

Report of Analysis

Client:	JACOBS Engineering Group, Inc.	Date Collected:	02/28/25
Project:	Former Schlumberger STC PTC Site # D3868221	Date Received:	02/28/25
Client Sample ID:	IDW-AQ-IW-02-COMP-022825	SDG No.:	Q1478
Lab Sample ID:	Q1478-05	Matrix:	Water
Analytical Method:	SW8270	% Solid:	0
Sample Wt/Vol:	940 Units: mL	Final Vol:	1000 uL
Soil Aliquot Vol:	uL	Test:	SVOC-TCL BNA -20
Extraction Type :	Decanted : N	Level :	LOW
Injection Volume :	GPC Factor : 1.0	GPC Cleanup :	N PH :
Prep Method :	SW3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BF141869.D	1	03/05/25 08:35	03/06/25 15:30	PB166982

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
100-52-7	Benzaldehyde	4.30	U	4.30	10.6	ug/L
108-95-2	Phenol	0.99	U	0.99	5.30	ug/L
111-44-4	bis(2-Chloroethyl)ether	1.30	U	1.30	5.30	ug/L
95-57-8	2-Chlorophenol	0.76	U	0.76	5.30	ug/L
95-48-7	2-Methylphenol	1.20	U	1.20	5.30	ug/L
108-60-1	2,2-oxybis(1-Chloropropane)	1.40	U	1.40	5.30	ug/L
98-86-2	Acetophenone	1.20	U	1.20	5.30	ug/L
65794-96-9	3+4-Methylphenols	1.20	U	1.20	10.6	ug/L
621-64-7	n-Nitroso-di-n-propylamine	1.60	U	1.60	2.70	ug/L
67-72-1	Hexachloroethane	1.10	U	1.10	5.30	ug/L
98-95-3	Nitrobenzene	1.40	U	1.40	5.30	ug/L
78-59-1	Isophorone	1.20	U	1.20	5.30	ug/L
88-75-5	2-Nitrophenol	2.10	U	2.10	5.30	ug/L
105-67-9	2,4-Dimethylphenol	1.60	U	1.60	5.30	ug/L
111-91-1	bis(2-Chloroethoxy)methane	1.10	U	1.10	5.30	ug/L
120-83-2	2,4-Dichlorophenol	0.94	U	0.94	5.30	ug/L
91-20-3	Naphthalene	1.10	U	1.10	5.30	ug/L
106-47-8	4-Chloroaniline	1.40	UQ	1.40	5.30	ug/L
87-68-3	Hexachlorobutadiene	1.40	U	1.40	5.30	ug/L
105-60-2	Caprolactam	1.80	U	1.80	10.6	ug/L
59-50-7	4-Chloro-3-methylphenol	0.89	U	0.89	5.30	ug/L
91-57-6	2-Methylnaphthalene	1.20	U	1.20	5.30	ug/L
77-47-4	Hexachlorocyclopentadiene	5.30	U	5.30	10.6	ug/L
88-06-2	2,4,6-Trichlorophenol	0.95	U	0.95	5.30	ug/L
95-95-4	2,4,5-Trichlorophenol	1.10	U	1.10	5.30	ug/L
92-52-4	1,1-Biphenyl	0.97	U	0.97	5.30	ug/L
91-58-7	2-Chloronaphthalene	1.00	U	1.00	5.30	ug/L
88-74-4	2-Nitroaniline	1.50	U	1.50	5.30	ug/L
131-11-3	Dimethylphthalate	0.99	U	0.99	5.30	ug/L

