

Report of Analysis

Client:	PSEG	Date Collected:	01/13/25
Project:	Hackensack Sub	Date Received:	01/13/25
Client Sample ID:	HACKENSACK-DGA	SDG No.:	Q1074
Lab Sample ID:	Q1074-02	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
BF141148.D	1	01/14/25 10:15	01/14/25 16:27	PB166052

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
110-86-1	Pyridine	0.050	U	0.016	0.050	mg/L
106-46-7	1,4-Dichlorobenzene	0.050	U	0.0084	0.050	mg/L
95-48-7	2-Methylphenol	0.050	U	0.011	0.050	mg/L
65794-96-9	3+4-Methylphenols	0.10	U	0.012	0.10	mg/L
67-72-1	Hexachloroethane	0.050	U	0.010	0.050	mg/L
98-95-3	Nitrobenzene	0.050	U	0.013	0.050	mg/L
87-68-3	Hexachlorobutadiene	0.050	U	0.013	0.050	mg/L
88-06-2	2,4,6-Trichlorophenol	0.050	U	0.0089	0.050	mg/L
95-95-4	2,4,5-Trichlorophenol	0.050	U	0.010	0.050	mg/L
121-14-2	2,4-Dinitrotoluene	0.050	U	0.015	0.050	mg/L
118-74-1	Hexachlorobenzene	0.050	U	0.011	0.050	mg/L
87-86-5	Pentachlorophenol	0.10	U	0.019	0.10	mg/L
SURROGATES						
367-12-4	2-Fluorophenol	146		10 - 139	97%	SPK: 150
13127-88-3	Phenol-d6	141		10 - 134	94%	SPK: 150
4165-60-0	Nitrobenzene-d5	106		49 - 133	106%	SPK: 100
321-60-8	2-Fluorobiphenyl	99.9		52 - 132	100%	SPK: 100
118-79-6	2,4,6-Tribromophenol	177		44 - 137	118%	SPK: 150
1718-51-0	Terphenyl-d14	87.4		48 - 125	87%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	151000	6.822			
1146-65-2	Naphthalene-d8	603000	8.098			
15067-26-2	Acenaphthene-d10	332000	9.851			
1517-22-2	Phenanthrene-d10	583000	11.339			
1719-03-5	Chrysene-d12	465000	13.98			
1520-96-3	Perylene-d12	371000	15.445			

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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