

Report of Analysis

Client:	G Environmental		Date Collected:	06/18/25	
Project:	Capra		Date Received:	06/18/25	
Client Sample ID:	GCAP1W		SDG No.:	Q2363	
Lab Sample ID:	Q2363-02		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:	Prep Date	Date Analyzed	Prep Batch ID
VX046768.D	1		06/19/25 13:36	VX061925

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
75-71-8	Dichlorodifluoromethane	0.22	U	0.22	1.00	ug/L
74-87-3	Chloromethane	0.32	U	0.32	1.00	ug/L
75-01-4	Vinyl Chloride	0.26	U	0.26	1.00	ug/L
74-83-9	Bromomethane	1.40	U	1.40	5.00	ug/L
75-00-3	Chloroethane	0.47	U	0.47	1.00	ug/L
75-69-4	Trichlorofluoromethane	0.33	U	0.33	1.00	ug/L
76-13-1	1,1,2-Trichlorotrifluoroethane	0.25	U	0.25	1.00	ug/L
75-35-4	1,1-Dichloroethene	0.23	U	0.23	1.00	ug/L
67-64-1	Acetone	1.50	U	1.50	5.00	ug/L
75-15-0	Carbon Disulfide	0.21	U	0.21	1.00	ug/L
1634-04-4	Methyl tert-butyl Ether	0.16	U	0.16	1.00	ug/L
79-20-9	Methyl Acetate	0.27	U	0.27	1.00	ug/L
75-09-2	Methylene Chloride	0.28	U	0.28	1.00	ug/L
156-60-5	trans-1,2-Dichloroethene	0.23	U	0.23	1.00	ug/L
75-34-3	1,1-Dichloroethane	0.23	U	0.23	1.00	ug/L
110-82-7	Cyclohexane	1200	E	1.50	5.00	ug/L
78-93-3	2-Butanone	0.98	U	0.98	5.00	ug/L
56-23-5	Carbon Tetrachloride	0.25	U	0.25	1.00	ug/L
156-59-2	cis-1,2-Dichloroethene	0.19	U	0.19	1.00	ug/L
74-97-5	Bromochloromethane	0.22	U	0.22	1.00	ug/L
67-66-3	Chloroform	0.25	U	0.25	1.00	ug/L
71-55-6	1,1,1-Trichloroethane	0.20	U	0.20	1.00	ug/L
108-87-2	Methylcyclohexane	1000	E	0.16	1.00	ug/L
71-43-2	Benzene	1.80		0.15	1.00	ug/L
107-06-2	1,2-Dichloroethane	0.22	U	0.22	1.00	ug/L
79-01-6	Trichloroethene	0.090	U	0.090	1.00	ug/L
78-87-5	1,2-Dichloropropane	0.20	U	0.20	1.00	ug/L
75-27-4	Bromodichloromethane	0.22	U	0.22	1.00	ug/L
108-10-1	4-Methyl-2-Pentanone	0.68	U	0.68	5.00	ug/L
108-88-3	Toluene	6.70		0.14	1.00	ug/L

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[illegible]

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CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
000078-78-4	Butane, 2-methyl-	170	J		1.77	ug/L
000109-66-0	Pentane	160	J		1.97	ug/L
000107-83-5	Pentane, 2-methyl-	110	J		2.84	ug/L
000096-14-0	Pentane, 3-methyl-	130	J		3.12	ug/L
000110-54-3	n-Hexane	130	J		3.46	ug/L
000096-37-7	Cyclopentane, methyl-	580	J		4.32	ug/L
000822-50-4	Cyclopentane, 1,2-dimethyl-, trans	130	J		6.37	ug/L
000591-49-1	Cyclohexene, 1-methyl-	200	J		8.51	ug/L
006876-23-9	Cyclohexane, 1,2-dimethyl-, trans-	160	J		8.98	ug/L
003073-66-3	Cyclohexane, 1,1,3-trimethyl-	130	J		9.62	ug/L
	unknown10.780	120	J		10.8	ug/L
103-65-1	n-propylbenzene	190	J		11.3	ug/L
135-98-8	sec-Butylbenzene	23.9	J		11.9	ug/L
99-87-6	p-Isopropyltoluene	20.8	J		12.0	ug/L
000496-11-7	Indane	220	J		12.2	ug/L
104-51-8	n-Butylbenzene	23.9	J		12.3	ug/L
002870-04-4	Benzene, 2-ethyl-1,3-dimethyl-	110	J		12.6	ug/L
007525-62-4	Benzene, 1-ethenyl-3-ethyl-	120	J		12.7	ug/L
000934-10-1	3-Phenylbut-1-ene	180	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