

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061025\
 Data File : VN086933.D
 Acq On : 10 Jun 2025 16:48
 Operator : JC\MD
 Sample : Q2251-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-VPB-182-EB-20250604

Quant Time: Jun 11 01:57:47 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 07 02:12:50 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	207035	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	385111	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	340856	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	162809	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	139054	50.164	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 100.320%		
35) Dibromofluoromethane	8.177	113	114361	50.109	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 100.220%		
50) Toluene-d8	10.570	98	469495	51.965	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 103.920%		
62) 4-Bromofluorobenzene	12.847	95	166962	49.740	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	= 99.480%		
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	88685	60.330	ug/l	100
20) Methylene Chloride	5.300	84	4441	1.614	ug/l #	83
25) 2-Butanone	7.500	43	23219	9.717	ug/l	95
29) Tetrahydrofuran	7.847	42	5759	3.700	ug/l	96

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

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