

## Report of Analysis

|                    |                                     |                 |          |
|--------------------|-------------------------------------|-----------------|----------|
| Client:            | ENTACT                              | Date Collected: | 05/13/25 |
| Project:           | 540 Degraw St, Brooklyn, NY - E9309 | Date Received:  | 05/19/25 |
| Client Sample ID:  | WC-A1-06A-GRE                       | SDG No.:        | Q2078    |
| Lab Sample ID:     | Q2078-09RE                          | Matrix:         | TCLP     |
| Analytical Method: | 8260D                               | % Solid:        | 0        |
| Sample Wt/Vol:     | 5 Units: mL                         | Final Vol:      | 5000 uL  |
| Soil Aliquot Vol:  | uL                                  | Test:           | TCLP VOA |
| GC Column:         | RXI-624 ID : 0.25                   | Level :         | LOW      |
| Prep Method :      | SW5035                              |                 |          |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VN086757.D        | 1         |           | 05/22/25 11:31 | VN052225      |

| CAS Number                | Parameter              | Conc.    | Qualifier | MDL                 | LOQ / CRQL | Units   |
|---------------------------|------------------------|----------|-----------|---------------------|------------|---------|
| <b>TARGETS</b>            |                        |          |           |                     |            |         |
| 75-01-4                   | Vinyl Chloride         | 0.00026  | U         | 0.00026             | 0.0050     | mg/L    |
| 75-35-4                   | 1,1-Dichloroethene     | 0.00023  | U         | 0.00023             | 0.0050     | mg/L    |
| 78-93-3                   | 2-Butanone             | 0.00098  | U         | 0.00098             | 0.025      | mg/L    |
| 56-23-5                   | Carbon Tetrachloride   | 0.00025  | U         | 0.00025             | 0.0050     | mg/L    |
| 67-66-3                   | Chloroform             | 0.00025  | U         | 0.00025             | 0.0050     | mg/L    |
| 71-43-2                   | Benzene                | 0.0073   |           | 0.00015             | 0.0050     | mg/L    |
| 107-06-2                  | 1,2-Dichloroethane     | 0.00022  | U         | 0.00022             | 0.0050     | mg/L    |
| 79-01-6                   | Trichloroethene        | 0.000090 | U         | 0.000090            | 0.0050     | mg/L    |
| 127-18-4                  | Tetrachloroethene      | 0.00023  | U         | 0.00023             | 0.0050     | mg/L    |
| 108-90-7                  | Chlorobenzene          | 0.00012  | U         | 0.00012             | 0.0050     | mg/L    |
| <b>SURROGATES</b>         |                        |          |           |                     |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4  | 43.7     |           | 70 (74) - 130 (125) | 87%        | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane   | 35.2     |           | 70 (75) - 130 (124) | 70%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8             | 50.4     |           | 70 (86) - 130 (113) | 101%       | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene   | 53.3     |           | 70 (77) - 130 (121) | 107%       | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                        |          |           |                     |            |         |
| 363-72-4                  | Pentafluorobenzene     | 236000   | 8.224     |                     |            |         |
| 540-36-3                  | 1,4-Difluorobenzene    | 434000   | 9.1       |                     |            |         |
| 3114-55-4                 | Chlorobenzene-d5       | 408000   | 11.865    |                     |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4 | 196000   | 13.788    |                     |            |         |

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