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208-96-8	Acenaphthylene	1.10	U	1.10	5.30	ug/L
606-20-2	2,6-Dinitrotoluene	1.30	U	1.30	5.30	ug/L
99-09-2	3-Nitroaniline	1.50	UQ	1.50	5.30	ug/L
83-32-9	Acenaphthene	0.86	U	0.86	5.30	ug/L
51-28-5	2,4-Dinitrophenol	6.80	U	6.80	10.6	ug/L
100-02-7	4-Nitrophenol	2.10	U	2.10	10.6	ug/L
132-64-9	Dibenzofuran	0.99	U	0.99	5.30	ug/L
121-14-2	2,4-Dinitrotoluene	1.60	U	1.60	5.30	ug/L
84-66-2	Diethylphthalate	1.10	U	1.10	5.30	ug/L
7005-72-3	4-Chlorophenyl-phenylether	1.00	U	1.00	5.30	ug/L
86-73-7	Fluorene	1.00	U	1.00	5.30	ug/L
100-01-6	4-Nitroaniline	2.20	U	2.20	5.30	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	3.30	U	3.30	10.6	ug/L
86-30-6	n-Nitrosodiphenylamine	0.95	U	0.95	5.30	ug/L
101-55-3	4-Bromophenyl-phenylether	1.00	U	1.00	5.30	ug/L
118-74-1	Hexachlorobenzene	1.20	U	1.20	5.30	ug/L
1912-24-9	Atrazine	1.30	U	1.30	5.30	ug/L
87-86-5	Pentachlorophenol	2.00	U	2.00	10.6	ug/L
85-01-8	Phenanthrene	0.95	U	0.95	5.30	ug/L
120-12-7	Anthracene	1.10	U	1.10	5.30	ug/L
86-74-8	Carbazole	1.20	U	1.20	5.30	ug/L
84-74-2	Di-n-butylphthalate	1.60	U	1.60	5.30	ug/L
206-44-0	Fluoranthene	1.40	U	1.40	5.30	ug/L
129-00-0	Pyrene	1.10	U	1.10	5.30	ug/L
85-68-7	Butylbenzylphthalate	2.20	U	2.20	5.30	ug/L
91-94-1	3,3-Dichlorobenzidine	1.40	UQ	1.40	10.6	ug/L
56-55-3	Benzo(a)anthracene	1.00	U	1.00	5.30	ug/L
218-01-9	Chrysene	0.91	U	0.91	5.30	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	2.00	U	2.00	5.30	ug/L
117-84-0	Di-n-octyl phthalate	2.70	U	2.70	10.6	ug/L
205-99-2	Benzo(b)fluoranthene	1.20	U	1.20	5.30	ug/L

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207-08-9	Benzo(k)fluoranthene	1.30	U	1.30	5.30	ug/L
50-32-8	Benzo(a)pyrene	1.80	U	1.80	5.30	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	1.10	U	1.10	5.30	ug/L
53-70-3	Dibenzo(a,h)anthracene	1.20	U	1.20	5.30	ug/L
191-24-2	Benzo(g,h,i)perylene	1.30	U	1.30	5.30	ug/L
95-94-3	1,2,4,5-Tetrachlorobenzene	1.20	U	1.20	5.30	ug/L
123-91-1	1,4-Dioxane	1.30	U	1.30	5.30	ug/L
58-90-2	2,3,4,6-Tetrachlorophenol	0.84	U	0.84	5.30	ug/L

SURROGATES

367-12-4	2-Fluorophenol	75.9		15 (10) - 110 (139)	51%	SPK: 150
13127-88-3	Phenol-d6	44.8		15 (10) - 110 (134)	30%	SPK: 150
4165-60-0	Nitrobenzene-d5	141	*	30 (49) - 130 (133)	141%	SPK: 100
321-60-8	2-Fluorobiphenyl	120		30 (52) - 130 (132)	120%	SPK: 100
118-79-6	2,4,6-Tribromophenol	222	*	15 (44) - 110 (137)	148%	SPK: 150
1718-51-0	Terphenyl-d14	109		30 (48) - 130 (125)	109%	SPK: 100

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	149000	6.886
1146-65-2	Naphthalene-d8	573000	8.163
15067-26-2	Acenaphthene-d10	305000	9.916
1517-22-2	Phenanthrene-d10	526000	11.404
1719-03-5	Chrysene-d12	408000	14.039
1520-96-3	Perylene-d12	419000	15.509

TENTATIVE IDENTIFIED COMPOUNDS

000994-05-8	Butane, 2-methoxy-2-methyl-	140	J	2.16	ug/L
000079-01-6	Trichloroethylene	34.2	J	2.53	ug/L
000057-55-6	Propylene Glycol	4.00	J	3.38	ug/L
000123-42-2	2-Pentanone, 4-hydroxy-4-methyl-	7.10	A	5.10	ug/L
000112-34-5	Ethanol, 2-(2-butoxyethoxy)-	3.70	J	8.06	ug/L
000119-61-9	Benzophenone	5.10	J	10.6	ug/L
000115-96-8	Tri(2-chloroethyl) phosphate	9.70	J	11.1	ug/L

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000057-10-3	n-Hexadecanoic acid	8.20	J		11.9	ug/L
000057-11-4	Octadecanoic acid	3.40	J		12.7	ug/L
007683-64-9	Supraene	2.40	J		14.9	ug/L

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