

Report of Analysis

Client:	RMJ Environomics, Inc.	Date Collected:	
Project:	245 Greenwood Ave	Date Received:	
Client Sample ID:	PB156921BL	SDG No.:	O5252
Lab Sample ID:	PB156921BL	Matrix:	SOIL
Analytical Method:	SW8270	% Solid:	100
Sample Wt/Vol:	30.01 Units: g	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 :	SW3541		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BF136177.D	1	11/06/23 09:48	11/07/23 15:47	PB156921

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
100-52-7	Benzaldehyde	0.15	U	0.15	0.33	mg/Kg
108-95-2	Phenol	0.074	U	0.074	0.17	mg/Kg
111-44-4	bis(2-Chloroethyl)ether	0.090	U	0.090	0.17	mg/Kg
95-57-8	2-Chlorophenol	0.072	U	0.072	0.17	mg/Kg
95-48-7	2-Methylphenol	0.11	U	0.11	0.17	mg/Kg
108-60-1	2,2-oxybis(1-Chloropropane)	0.10	U	0.10	0.17	mg/Kg
98-86-2	Acetophenone	0.087	U	0.087	0.17	mg/Kg
65794-96-9	3+4-Methylphenols	0.11	U	0.11	0.33	mg/Kg
621-64-7	n-Nitroso-di-n-propylamine	0.059	U	0.059	0.080	mg/Kg
67-72-1	Hexachloroethane	0.074	U	0.074	0.17	mg/Kg
98-95-3	Nitrobenzene	0.075	U	0.075	0.17	mg/Kg
78-59-1	Isophorone	0.068	U	0.068	0.17	mg/Kg
88-75-5	2-Nitrophenol	0.095	U	0.095	0.17	mg/Kg
105-67-9	2,4-Dimethylphenol	0.100	U	0.100	0.17	mg/Kg
111-91-1	bis(2-Chloroethoxy)methane	0.11	U	0.11	0.17	mg/Kg
120-83-2	2,4-Dichlorophenol	0.078	U	0.078	0.17	mg/Kg
91-20-3	Naphthalene	0.081	U	0.081	0.17	mg/Kg
106-47-8	4-Chloroaniline	0.10	U	0.10	0.17	mg/Kg
87-68-3	Hexachlorobutadiene	0.084	U	0.084	0.17	mg/Kg
105-60-2	Caprolactam	0.12	U	0.12	0.33	mg/Kg
59-50-7	4-Chloro-3-methylphenol	0.083	U	0.083	0.17	mg/Kg
91-57-6	2-Methylnaphthalene	0.095	U	0.095	0.17	mg/Kg
77-47-4	Hexachlorocyclopentadiene	0.21	U	0.21	0.33	mg/Kg
88-06-2	2,4,6-Trichlorophenol	0.077	U	0.077	0.17	mg/Kg
95-95-4	2,4,5-Trichlorophenol	0.088	U	0.088	0.17	mg/Kg
92-52-4	1,1-Biphenyl	0.091	U	0.091	0.17	mg/Kg
91-58-7	2-Chloronaphthalene	0.085	U	0.085	0.17	mg/Kg
88-74-4	2-Nitroaniline	0.099	U	0.099	0.17	mg/Kg
131-11-3	Dimethylphthalate	0.089	U	0.089	0.17	mg/Kg
208-96-8	Acenaphthylene	0.088	U	0.088	0.17	mg/Kg
606-20-2	2,6-Dinitrotoluene	0.090	U	0.090	0.17	mg/Kg

