

Report of Analysis

Client:	AECOM Technical Services, Inc.		Date Collected:	12/10/24	
Project:	NAVFAC NWIRP Bethpage, NY Site 1 OU-2 - 32258		Date Received:	12/11/24	
Client Sample ID:	RE120D3-20241210DL		SDG No.:	P5261	
Lab Sample ID:	P5261-19DL		Matrix:	Water	
Analytical Method:	SW8260		% Solid:	0	
Sample Wt/Vol:	5	Units: mL	Final Vol:	5000	uL
Soil Aliquot Vol:		uL	Test:	VOCMS Group1	
GC Column:	DB-624UI	ID : 0.18	Level :	LOW	
Prep Method :					

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX044380.D	20		12/18/24 14:01	VX121824

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
TARGETS							
75-71-8	Dichlorodifluoromethane	10.0	UD	4.20	10.0	20.0	ug/L
74-87-3	Chloromethane	10.0	UD	7.00	10.0	20.0	ug/L
75-01-4	Vinyl Chloride	15.0	UD	6.80	15.0	20.0	ug/L
74-83-9	Bromomethane	75.0	UD	27.2	75.0	100	ug/L
75-00-3	Chloroethane	15.0	UD	11.2	15.0	20.0	ug/L
75-69-4	Trichlorofluoromethane	10.0	UD	6.80	10.0	20.0	ug/L
76-13-1	1,1,2-Trichlorotrifluoroethane	14.6	JD	5.00	10.0	20.0	ug/L
75-35-4	1,1-Dichloroethene	8.00	JD	5.20	15.0	20.0	ug/L
67-64-1	Acetone	75.0	UD	27.8	75.0	100	ug/L
75-15-0	Carbon Disulfide	15.0	UD	6.40	15.0	20.0	ug/L
1634-04-4	Methyl tert-butyl Ether	10.0	UD	3.20	10.0	20.0	ug/L
79-20-9	Methyl Acetate	15.0	UD	12.0	15.0	20.0	ug/L
75-09-2	Methylene Chloride	10.0	UD	6.40	10.0	20.0	ug/L
156-60-5	trans-1,2-Dichloroethene	10.0	UD	5.00	10.0	20.0	ug/L
75-34-3	1,1-Dichloroethane	10.0	UD	4.60	10.0	20.0	ug/L
110-82-7	Cyclohexane	50.0	UD	32.4	50.0	100	ug/L
78-93-3	2-Butanone	50.0	UD	26.0	50.0	100	ug/L
56-23-5	Carbon Tetrachloride	10.0	UD	5.00	10.0	20.0	ug/L
156-59-2	cis-1,2-Dichloroethene	15.0	UD	5.00	15.0	20.0	ug/L
74-97-5	Bromochloromethane	10.0	UD	3.60	10.0	20.0	ug/L
67-66-3	Chloroform	10.0	UD	5.20	10.0	20.0	ug/L
71-55-6	1,1,1-Trichloroethane	10.0	UD	3.80	10.0	20.0	ug/L
108-87-2	Methylcyclohexane	10.0	UD	3.80	10.0	20.0	ug/L
71-43-2	Benzene	10.0	UD	3.20	10.0	20.0	ug/L
107-06-2	1,2-Dichloroethane	10.0	UD	4.80	10.0	20.0	ug/L
79-01-6	Trichloroethene	730	D	6.40	15.0	20.0	ug/L
78-87-5	1,2-Dichloropropane	10.0	UD	3.80	10.0	20.0	ug/L
75-27-4	Bromodichloromethane	10.0	UD	4.80	10.0	20.0	ug/L
108-10-1	4-Methyl-2-Pentanone	50.0	UD	15.0	50.0	100	ug/L
108-88-3	Toluene	10.0	UD	3.60	10.0	20.0	ug/L

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CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
10061-02-6	t-1,3-Dichloropropene	10.0	UD	4.20	10.0	20.0	ug/L
10061-01-5	cis-1,3-Dichloropropene	10.0	UD	3.60	10.0	20.0	ug/L
79-00-5	1,1,2-Trichloroethane	10.0	UD	4.20	10.0	20.0	ug/L
591-78-6	2-Hexanone	50.0	UD	22.6	50.0	100	ug/L
124-48-1	Dibromochloromethane	10.0	UD	3.60	10.0	20.0	ug/L
106-93-4	1,2-Dibromoethane	10.0	UD	3.20	10.0	20.0	ug/L
127-18-4	Tetrachloroethene	12.1	JD	5.00	10.0	20.0	ug/L
108-90-7	Chlorobenzene	10.0	UD	2.60	10.0	20.0	ug/L
100-41-4	Ethyl Benzene	10.0	UD	3.20	10.0	20.0	ug/L
179601-23-1	m/p-Xylenes	20.0	UD	6.20	20.0	40.0	ug/L
95-47-6	o-Xylene	10.0	UD	2.80	10.0	20.0	ug/L
100-42-5	Styrene	10.0	UD	3.20	10.0	20.0	ug/L
75-25-2	Bromoform	10.0	UD	4.20	10.0	20.0	ug/L
98-82-8	Isopropylbenzene	10.0	UD	2.60	10.0	20.0	ug/L
79-34-5	1,1,2,2-Tetrachloroethane	10.0	UD	5.40	10.0	20.0	ug/L
541-73-1	1,3-Dichlorobenzene	10.0	UD	4.80	10.0	20.0	ug/L
106-46-7	1,4-Dichlorobenzene	10.0	UD	5.40	10.0	20.0	ug/L
95-50-1	1,2-Dichlorobenzene	10.0	UD	3.80	10.0	20.0	ug/L
96-12-8	1,2-Dibromo-3-Chloropropane	15.0	UD	9.20	15.0	20.0	ug/L
120-82-1	1,2,4-Trichlorobenzene	10.0	UD	8.40	10.0	20.0	ug/L
87-61-6	1,2,3-Trichlorobenzene	15.0	UD	10.2	15.0	20.0	ug/L
SURROGATES							
17060-07-0	1,2-Dichloroethane-d4	45.1		81 - 118		90%	SPK: 50
1868-53-7	Dibromofluoromethane	46.7		80 - 119		93%	SPK: 50
2037-26-5	Toluene-d8	50.7		89 - 112		101%	SPK: 50
460-00-4	4-Bromofluorobenzene	49.4		85 - 114		99%	SPK: 50
INTERNAL STANDARDS							
363-72-4	Pentafluorobenzene	115000	5.55				
540-36-3	1,4-Difluorobenzene	215000	6.757				
3114-55-4	Chlorobenzene-d5	187000	10.055				
3855-82-1	1,4-Dichlorobenzene-d4	83900	12.024				

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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