

Report of Analysis

Client:	Tetra Tech NUS, Inc.		Date Collected:	09/22/25	
Project:	NWIRP Bethpage 112G08005-WE13		Date Received:	09/22/25	
Client Sample ID:	BP-VPB-184-GW-320-322MSD		SDG No.:	Q3165	
Lab Sample ID:	Q3165-10MSD		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
VX047822.D	1	09/25/25 18:48	VX092525

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
TARGETS							
75-71-8	Dichlorodifluoromethane	53.8		0.22	0.50	1.00	ug/L
74-87-3	Chloromethane	68.2		0.32	0.50	1.00	ug/L
75-01-4	Vinyl Chloride	63.8		0.26	0.75	1.00	ug/L
74-83-9	Bromomethane	65.0		1.40	3.80	5.00	ug/L
75-00-3	Chloroethane	66.2		0.47	0.75	1.00	ug/L
75-69-4	Trichlorofluoromethane	52.9		0.33	0.50	1.00	ug/L
76-13-1	1,1,2-Trichlorotrifluoroethane	41.7		0.25	0.50	1.00	ug/L
75-35-4	1,1-Dichloroethene	61.8		0.23	0.75	1.00	ug/L
67-64-1	Acetone	300		1.50	3.80	5.00	ug/L
75-15-0	Carbon Disulfide	44.3		0.21	0.75	1.00	ug/L
1634-04-4	Methyl tert-butyl Ether	68.1		0.16	0.50	1.00	ug/L
79-20-9	Methyl Acetate	56.1		0.27	0.75	1.00	ug/L
75-09-2	Methylene Chloride	67.0		0.28	0.50	1.00	ug/L
156-60-5	trans-1,2-Dichloroethene	57.7		0.23	0.50	1.00	ug/L
75-34-3	1,1-Dichloroethane	65.2		0.23	0.50	1.00	ug/L
110-82-7	Cyclohexane	37.5		1.50	2.50	5.00	ug/L
78-93-3	2-Butanone	360		0.98	2.50	5.00	ug/L
56-23-5	Carbon Tetrachloride	35.6		0.25	0.50	1.00	ug/L
156-59-2	cis-1,2-Dichloroethene	63.6		0.19	0.75	1.00	ug/L
74-97-5	Bromochloromethane	66.6		0.22	0.50	1.00	ug/L
67-66-3	Chloroform	69.3		0.25	0.50	1.00	ug/L
71-55-6	1,1,1-Trichloroethane	58.7		0.20	0.50	1.00	ug/L
108-87-2	Methylcyclohexane	27.2		0.16	0.50	1.00	ug/L
71-43-2	Benzene	58.2		0.15	0.50	1.00	ug/L
107-06-2	1,2-Dichloroethane	59.7		0.22	0.50	1.00	ug/L
79-01-6	Trichloroethene	56.2		0.090	0.75	1.00	ug/L
78-87-5	1,2-Dichloropropane	61.1		0.20	0.50	1.00	ug/L
75-27-4	Bromodichloromethane	44.8		0.22	0.50	1.00	ug/L
108-10-1	4-Methyl-2-Pentanone	350		0.68	2.50	5.00	ug/L
108-88-3	Toluene	53.2		0.14	0.50	1.00	ug/L

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10061-02-6	t-1,3-Dichloropropene	42.1		0.17	0.50	1.00	ug/L
10061-01-5	cis-1,3-Dichloropropene	42.4		0.16	0.50	1.00	ug/L
79-00-5	1,1,2-Trichloroethane	64.7		0.21	0.50	1.00	ug/L
591-78-6	2-Hexanone	340		0.89	2.50	5.00	ug/L
124-48-1	Dibromochloromethane	44.0		0.18	0.50	1.00	ug/L
106-93-4	1,2-Dibromoethane	63.4		0.15	0.50	1.00	ug/L
127-18-4	Tetrachloroethene	49.7		0.23	0.50	1.00	ug/L
108-90-7	Chlorobenzene	53.9		0.12	0.50	1.00	ug/L
100-41-4	Ethyl Benzene	49.0		0.13	0.50	1.00	ug/L
179601-23-1	m/p-Xylenes	97.5		0.24	1.00	2.00	ug/L
95-47-6	o-Xylene	51.7		0.12	0.50	1.00	ug/L
100-42-5	Styrene	42.3		0.15	0.50	1.00	ug/L
75-25-2	Bromoform	43.1		0.19	0.50	1.00	ug/L
98-82-8	Isopropylbenzene	50.2		0.12	0.50	1.00	ug/L
79-34-5	1,1,2,2-Tetrachloroethane	76.6		0.26	0.50	1.00	ug/L
541-73-1	1,3-Dichlorobenzene	51.1		0.16	0.50	1.00	ug/L
106-46-7	1,4-Dichlorobenzene	50.8		0.19	0.50	1.00	ug/L
95-50-1	1,2-Dichlorobenzene	55.9		0.16	0.50	1.00	ug/L
96-12-8	1,2-Dibromo-3-Chloropropane	68.7		0.53	0.75	1.00	ug/L
120-82-1	1,2,4-Trichlorobenzene	42.8		0.20	0.50	1.00	ug/L
87-61-6	1,2,3-Trichlorobenzene	48.3		0.20	0.75	1.00	ug/L
SURROGATES							
17060-07-0	1,2-Dichloroethane-d4	61.8	*	81 - 118		124%	SPK: 50
1868-53-7	Dibromofluoromethane	56.6		80 - 119		113%	SPK: 50
2037-26-5	Toluene-d8	49.9		89 - 112		100%	SPK: 50
460-00-4	4-Bromofluorobenzene	46.0		85 - 114		92%	SPK: 50
INTERNAL STANDARDS							
363-72-4	Pentafluorobenzene	108000	5.544				
540-36-3	1,4-Difluorobenzene	211000	6.751				
3114-55-4	Chlorobenzene-d5	180000	10.037				
3855-82-1	1,4-Dichlorobenzene-d4	77400	12.006				

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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