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99-09-2	3-Nitroaniline	0.093	U	0.093	0.17	mg/Kg
83-32-9	Acenaphthene	0.080	U	0.080	0.17	mg/Kg
51-28-5	2,4-Dinitrophenol	0.18	U	0.18	0.33	mg/Kg
100-02-7	4-Nitrophenol	0.11	U	0.11	0.33	mg/Kg
132-64-9	Dibenzofuran	0.077	U	0.077	0.17	mg/Kg
121-14-2	2,4-Dinitrotoluene	0.100	U	0.100	0.17	mg/Kg
84-66-2	Diethylphthalate	0.086	U	0.086	0.17	mg/Kg
7005-72-3	4-Chlorophenyl-phenylether	0.090	U	0.090	0.17	mg/Kg
86-73-7	Fluorene	0.085	U	0.085	0.17	mg/Kg
100-01-6	4-Nitroaniline	0.10	U	0.10	0.17	mg/Kg
534-52-1	4,6-Dinitro-2-methylphenol	0.088	U	0.088	0.33	mg/Kg
86-30-6	n-Nitrosodiphenylamine	0.093	U	0.093	0.17	mg/Kg
101-55-3	4-Bromophenyl-phenylether	0.097	U	0.097	0.17	mg/Kg
118-74-1	Hexachlorobenzene	0.099	U	0.099	0.17	mg/Kg
1912-24-9	Atrazine	0.095	U	0.095	0.17	mg/Kg
87-86-5	Pentachlorophenol	0.11	U	0.11	0.33	mg/Kg
85-01-8	Phenanthrene	0.093	U	0.093	0.17	mg/Kg
120-12-7	Anthracene	0.10	U	0.10	0.17	mg/Kg
86-74-8	Carbazole	0.085	U	0.085	0.17	mg/Kg
84-74-2	Di-n-butylphthalate	0.10	U	0.10	0.17	mg/Kg
206-44-0	Fluoranthene	0.094	U	0.094	0.17	mg/Kg
129-00-0	Pyrene	0.084	U	0.084	0.17	mg/Kg
85-68-7	Butylbenzylphthalate	0.10	U	0.10	0.17	mg/Kg
91-94-1	3,3-Dichlorobenzidine	0.16	U	0.16	0.33	mg/Kg
56-55-3	Benzo(a)anthracene	0.084	U	0.084	0.17	mg/Kg
218-01-9	Chrysene	0.087	U	0.087	0.17	mg/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	0.11	U	0.11	0.17	mg/Kg
117-84-0	Di-n-octyl phthalate	0.12	U	0.12	0.33	mg/Kg
205-99-2	Benzo(b)fluoranthene	0.080	U	0.080	0.17	mg/Kg
207-08-9	Benzo(k)fluoranthene	0.088	U	0.088	0.17	mg/Kg
50-32-8	Benzo(a)pyrene	0.094	U	0.094	0.17	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	0.11	U	0.11	0.17	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	0.097	U	0.097	0.17	mg/Kg

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191-24-2	Benzo(g,h,i)perylene	0.093	U	0.093	0.17	mg/Kg
95-94-3	1,2,4,5-Tetrachlorobenzene	0.089	U	0.089	0.17	mg/Kg
123-91-1	1,4-Dioxane	0.12	U	0.12	0.17	mg/Kg
58-90-2	2,3,4,6-Tetrachlorophenol	0.083	U	0.083	0.17	mg/Kg

SURROGATES

367-12-4	2-Fluorophenol	126		30 (18) - 130 (112)	84%	SPK: 150
13127-88-3	Phenol-d6	125		30 (15) - 130 (107)	84%	SPK: 150
4165-60-0	Nitrobenzene-d5	86.6		30 (18) - 130 (107)	87%	SPK: 100
321-60-8	2-Fluorobiphenyl	85.1		30 (20) - 130 (109)	85%	SPK: 100
118-79-6	2,4,6-Tribromophenol	131		30 (10) - 130 (110)	87%	SPK: 150
1718-51-0	Terphenyl-d14	84.9		30 (14) - 130 (112)	85%	SPK: 100

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	92000	6.816
1146-65-2	Naphthalene-d8	368000	8.098
15067-26-2	Acenaphthene-d10	189000	9.857
1517-22-2	Phenanthrene-d10	342000	11.351
1719-03-5	Chrysene-d12	196000	14.004
1520-96-3	Perylene-d12	181000	15.492

TENTATIVE IDENTIFIED COMPOUNDS

000556-71-8	Cyclononasiloxane, octadecamethyl-	160	J	14.5	ug/Kg
000107-52-8	Hexasiloxane, tetradecamethyl-	300	J	14.9	ug/Kg
	unknown15.351	370	J	15.4	ug/Kg
	unknown15.880	380	J	15.9	ug/Kg
	unknown16.510	350	J	16.5	ug/Kg
	unknown17.274	300	J	17.3	ug/Kg
000556-70-7	1,1,1,3,3,5,5,7,7,9,9,11,11,13,13,	280	J	18.2	ug/Kg

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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