

## Report of Analysis

|                    |                                    |                 |          |
|--------------------|------------------------------------|-----------------|----------|
| Client:            | Walsh Construction Company II, LLC | Date Collected: | 04/28/25 |
| Project:           | Walsh CO-032 Sampling              | Date Received:  | 04/28/25 |
| Client Sample ID:  | CO-8R-WC                           | SDG No.:        | Q1907    |
| Lab Sample ID:     | Q1907-02                           | Matrix:         | TCLP     |
| Analytical Method: | SW8270                             | % Solid:        | 0        |
| Sample Wt/Vol:     | 100 Units: mL                      | Final Vol:      | 1000 uL  |
| Soil Aliquot Vol:  | uL                                 | Test:           | TCLP BNA |
| 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 |
| BF142282.D        | 1         | 04/30/25 13:15 | 05/02/25 14:51 | PB167810      |

| CAS Number                | Parameter              | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------|------------------------|--------|-----------|----------|------------|----------|
| <b>TARGETS</b>            |                        |        |           |          |            |          |
| 110-86-1                  | Pyridine               | 12.8   | U         | 12.8     | 50.0       | ug/L     |
| 106-46-7                  | 1,4-Dichlorobenzene    | 5.30   | U         | 5.30     | 50.0       | ug/L     |
| 95-48-7                   | 2-Methylphenol         | 11.2   | U         | 11.2     | 50.0       | ug/L     |
| 65794-96-9                | 3+4-Methylphenols      | 11.0   | U         | 11.0     | 100        | ug/L     |
| 67-72-1                   | Hexachloroethane       | 6.50   | U         | 6.50     | 50.0       | ug/L     |
| 98-95-3                   | Nitrobenzene           | 7.60   | U         | 7.60     | 50.0       | ug/L     |
| 87-68-3                   | Hexachlorobutadiene    | 5.40   | U         | 5.40     | 50.0       | ug/L     |
| 88-06-2                   | 2,4,6-Trichlorophenol  | 5.10   | U         | 5.10     | 50.0       | ug/L     |
| 95-95-4                   | 2,4,5-Trichlorophenol  | 6.20   | U         | 6.20     | 50.0       | ug/L     |
| 121-14-2                  | 2,4-Dinitrotoluene     | 12.2   | U         | 12.2     | 50.0       | ug/L     |
| 118-74-1                  | Hexachlorobenzene      | 5.20   | U         | 5.20     | 50.0       | ug/L     |
| 87-86-5                   | Pentachlorophenol      | 15.8   | U         | 15.8     | 100        | ug/L     |
| <b>SURROGATES</b>         |                        |        |           |          |            |          |
| 367-12-4                  | 2-Fluorophenol         | 94.6   |           | 10 - 139 | 63%        | SPK: 150 |
| 13127-88-3                | Phenol-d6              | 33.9   |           | 10 - 134 | 23%        | SPK: 150 |
| 4165-60-0                 | Nitrobenzene-d5        | 91.6   |           | 49 - 133 | 92%        | SPK: 100 |
| 321-60-8                  | 2-Fluorobiphenyl       | 253    | *         | 52 - 132 | 253%       | SPK: 100 |
| 118-79-6                  | 2,4,6-Tribromophenol   | 483    | *         | 44 - 137 | 322%       | SPK: 150 |
| 1718-51-0                 | Terphenyl-d14          | 419    | *         | 48 - 125 | 419%       | SPK: 100 |
| <b>INTERNAL STANDARDS</b> |                        |        |           |          |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4 | 205000 |           | 6.904    |            |          |
| 1146-65-2                 | Naphthalene-d8         | 707000 |           | 8.186    |            |          |
| 15067-26-2                | Acenaphthene-d10       | 110000 |           | 9.939    |            |          |
| 1517-22-2                 | Phenanthrene-d10       | 264000 |           | 11.427   |            |          |
| 1719-03-5                 | Chrysene-d12           | 22800  |           | 14.063   |            |          |
| 1520-96-3                 | Perylene-d12           | 209    |           | 15.557   |            |          |

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|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

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