



Client:	JACOBS Engineering Group, Inc.	Date Collected:	08/01/24
Project:	Former Schlumberger Site Princeton NJ	Date Received:	08/01/24
Client Sample ID:	923-K1-WS-080124MS	SDG No.:	P3440
Lab Sample ID:	P3440-02MS	Matrix:	Water
Analytical Method:	SW8260	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group6
GC Column:	RXI-624 ID : 0.25	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VN083252.D	1		08/13/24 12:55	VN081324

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
75-71-8	Dichlorodifluoromethane	48.6		0.21	1.00	ug/L
74-87-3	Chloromethane	51.0		0.35	1.00	ug/L
75-01-4	Vinyl Chloride	52.9		0.34	1.00	ug/L
74-83-9	Bromomethane	50.3		1.40	5.00	ug/L
75-00-3	Chloroethane	53.4		0.56	1.00	ug/L
76-13-1	1,1,2-Trichlorotrifluoroethane	50.2		0.25	1.00	ug/L
75-35-4	1,1-Dichloroethene	49.8		0.26	1.00	ug/L
67-64-1	Acetone	240		1.40	5.00	ug/L
75-15-0	Carbon Disulfide	44.2		0.32	1.00	ug/L
1634-04-4	Methyl tert-butyl Ether	52.8		0.16	1.00	ug/L
75-09-2	Methylene Chloride	49.1		0.32	1.00	ug/L
156-60-5	trans-1,2-Dichloroethene	49.7		0.25	1.00	ug/L
75-34-3	1,1-Dichloroethane	53.9		0.23	1.00	ug/L
110-82-7	Cyclohexane	46.8		1.60	5.00	ug/L
78-93-3	2-Butanone	240		1.30	5.00	ug/L
56-23-5	Carbon Tetrachloride	51.0		0.25	1.00	ug/L
156-59-2	cis-1,2-Dichloroethene	52.3		0.25	1.00	ug/L
67-66-3	Chloroform	55.8		0.26	1.00	ug/L
71-55-6	1,1,1-Trichloroethane	54.1		0.19	1.00	ug/L
108-87-2	Methylcyclohexane	43.7		0.19	1.00	ug/L
71-43-2	Benzene	50.3		0.16	1.00	ug/L
107-06-2	1,2-Dichloroethane	52.0		0.24	1.00	ug/L
79-01-6	Trichloroethene	48.7		0.32	1.00	ug/L
75-27-4	Bromodichloromethane	50.7		0.24	1.00	ug/L
108-88-3	Toluene	51.0		0.18	1.00	ug/L
79-00-5	1,1,2-Trichloroethane	52.3		0.21	1.00	ug/L
124-48-1	Dibromochloromethane	52.7		0.18	1.00	ug/L
127-18-4	Tetrachloroethene	46.0		0.25	1.00	ug/L
108-90-7	Chlorobenzene	49.6		0.13	1.00	ug/L
100-41-4	Ethyl Benzene	49.7		0.16	1.00	ug/L

Report of Analysis

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CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
179601-23-1	m/p-Xylenes	99.8		0.31	2.00	ug/L
1330-20-7	Total Xylenes	150		0.45	3.00	ug/L
95-47-6	o-Xylene	49.7		0.14	1.00	ug/L
98-82-8	Isopropylbenzene	48.6		0.13	1.00	ug/L
106-46-7	1,4-Dichlorobenzene	48.4		0.27	1.00	ug/L
95-50-1	1,2-Dichlorobenzene	49.8		0.19	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	56.0		70 (74) - 130 (125)	112%	SPK: 50
1868-53-7	Dibromofluoromethane	50.6		70 (75) - 130 (124)	101%	SPK: 50
2037-26-5	Toluene-d8	49.9		70 (86) - 130 (113)	100%	SPK: 50
460-00-4	4-Bromofluorobenzene	50.4		70 (77) - 130 (121)	101%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	160000	8.229			
540-36-3	1,4-Difluorobenzene	294000	9.1			
3114-55-4	Chlorobenzene-d5	257000	11.865			
3855-82-1	1,4-Dichlorobenzene-d4	121000	13.794			

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