Anthracene             | 0.20 | U | 0.052 | 0.20 | 0.26 | mg/Kg |
| 206-44-0 | Fluoranthene           | 0.20 | U | 0.046 | 0.20 | 0.26 | mg/Kg |
| 129-00-0 | Pyrene                 | 0.20 | U | 0.056 | 0.20 | 0.26 | mg/Kg |
| 56-55-3  | Benzo(a)anthracene     | 0.20 | U | 0.036 | 0.20 | 0.26 | mg/Kg |
| 218-01-9 | Chrysene               | 0.20 | U | 0.031 | 0.20 | 0.26 | mg/Kg |
| 205-99-2 | Benzo(b)fluoranthene   | 0.20 | U | 0.029 | 0.20 | 0.26 | mg/Kg |
| 207-08-9 | Benzo(k)fluoranthene   | 0.20 | U | 0.035 | 0.20 | 0.26 | mg/Kg |
| 50-32-8  | Benzo(a)pyrene         | 0.20 | U | 0.046 | 0.20 | 0.26 | mg/Kg |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.20 | U | 0.045 | 0.20 | 0.26 | mg/Kg |
| 53-70-3  | Dibenzo(a,h)anthracene | 0.20 | U | 0.042 | 0.20 | 0.26 | mg/Kg |
| 191-24-2 | Benzo(g,h,i)perylene   | 0.20 | U | 0.040 | 0.20 | 0.26 | mg/Kg |

### SURROGATES

|           |                  |      |   |          |     |          |
|-----------|------------------|------|---|----------|-----|----------|
| 4165-60-0 | Nitrobenzene-d5  | 51.3 |   | 37 - 122 | 51% | SPK: 100 |
| 321-60-8  | 2-Fluorobiphenyl | 44.6 |   | 44 - 115 | 45% | SPK: 100 |
| 1718-51-0 | Terphenyl-d14    | 43.8 | * | 54 - 127 | 44% | SPK: 100 |

### INTERNAL STANDARDS

|            |                        |         |        |
|------------|------------------------|---------|--------|
| 3855-82-1  | 1,4-Dichlorobenzene-d4 | 238000  | 7.658  |
| 1146-65-2  | Naphthalene-d8         | 987000  | 10.428 |
| 15067-26-2 | Acenaphthene-d10       | 607000  | 14.293 |
| 1517-22-2  | Phenanthrene-d10       | 1030000 | 17.092 |
| 1719-03-5  | Chrysene-d12           | 847000  | 21.527 |
| 1520-96-3  | Perylene-d12           | 1080000 | 24.81  |

### TENTATIVE IDENTIFIED COMPOUNDS

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|------------|------------|-------|-----------|-----|-----|------------|-------------------|
| 15972608   | Alachlor   | N.D   |           |     |     |            |                   |
| 82-68-8    | Quintozine | N.D   |           |     |     |            |                   |

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