

Report of Analysis

Client:	RU2 Engineering, LLC		Date Collected:	01/29/25	
Project:	NYCDDC SANTWOBR Brooklyn Bridge BBMCR		Date Received:	01/30/25	
Client Sample ID:	JPP-51.2-012925		SDG No.:	Q1235	
Lab Sample ID:	Q1235-01		Matrix:	SOIL	
Analytical Method:	SW8260		% Solid:	91.2	
Sample Wt/Vol:	9.09	Units: g	Final Vol:	5000	uL
Soil Aliquot Vol:		uL	Test:	VOCMS Group1	
GC Column:	RXI-624	ID : 0.25	Level :	LOW	
Prep Method :					

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021061.D	1		02/04/25 16:13	VY020425

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	3.00	U	1.00	3.00	ug/Kg
74-87-3	Chloromethane	3.00	U	0.70	3.00	ug/Kg
75-01-4	Vinyl Chloride	3.00	U	0.46	3.00	ug/Kg
74-83-9	Bromomethane	3.00	U	0.62	3.00	ug/Kg
75-00-3	Chloroethane	3.00	U	0.61	3.00	ug/Kg
75-69-4	Trichlorofluoromethane	3.00	U	0.55	3.00	ug/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	3.00	U	0.65	3.00	ug/Kg
75-65-0	Tert butyl alcohol	15.1	U	9.40	15.1	ug/Kg
75-35-4	1,1-Dichloroethene	3.00	U	0.47	3.00	ug/Kg
67-64-1	Acetone	15.6		3.80	15.1	ug/Kg
75-15-0	Carbon Disulfide	3.00	U	0.77	3.00	ug/Kg
1634-04-4	Methyl tert-butyl Ether	3.00	U	0.40	3.00	ug/Kg
79-20-9	Methyl Acetate	3.00	U	1.10	3.00	ug/Kg
75-09-2	Methylene Chloride	6.00	U	2.10	6.00	ug/Kg
156-60-5	trans-1,2-Dichloroethene	3.00	U	0.51	3.00	ug/Kg
75-34-3	1,1-Dichloroethane	3.00	U	0.38	3.00	ug/Kg
110-82-7	Cyclohexane	3.00	U	0.42	3.00	ug/Kg
78-93-3	2-Butanone	15.1	U	3.40	15.1	ug/Kg
56-23-5	Carbon Tetrachloride	3.00	U	0.52	3.00	ug/Kg
156-59-2	cis-1,2-Dichloroethene	3.00	U	0.37	3.00	ug/Kg
74-97-5	Bromochloromethane	3.00	U	1.50	3.00	ug/Kg
67-66-3	Chloroform	3.00	U	0.40	3.00	ug/Kg
71-55-6	1,1,1-Trichloroethane	3.00	U	0.47	3.00	ug/Kg
108-87-2	Methylcyclohexane	3.00	U	0.52	3.00	ug/Kg
71-43-2	Benzene	3.00	U	0.43	3.00	ug/Kg
107-06-2	1,2-Dichloroethane	3.00	U	0.37	3.00	ug/Kg
79-01-6	Trichloroethene	3.00	U	0.45	3.00	ug/Kg
78-87-5	1,2-Dichloropropane	3.00	U	0.40	3.00	ug/Kg
75-27-4	Bromodichloromethane	3.00	U	0.34	3.00	ug/Kg
108-10-1	4-Methyl-2-Pentanone	15.1	U	2.60	15.1	ug/Kg

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108-88-3	Toluene	3.00	U	0.40	3.00	ug/Kg
10061-02-6	t-1,3-Dichloropropene	3.00	U	0.36	3.00	ug/Kg
10061-01-5	cis-1,3-Dichloropropene	3.00	U	0.34	3.00	ug/Kg
79-00-5	1,1,2-Trichloroethane	3.00	U	0.51	3.00	ug/Kg
591-78-6	2-Hexanone	15.1	U	2.90	15.1	ug/Kg
124-48-1	Dibromochloromethane	3.00	U	0.39	3.00	ug/Kg
106-93-4	1,2-Dibromoethane	3.00	U	0.48	3.00	ug/Kg
127-18-4	Tetrachloroethene	3.00	U	0.54	3.00	ug/Kg
108-90-7	Chlorobenzene	3.00	U	0.45	3.00	ug/Kg
100-41-4	Ethyl Benzene	3.00	U	0.37	3.00	ug/Kg
179601-23-1	m/p-Xylenes	6.00	U	0.81	6.00	ug/Kg
95-47-6	o-Xylene	3.00	U	0.42	3.00	ug/Kg
100-42-5	Styrene	3.00	U	0.36	3.00	ug/Kg
75-25-2	Bromoform	3.00	U	0.49	3.00	ug/Kg
98-82-8	Isopropylbenzene	3.00	U	0.40	3.00	ug/Kg
79-34-5	1,1,2,2-Tetrachloroethane	3.00	U	0.66	3.00	ug/Kg
541-73-1	1,3-Dichlorobenzene	3.00	U	0.45	3.00	ug/Kg
106-46-7	1,4-Dichlorobenzene	3.00	U	0.48	3.00	ug/Kg
95-50-1	1,2-Dichlorobenzene	3.00	U	0.36	3.00	ug/Kg
96-12-8	1,2-Dibromo-3-Chloropropane	3.00	U	0.94	3.00	ug/Kg
120-82-1	1,2,4-Trichlorobenzene	3.00	U	0.48	3.00	ug/Kg
87-61-6	1,2,3-Trichlorobenzene	3.00	U	0.47	3.00	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	64.2		50 - 163	128%	SPK: 50
1868-53-7	Dibromofluoromethane	39.5		54 - 147	79%	SPK: 50
2037-26-5	Toluene-d8	50.4		58 - 134	101%	SPK: 50
460-00-4	4-Bromofluorobenzene	51.2		29 - 146	102%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	149000	7.707			
540-36-3	1,4-Difluorobenzene	281000	8.616			
3114-55-4	Chlorobenzene-d5	265000	11.414			
3855-82-1	1,4-Dichlorobenzene-d4	112000	13.346			

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TENTATIVE IDENTIFIED COMPOUNDS						
95-63-6	1,2,4-Trimethylbenzene	0.95	J		13.0	ug/Kg
91-20-3	Naphthalene	6.50	J		15.1	ug/Kg

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