

Report of Analysis

Client:	G Environmental	Date Collected:	03/13/25
Project:	Communipaw	Date Received:	03/14/25
Client Sample ID:	MW3	SDG No.:	Q1575
Lab Sample ID:	Q1575-01	Matrix:	Water
Analytical Method:	SW8260	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOC-TCLVOA-10
GC Column:	RXI-624 ID : 0.25	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VN086031.D	1		03/20/25 17:57	VN032025

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	4.90	J	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	1.50	U	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	6.90		0.16	1.00	ug/L
71-43-2	Benzene	0.15	U	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	0.14	U	0.14	1.00	ug/L

Report of Analysis

Client:	G Environmental	Date Collected:	03/13/25
Project:	Communipaw	Date Received:	03/14/25
Client Sample ID:	MW3	SDG No.:	Q1575
Lab Sample ID:	Q1575-01	Matrix:	Water
Analytical Method:	SW8260	% Solid:	0
Sample Wt/Vol:	5	Units:	mL
Soil Aliquot Vol:			uL
GC Column:	RXI-624	ID :	0.25
Prep Method :		Level :	LOW

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VN086031.D	1		03/20/25 17:57	VN032025

[illegible]

Report of Analysis

Client:	G Environmental	Date Collected:	03/13/25
Project:	Communipaw	Date Received:	03/14/25
Client Sample ID:	MW3	SDG No.:	Q1575
Lab Sample ID:	Q1575-01	Matrix:	Water
Analytical Method:	SW8260	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOC-TCLVOA-10
GC Column:	RXI-624 ID : 0.25	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VN086031.D	1		03/20/25 17:57	VN032025

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
000079-29-8	Butane, 2,3-dimethyl-	11.8	J		5.28	ug/L
000108-08-7	Pentane, 2,4-dimethyl-	22.6	J		7.22	ug/L
000565-59-3	Pentane, 2,3-dimethyl-	59.9	J		8.32	ug/L
000594-82-1	Butane, 2,2,3,3-tetramethyl-	240	J		8.76	ug/L
013475-82-6	Heptane, 2,2,4,6,6-pentamethyl-	16.2	J		9.72	ug/L
000565-75-3	Pentane, 2,3,4-trimethyl-	130	J		10.0	ug/L
000560-21-4	Pentane, 2,3,3-trimethyl-	210	J		10.1	ug/L
	unknown10.318	11.8	J		10.3	ug/L
003522-94-9	Hexane, 2,2,5-trimethyl-	16.1	J		10.5	ug/L
	unknown11.006	5.80	J		11.0	ug/L
000565-80-0	3-Pentanone, 2,4-dimethyl-	8.90	J		11.2	ug/L
000564-04-5	3-Pentanone, 2,2-dimethyl-	50.9	J		11.2	ug/L
103-65-1	n-propylbenzene	2.70	J		13.0	ug/L
95-63-6	1,2,4-Trimethylbenzene	0.37	J		13.5	ug/L
1000309-17-6	Sulfurous acid, butyl nonyl ester	8.10	J		13.6	ug/L
000496-11-7	Indane	6.20	J		14.0	ug/L
104-51-8	n-Butylbenzene	0.90	J		14.1	ug/L
007525-62-4	Benzene, 1-ethenyl-3-ethyl-	5.40	J		15.0	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