

Report of Analysis

Client:	ENTACT		Date Collected:	
Project:	540 Degraw St, Brooklyn, NY - E9309		Date Received:	
Client Sample ID:	PB167089BS		SDG No.:	Q1523
Lab Sample ID:	PB167089BS		Matrix:	TCLP
Analytical Method:	SW8270		% Solid:	0
Sample Wt/Vol:	1000	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 :	SW3510C			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BF141933.D	1	03/11/25 11:15	03/12/25 11:06	PB167089

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
110-86-1	Pyridine	41.1		1.30	5.00	ug/Kg
106-46-7	1,4-Dichlorobenzene	45.6		0.53	5.00	ug/Kg
95-48-7	2-Methylphenol	45.8		1.10	5.00	ug/Kg
65794-96-9	3+4-Methylphenols	45.0		1.10	10.0	ug/Kg
67-72-1	Hexachloroethane	45.4		0.65	5.00	ug/Kg
98-95-3	Nitrobenzene	46.1		0.76	5.00	ug/Kg
87-68-3	Hexachlorobutadiene	46.4		0.54	5.00	ug/Kg
88-06-2	2,4,6-Trichlorophenol	45.5		0.51	5.00	ug/Kg
95-95-4	2,4,5-Trichlorophenol	45.6		0.62	5.00	ug/Kg
121-14-2	2,4-Dinitrotoluene	48.3		1.20	5.00	ug/Kg
118-74-1	Hexachlorobenzene	47.3		0.52	5.00	ug/Kg
87-86-5	Pentachlorophenol	97.3	E	1.60	10.0	ug/Kg
SURROGATES						
367-12-4	2-Fluorophenol	138		15 (10) - 110 (139)	92%	SPK: 150
13127-88-3	Phenol-d6	134		15 (10) - 110 (134)	89%	SPK: 150
4165-60-0	Nitrobenzene-d5	99.0		30 (49) - 130 (133)	99%	SPK: 100
321-60-8	2-Fluorobiphenyl	98.1		30 (52) - 130 (132)	98%	SPK: 100
118-79-6	2,4,6-Tribromophenol	147		15 (44) - 110 (137)	98%	SPK: 150
1718-51-0	Terphenyl-d14	96.6		30 (48) - 130 (125)	97%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	148000		6.881		
1146-65-2	Naphthalene-d8	575000		8.163		
15067-26-2	Acenaphthene-d10	321000		9.916		
1517-22-2	Phenanthrene-d10	536000		11.404		
1719-03-5	Chrysene-d12	372000		14.039		
1520-96-3	Perylene-d12	339000		15.509		

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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