

Report of Analysis

Client:	Garden State Laboratories, Inc.		Date Collected:	
Project:	Waste Water 2025		Date Received:	
Client Sample ID:	VN0612WBSD01		SDG No.:	Q2302
Lab Sample ID:	VN0612WBSD01		Matrix:	Water
Analytical Method:	8260D		% Solid:	0
Sample Wt/Vol:	5	Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:		uL	Test:	VOCMS Group2
GC Column:	RXI-624	ID : 0.25	Level :	LOW
Prep Method :				

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VN086971.D	1		06/12/25 12:53	VN061225

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
75-71-8	Dichlorodifluoromethane	20.7		0.22	1.00	ug/L
74-87-3	Chloromethane	16.8		0.32	1.00	ug/L
75-01-4	Vinyl Chloride	20.9		0.26	1.00	ug/L
74-83-9	Bromomethane	16.7		1.40	5.00	ug/L
75-00-3	Chloroethane	21.2		0.47	1.00	ug/L
75-69-4	Trichlorofluoromethane	21.0		0.33	1.00	ug/L
76-13-1	1,1,2-Trichlorotrifluoroethane	21.6		0.25	1.00	ug/L
75-35-4	1,1-Dichloroethene	20.7		0.23	1.00	ug/L
67-64-1	Acetone	100		1.50	5.00	ug/L
75-15-0	Carbon Disulfide	18.8		0.21	1.00	ug/L
1634-04-4	Methyl tert-butyl Ether	22.7		0.16	1.00	ug/L
79-20-9	Methyl Acetate	22.9		0.27	1.00	ug/L
75-09-2	Methylene Chloride	21.7		0.28	1.00	ug/L
156-60-5	trans-1,2-Dichloroethene	20.6		0.23	1.00	ug/L
75-34-3	1,1-Dichloroethane	21.5		0.23	1.00	ug/L
110-82-7	Cyclohexane	18.8		1.50	5.00	ug/L
78-93-3	2-Butanone	110		0.98	5.00	ug/L
56-23-5	Carbon Tetrachloride	21.2		0.25	1.00	ug/L
156-59-2	cis-1,2-Dichloroethene	21.7		0.19	1.00	ug/L
74-97-5	Bromochloromethane	24.5		0.22	1.00	ug/L
67-66-3	Chloroform	22.1		0.25	1.00	ug/L
71-55-6	1,1,1-Trichloroethane	21.2		0.20	1.00	ug/L
108-87-2	Methylcyclohexane	17.9		0.16	1.00	ug/L
71-43-2	Benzene	21.6		0.15	1.00	ug/L
107-06-2	1,2-Dichloroethane	23.0		0.22	1.00	ug/L
79-01-6	Trichloroethene	21.5		0.090	1.00	ug/L
78-87-5	1,2-Dichloropropane	22.4		0.20	1.00	ug/L
75-27-4	Bromodichloromethane	23.2		0.22	1.00	ug/L
108-10-1	4-Methyl-2-Pentanone	120		0.68	5.00	ug/L
108-88-3	Toluene	21.8		0.14	1.00	ug/L

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10061-02-6	t-1,3-Dichloropropene	23.1		0.17	1.00	ug/L
10061-01-5	cis-1,3-Dichloropropene	22.5		0.16	1.00	ug/L
79-00-5	1,1,2-Trichloroethane	23.2		0.21	1.00	ug/L
591-78-6	2-Hexanone	100		0.89	5.00	ug/L
124-48-1	Dibromochloromethane	23.3		0.18	1.00	ug/L
106-93-4	1,2-Dibromoethane	23.0		0.15	1.00	ug/L
127-18-4	Tetrachloroethene	20.7		0.23	1.00	ug/L
108-90-7	Chlorobenzene	21.6		0.12	1.00	ug/L
100-41-4	Ethyl Benzene	20.7		0.13	1.00	ug/L
179601-23-1	m/p-Xylenes	42.3		0.24	2.00	ug/L
95-47-6	o-Xylene	21.6		0.12	1.00	ug/L
100-42-5	Styrene	21.8		0.15	1.00	ug/L
75-25-2	Bromoform	24.5		0.19	1.00	ug/L
98-82-8	Isopropylbenzene	20.3		0.12	1.00	ug/L
79-34-5	1,1,2,2-Tetrachloroethane	23.8		0.26	1.00	ug/L
541-73-1	1,3-Dichlorobenzene	21.5		0.16	1.00	ug/L
106-46-7	1,4-Dichlorobenzene	21.7		0.19	1.00	ug/L
95-50-1	1,2-Dichlorobenzene	22.4		0.16	1.00	ug/L
96-12-8	1,2-Dibromo-3-Chloropropane	21.6		0.53	1.00	ug/L
120-82-1	1,2,4-Trichlorobenzene	19.9		0.20	1.00	ug/L
87-61-6	1,2,3-Trichlorobenzene	19.6		0.20	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	54.6		74 - 125	109%	SPK: 50
1868-53-7	Dibromofluoromethane	57.4		75 - 124	115%	SPK: 50
2037-26-5	Toluene-d8	54.1		86 - 113	108%	SPK: 50
460-00-4	4-Bromofluorobenzene	55.3		77 - 121	111%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	164000	8.23			
540-36-3	1,4-Difluorobenzene	293000	9.106			
3114-55-4	Chlorobenzene-d5	261000	11.865			
3855-82-1	1,4-Dichlorobenzene-d4	129000	13.794			

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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