

Report of Analysis

Client:	First Environment, Inc.	Date Collected:	07/09/25
Project:	White Plains Housing Authority - WPHAU006	Date Received:	07/09/25
Client Sample ID:	AOC-201	SDG No.:	Q2553
Lab Sample ID:	Q2553-01	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOC-TCLVOA-10
GC Column:	DB-624UI ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Date Analyzed	Prep Batch ID
VX046943.D	1	07/10/25 14:43	VX071025

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
75-71-8	Dichlorodifluoromethane	5.00	U	0.22	5.00	ug/L
74-87-3	Chloromethane	5.00	U	0.32	5.00	ug/L
75-01-4	Vinyl Chloride	5.00	U	0.26	5.00	ug/L
74-83-9	Bromomethane	5.00	U	1.40	5.00	ug/L
75-00-3	Chloroethane	5.00	U	0.47	5.00	ug/L
75-69-4	Trichlorofluoromethane	5.00	U	0.33	5.00	ug/L
76-13-1	1,1,2-Trichlorotrifluoroethane	5.00	U	0.25	5.00	ug/L
75-35-4	1,1-Dichloroethene	5.00	U	0.23	5.00	ug/L
67-64-1	Acetone	8.10	J	1.50	25.0	ug/L
75-15-0	Carbon Disulfide	0.94	J	0.21	5.00	ug/L
1634-04-4	Methyl tert-butyl Ether	5.00	U	0.16	5.00	ug/L
79-20-9	Methyl Acetate	5.00	U	0.27	5.00	ug/L
75-09-2	Methylene Chloride	5.00	U	0.28	5.00	ug/L
156-60-5	trans-1,2-Dichloroethene	5.00	U	0.23	5.00	ug/L
75-34-3	1,1-Dichloroethane	5.00	U	0.23	5.00	ug/L
110-82-7	Cyclohexane	19.5		1.50	5.00	ug/L
78-93-3	2-Butanone	25.0	U	0.98	25.0	ug/L
56-23-5	Carbon Tetrachloride	5.00	U	0.25	5.00	ug/L
156-59-2	cis-1,2-Dichloroethene	5.00	U	0.19	5.00	ug/L
74-97-5	Bromochloromethane	5.00	U	0.22	5.00	ug/L
67-66-3	Chloroform	5.00	U	0.25	5.00	ug/L
71-55-6	1,1,1-Trichloroethane	5.00	U	0.20	5.00	ug/L
108-87-2	Methylcyclohexane	48.2		0.16	5.00	ug/L
71-43-2	Benzene	5.00	U	0.15	5.00	ug/L
107-06-2	1,2-Dichloroethane	5.00	U	0.22	5.00	ug/L
79-01-6	Trichloroethene	5.00	U	0.090	5.00	ug/L
78-87-5	1,2-Dichloropropane	5.00	U	0.20	5.00	ug/L
75-27-4	Bromodichloromethane	5.00	U	0.22	5.00	ug/L
108-10-1	4-Methyl-2-Pentanone	25.0	U	0.68	25.0	ug/L
108-88-3	Toluene	5.00	U	0.14	5.00	ug/L

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000107-83-5	Pentane, 2-methyl-	92.1	J		2.84	ug/L
000565-59-3	Pentane, 2,3-dimethyl-	86.1	J		5.63	ug/L
000589-34-4	Hexane, 3-methyl-	91.5	J		5.84	ug/L
103-65-1	n-propylbenzene	280	J		11.3	ug/L
000611-14-3	Benzene, 1-ethyl-2-methyl-	150	J		11.4	ug/L
108-67-8	1,3,5-Trimethylbenzene	46.3	J		11.5	ug/L
000622-96-8	Benzene, 1-ethyl-4-methyl-	120	J		11.6	ug/L
95-63-6	1,2,4-Trimethylbenzene	680	J		11.8	ug/L
135-98-8	sec-Butylbenzene	25.0	J		11.9	ug/L
99-87-6	p-Isopropyltoluene	7.80	J		12.0	ug/L
000526-73-8	Benzene, 1,2,3-trimethyl-	170	J		12.1	ug/L
000300-57-2	Benzene, 2-propenyl-	190	J		12.2	ug/L
104-51-8	n-Butylbenzene	34.4	J		12.3	ug/L
001758-88-9	Benzene, 2-ethyl-1,4-dimethyl-	140	J		12.5	ug/L
000934-80-5	Benzene, 4-ethyl-1,2-dimethyl-	280	J		12.6	ug/L
000767-58-8	Indan, 1-methyl-	190	J		12.7	ug/L
000095-93-2	Benzene, 1,2,4,5-tetramethyl-	160	J		12.9	ug/L
000488-23-3	Benzene, 1,2,3,4-tetramethyl-	200	J		13.0	ug/L
003454-07-7	Benzene, 1-ethenyl-4-ethyl-	150	J		13.2	ug/L
000824-22-6	1H-Indene, 2,3-dihydro-4-methyl-	340	J		13.3	ug/L

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