

Report of Analysis

Client:	R3M Engineering, Inc.		Date Collected:	
Project:	MC054.36 23-8-1 LM - Landing Lane New Brunswick		Date Received:	
Client Sample ID:	PB165432BSD		SDG No.:	P5093
Lab Sample ID:	PB165432BSD		Matrix:	Water
Analytical Method:	SW8270		% 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 :	SW3510C			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BF140786.D	1	12/06/24 09:33	12/09/24 13:39	PB165432

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
100-52-7	Benzaldehyde	13.7		4.00	10.0	ug/L
108-95-2	Phenol	51.0		0.93	5.00	ug/L
111-44-4	bis(2-Chloroethyl)ether	50.5		1.20	5.00	ug/L
95-57-8	2-Chlorophenol	50.9		0.71	5.00	ug/L
95-48-7	2-Methylphenol	50.0		1.10	5.00	ug/L
108-60-1	2,2-oxybis(1-Chloropropane)	49.7		1.40	5.00	ug/L
98-86-2	Acetophenone	48.2		1.10	5.00	ug/L
65794-96-9	3+4-Methylphenols	51.4		1.20	10.0	ug/L
621-64-7	n-Nitroso-di-n-propylamine	49.7		1.50	2.50	ug/L
67-72-1	Hexachloroethane	47.2		1.00	5.00	ug/L
98-95-3	Nitrobenzene	45.6		1.30	5.00	ug/L
78-59-1	Isophorone	49.4		1.10	5.00	ug/L
88-75-5	2-Nitrophenol	50.2		2.00	5.00	ug/L
105-67-9	2,4-Dimethylphenol	62.4		1.50	5.00	ug/L
111-91-1	bis(2-Chloroethoxy)methane	49.0		1.00	5.00	ug/L
120-83-2	2,4-Dichlorophenol	50.4		0.88	5.00	ug/L
91-20-3	Naphthalene	47.8		1.00	5.00	ug/L
106-47-8	4-Chloroaniline	24.4		1.30	5.00	ug/L
87-68-3	Hexachlorobutadiene	45.5		1.30	5.00	ug/L
105-60-2	Caprolactam	51.8		1.70	10.0	ug/L
59-50-7	4-Chloro-3-methylphenol	49.5		0.84	5.00	ug/L
91-57-6	2-Methylnaphthalene	49.9		1.10	5.00	ug/L
77-47-4	Hexachlorocyclopentadiene	130	E	5.00	10.0	ug/L
88-06-2	2,4,6-Trichlorophenol	47.6		0.89	5.00	ug/L
95-95-4	2,4,5-Trichlorophenol	47.4		1.00	5.00	ug/L
92-52-4	1,1-Biphenyl	48.1		0.91	5.00	ug/L
91-58-7	2-Chloronaphthalene	47.3		0.97	5.00	ug/L
88-74-4	2-Nitroaniline	48.5		1.40	5.00	ug/L
131-11-3	Dimethylphthalate	49.9		0.93	5.00	ug/L

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208-96-8	Acenaphthylene	52.0		1.00	5.00	ug/L
606-20-2	2,6-Dinitrotoluene	47.6		1.20	5.00	ug/L
99-09-2	3-Nitroaniline	31.1		1.40	5.00	ug/L
83-32-9	Acenaphthene	48.9		0.81	5.00	ug/L
51-28-5	2,4-Dinitrophenol	94.0	E	6.40	10.0	ug/L
100-02-7	4-Nitrophenol	88.6	E	2.00	10.0	ug/L
132-64-9	Dibenzofuran	49.2		0.93	5.00	ug/L
121-14-2	2,4-Dinitrotoluene	48.8		1.50	5.00	ug/L
84-66-2	Diethylphthalate	49.1		1.00	5.00	ug/L
7005-72-3	4-Chlorophenyl-phenylether	49.3		0.98	5.00	ug/L
86-73-7	Fluorene	49.7		0.96	5.00	ug/L
100-01-6	4-Nitroaniline	49.2		2.00	5.00	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	56.0		3.10	10.0	ug/L
86-30-6	n-Nitrosodiphenylamine	50.6		0.89	5.00	ug/L
101-55-3	4-Bromophenyl-phenylether	49.1		0.95	5.00	ug/L
118-74-1	Hexachlorobenzene	46.9		1.10	5.00	ug/L
1912-24-9	Atrazine	57.1		1.30	5.00	ug/L
87-86-5	Pentachlorophenol	84.9	E	1.90	10.0	ug/L
85-01-8	Phenanthrene	51.0		0.89	5.00	ug/L
120-12-7	Anthracene	53.5		1.10	5.00	ug/L
86-74-8	Carbazole	51.1		1.20	5.00	ug/L
84-74-2	Di-n-butylphthalate	50.2		1.50	5.00	ug/L
206-44-0	Fluoranthene	52.3		1.30	5.00	ug/L
129-00-0	Pyrene	46.8		1.10	5.00	ug/L
85-68-7	Butylbenzylphthalate	47.4		2.10	5.00	ug/L
91-94-1	3,3-Dichlorobenzidine	34.8		1.30	10.0	ug/L
56-55-3	Benzo(a)anthracene	51.9		0.94	5.00	ug/L
218-01-9	Chrysene	50.8		0.86	5.00	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	44.5		1.90	5.00	ug/L
117-84-0	Di-n-octyl phthalate	42.2		2.50	10.0	ug/L
205-99-2	Benzo(b)fluoranthene	49.8		1.10	5.00	ug/L

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207-08-9	Benzo(k)fluoranthene	57.9		1.20	5.00	ug/L
50-32-8	Benzo(a)pyrene	55.5		1.70	5.00	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	50.5		1.00	5.00	ug/L
53-70-3	Dibenzo(a,h)anthracene	50.0		1.20	5.00	ug/L
191-24-2	Benzo(g,h,i)perylene	47.2		1.20	5.00	ug/L
95-94-3	1,2,4,5-Tetrachlorobenzene	48.0		1.10	5.00	ug/L
123-91-1	1,4-Dioxane	43.3		1.30	5.00	ug/L
58-90-2	2,3,4,6-Tetrachlorophenol	55.3		0.79	5.00	ug/L
SURROGATES						
367-12-4	2-Fluorophenol	145		15 (10) - 110 (139)	97%	SPK: 150
13127-88-3	Phenol-d6	142		15 (10) - 110 (134)	94%	SPK: 150
4165-60-0	Nitrobenzene-d5	95.3		30 (49) - 130 (133)	95%	SPK: 100
321-60-8	2-Fluorobiphenyl	96.4		30 (52) - 130 (132)	96%	SPK: 100
118-79-6	2,4,6-Tribromophenol	150		15 (44) - 110 (137)	100%	SPK: 150
1718-51-0	Terphenyl-d14	96.1		30 (48) - 130 (125)	96%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	73200	6.863			
1146-65-2	Naphthalene-d8	279000	8.145			
15067-26-2	Acenaphthene-d10	158000	9.904			
1517-22-2	Phenanthrene-d10	303000	11.398			
1719-03-5	Chrysene-d12	197000	14.051			
1520-96-3	Perylene-d12	145000	15.551			

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