

Report of Analysis

Client:	AECOM		Date Collected:		
Project:	National Grid Equity - Brooklyn NY		Date Received:		
Client Sample ID:	PB169313BSD		SDG No.:	Q2900	
Lab Sample ID:	PB169313BSD		Matrix:	Water	
Analytical Method:	8270E		% Solid:	0	
Sample Wt/Vol:	1000	Units: mL	Final Vol:	1000	uL
Soil Aliquot Vol:		uL	Test:	SVOC-TCL BNA -20	
Extraction Type :		Decanted : N	Level :	LOW	
Injection Volume :		GPC Factor : 1.0	GPC Cleanup :	N	PH :
Prep Method :					

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BF143489.D	1	08/19/25 09:29	08/20/25 17:13	PB169313

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
100-52-7	Benzaldehyde	26.7		3.90	10.0	ug/L
108-95-2	Phenol	43.9		0.91	5.00	ug/L
111-44-4	bis(2-Chloroethyl)ether	46.1		0.81	5.00	ug/L
95-57-8	2-Chlorophenol	46.9		0.58	5.00	ug/L
95-48-7	2-Methylphenol	47.2		1.10	5.00	ug/L
108-60-1	2,2-oxybis(1-Chloropropane)	46.8		1.30	5.00	ug/L
98-86-2	Acetophenone	47.0		0.74	5.00	ug/L
65794-96-9	3+4-Methylphenols	46.3		1.10	10.0	ug/L
621-64-7	n-Nitroso-di-n-propylamine	47.8		1.40	2.50	ug/L
67-72-1	Hexachloroethane	46.8		0.65	5.00	ug/L
98-95-3	Nitrobenzene	46.8		0.76	5.00	ug/L
78-59-1	Isophorone	48.7		0.75	5.00	ug/L
88-75-5	2-Nitrophenol	48.9		1.80	5.00	ug/L
105-67-9	2,4-Dimethylphenol	49.4		1.90	5.00	ug/L
111-91-1	bis(2-Chloroethoxy)methane	46.6		0.68	5.00	ug/L
120-83-2	2,4-Dichlorophenol	45.3		0.52	5.00	ug/L
91-20-3	Naphthalene	44.8		0.50	5.00	ug/L
106-47-8	4-Chloroaniline	12.9		0.84	5.00	ug/L
87-68-3	Hexachlorobutadiene	46.5		0.54	5.00	ug/L
105-60-2	Caprolactam	52.1		1.10	10.0	ug/L
59-50-7	4-Chloro-3-methylphenol	46.1		0.59	5.00	ug/L
91-57-6	2-Methylnaphthalene	41.2		0.56	5.00	ug/L
77-47-4	Hexachlorocyclopentadiene	120	E	3.60	10.0	ug/L
88-06-2	2,4,6-Trichlorophenol	45.7		0.51	5.00	ug/L
95-95-4	2,4,5-Trichlorophenol	44.0		0.62	5.00	ug/L
92-52-4	1,1-Biphenyl	44.9		0.53	5.00	ug/L
91-58-7	2-Chloronaphthalene	44.4		0.61	5.00	ug/L
88-74-4	2-Nitroaniline	47.5		1.30	5.00	ug/L
131-11-3	Dimethylphthalate	47.0		0.61	5.00	ug/L

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208-96-8	Acenaphthylene	50.0		0.75	5.00	ug/L
606-20-2	2,6-Dinitrotoluene	51.2		0.92	5.00	ug/L
99-09-2	3-Nitroaniline	24.3		1.10	5.00	ug/L
83-32-9	Acenaphthene	49.3		0.55	5.00	ug/L
51-28-5	2,4-Dinitrophenol	93.0	E	6.00	10.0	ug/L
100-02-7	4-Nitrophenol	100	E	2.40	10.0	ug/L
132-64-9	Dibenzofuran	44.9		0.61	5.00	ug/L
121-14-2	2,4-Dinitrotoluene	52.3		1.20	5.00	ug/L
84-66-2	Diethylphthalate	48.4		0.69	5.00	ug/L
7005-72-3	4-Chlorophenyl-phenylether	45.5		0.68	5.00	ug/L
86-73-7	Fluorene	45.8		0.63	5.00	ug/L
100-01-6	4-Nitroaniline	48.3		1.50	5.00	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	53.3		2.90	10.0	ug/L
86-30-6	n-Nitrosodiphenylamine	48.3		0.58	5.00	ug/L
101-55-3	4-Bromophenyl-phenylether	46.8		0.40	5.00	ug/L
118-74-1	Hexachlorobenzene	47.1		0.52	5.00	ug/L
1912-24-9	Atrazine	52.6		1.00	5.00	ug/L
87-86-5	Pentachlorophenol	100	E	1.60	10.0	ug/L
85-01-8	Phenanthrene	46.8		0.50	5.00	ug/L
120-12-7	Anthracene	47.1		0.61	5.00	ug/L
86-74-8	Carbazole	47.8		0.72	5.00	ug/L
84-74-2	Di-n-butylphthalate	51.4		1.20	5.00	ug/L
206-44-0	Fluoranthene	46.9		0.82	5.00	ug/L
129-00-0	Pyrene	48.7		0.50	5.00	ug/L
85-68-7	Butylbenzylphthalate	53.8		1.90	5.00	ug/L
91-94-1	3,3-Dichlorobenzidine	20.5		0.93	10.0	ug/L
56-55-3	Benzo(a)anthracene	46.9		0.45	5.00	ug/L
218-01-9	Chrysene	49.3		0.44	5.00	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	50.7		1.60	5.00	ug/L
117-84-0	Di-n-octyl phthalate	51.3		2.30	10.0	ug/L
205-99-2	Benzo(b)fluoranthene	47.0		0.49	5.00	ug/L

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207-08-9	Benzo(k)fluoranthene	49.1		0.48	5.00	ug/L
50-32-8	Benzo(a)pyrene	50.8		0.55	5.00	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	47.4		0.59	5.00	ug/L
53-70-3	Dibenzo(a,h)anthracene	48.4		0.67	5.00	ug/L
191-24-2	Benzo(g,h,i)perylene	49.0		0.69	5.00	ug/L
95-94-3	1,2,4,5-Tetrachlorobenzene	43.6		0.52	5.00	ug/L
123-91-1	1,4-Dioxane	39.4		1.00	5.00	ug/L
58-90-2	2,3,4,6-Tetrachlorophenol	52.0		0.72	5.00	ug/L
SURROGATES						
367-12-4	2-Fluorophenol	118		23 - 138	79%	SPK: 150
13127-88-3	Phenol-d6	120		10 - 134	80%	SPK: 150
4165-60-0	Nitrobenzene-d5	87.2		67 - 132	87%	SPK: 100
321-60-8	2-Fluorobiphenyl	81.4		52 - 132	81%	SPK: 100
118-79-6	2,4,6-Tribromophenol	127		44 - 137	84%	SPK: 150
1718-51-0	Terphenyl-d14	91.2		42 - 152	91%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	134000	6.928			
1146-65-2	Naphthalene-d8	510000	8.21			
15067-26-2	Acenaphthene-d10	281000	9.969			
1517-22-2	Phenanthrene-d10	430000	11.457			
1719-03-5	Chrysene-d12	235000	14.098			
1520-96-3	Perylene-d12	275000	15.598			

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