

Report of Analysis

Client:	Louis Berger U.S., Inc., A WSP Company	Date Collected:	01/18/23
Project:	NYCDDC Phase II SCI Arthur Kill Road CEQR	Date Received:	01/19/23
Client Sample ID:	B-P38B	SDG No.:	O1233
Lab Sample ID:	O1233-09	Matrix:	SOIL
Analytical Method:	SW8260	% Solid:	80.7
Sample Wt/Vol:	5.01 Units: g	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOC-TCLVOA-10
GC Column:	RTX-VMS ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VD075162.D	1		01/20/23 19:17	VD012023

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	1.10	U	1.10	6.20	ug/Kg
74-87-3	Chloromethane	1.30	U	1.30	6.20	ug/Kg
75-01-4	Vinyl Chloride	1.10	U	1.10	6.20	ug/Kg
74-83-9	Bromomethane	1.40	U	1.40	6.20	ug/Kg
75-00-3	Chloroethane	1.10	U	1.10	6.20	ug/Kg
75-69-4	Trichlorofluoromethane	1.20	U	1.20	6.20	ug/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	0.89	U	0.89	6.20	ug/Kg
75-35-4	1,1-Dichloroethene	1.10	U	1.10	6.20	ug/Kg
67-64-1	Acetone	15.1	U	15.1	30.9	ug/Kg
75-15-0	Carbon Disulfide	0.93	U	0.93	6.20	ug/Kg
1634-04-4	Methyl tert-butyl Ether	1.20	U	1.20	6.20	ug/Kg
79-20-9	Methyl Acetate	1.60	U	1.60	6.20	ug/Kg
75-09-2	Methylene Chloride	7.40	U	7.40	12.4	ug/Kg
156-60-5	trans-1,2-Dichloroethene	0.84	U	0.84	6.20	ug/Kg
75-34-3	1,1-Dichloroethane	0.87	U	0.87	6.20	ug/Kg
110-82-7	Cyclohexane	1.00	U	1.00	6.20	ug/Kg
78-93-3	2-Butanone	9.00	U	9.00	30.9	ug/Kg
56-23-5	Carbon Tetrachloride	0.98	U	0.98	6.20	ug/Kg
156-59-2	cis-1,2-Dichloroethene	0.84	U	0.84	6.20	ug/Kg
74-97-5	Bromochloromethane	1.00	U	1.00	6.20	ug/Kg
67-66-3	Chloroform	0.83	U	0.83	6.20	ug/Kg
71-55-6	1,1,1-Trichloroethane	0.93	U	0.93	6.20	ug/Kg
108-87-2	Methylcyclohexane	0.99	U	0.99	6.20	ug/Kg
71-43-2	Benzene	0.82	U	0.82	6.20	ug/Kg
107-06-2	1,2-Dichloroethane	1.00	U	1.00	6.20	ug/Kg
79-01-6	Trichloroethene	0.90	U	0.90	6.20	ug/Kg
78-87-5	1,2-Dichloropropane	0.80	U	0.80	6.20	ug/Kg
75-27-4	Bromodichloromethane	0.87	U	0.87	6.20	ug/Kg
108-10-1	4-Methyl-2-Pentanone	5.70	U	5.70	30.9	ug/Kg
108-88-3	Toluene	0.78	U	0.78	6.20	ug/Kg
10061-02-6	t-1,3-Dichloropropene	0.92	U	0.92	6.20	ug/Kg
10061-01-5	cis-1,3-Dichloropropene	0.88	U	0.88	6.20	ug/Kg

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79-00-5	1,1,2-Trichloroethane	1.10	U	1.10	6.20	ug/Kg
591-78-6	2-Hexanone	5.80	U	5.80	30.9	ug/Kg
124-48-1	Dibromochloromethane	0.93	U	0.93	6.20	ug/Kg
106-93-4	1,2-Dibromoethane	0.93	U	0.93	6.20	ug/Kg
127-18-4	Tetrachloroethene	0.94	U	0.94	6.20	ug/Kg
108-90-7	Chlorobenzene	0.80	U	0.80	6.20	ug/Kg
100-41-4	Ethyl Benzene	0.87	U	0.87	6.20	ug/Kg
179601-23-1	m/p-Xylenes	1.80	U	1.80	12.4	ug/Kg
95-47-6	o-Xylene	0.98	U	0.98	6.20	ug/Kg
100-42-5	Styrene	0.98	U	0.98	6.20	ug/Kg
75-25-2	Bromoform	1.00	U	1.00	6.20	ug/Kg
98-82-8	Isopropylbenzene	0.89	U	0.89	6.20	ug/Kg
79-34-5	1,1,2,2-Tetrachloroethane	1.40	U	1.40	6.20	ug/Kg
541-73-1	1,3-Dichlorobenzene	0.83	U	0.83	6.20	ug/Kg
106-46-7	1,4-Dichlorobenzene	0.78	U	0.78	6.20	ug/Kg
95-50-1	1,2-Dichlorobenzene	0.79	U	0.79	6.20	ug/Kg
96-12-8	1,2-Dibromo-3-Chloropropane	1.50	U	1.50	6.20	ug/Kg
120-82-1	1,2,4-Trichlorobenzene	1.20	U	1.20	6.20	ug/Kg
87-61-6	1,2,3-Trichlorobenzene	1.20	U	1.20	6.20	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	74.0		50 - 163	148%	SPK: 50
1868-53-7	Dibromofluoromethane	54.8		54 - 147	110%	SPK: 50
2037-26-5	Toluene-d8	54.3		58 - 134	109%	SPK: 50
460-00-4	4-Bromofluorobenzene	60.4		30 - 143	121%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	41900	7.881			
540-36-3	1,4-Difluorobenzene	88400	8.781			
3114-55-4	Chlorobenzene-d5	88300	11.586			
3855-82-1	1,4-Dichlorobenzene-d4	40700	13.522			
TENTATIVE IDENTIFIED COMPOUNDS						
000111-65-9	Octane	6.70	J		10.5	ug/Kg

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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