

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072125\
 Data File : VN087378.D
 Acq On : 21 Jul 2025 14:38
 Operator : JC\MD
 Sample : Q2649-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-1

Quant Time: Jul 22 03:08:18 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071625W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 17 02:56:13 2025
 Response via : Initial Calibration

Compound	R.T. QIon		Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.206	168	162721	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.088	114	334567	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.847	117	329066	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	149766	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.565	65	159643	57.820	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	115.640%
35) Dibromofluoromethane	8.153	113	117660	50.983	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.960%
50) Toluene-d8	10.547	98	440507	53.509	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	107.020%
62) 4-Bromofluorobenzene	12.829	95	151566	49.833	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	99.660%
Target Compounds						
					Qvalue	
16) Acetone	4.424	43	46099	35.610	ug/l	90
18) Methyl Acetate	5.030	43	27161	8.351	ug/l	99
20) Methylene Chloride	5.277	84	8559	3.313	ug/l #	82
37) Ethyl Acetate	7.559	43	12292	2.791	ug/l #	75
43) Isopropyl Acetate	8.677	43	9074	1.327	ug/l #	87

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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