

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070125\
 Data File : VN087256.D
 Acq On : 01 Jul 2025 12:39
 Operator : JC\MD
 Sample : Q2459-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-1

Quant Time: Jul 02 02:01:35 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N063025W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 01 03:15:00 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	176485	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	329684	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	313203	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	155219	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	122060	55.106	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 110.220%		
35) Dibromofluoromethane	8.177	113	108309	51.307	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 102.620%		
50) Toluene-d8	10.565	98	409435	49.760	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 99.520%		
62) 4-Bromofluorobenzene	12.847	95	143945	49.167	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	= 98.340%		
Target Compounds						
						Qvalue
16) Acetone	4.465	43	17198	20.309	ug/l	93
18) Methyl Acetate	5.047	43	17277	8.912	ug/l	98
20) Methylene Chloride	5.300	84	8221	4.382	ug/l	95
25) 2-Butanone	7.500	43	7203	5.443	ug/l #	78
37) Ethyl Acetate	7.577	43	11941	3.869	ug/l #	89
43) Isopropyl Acetate	8.694	43	7098	1.483	ug/l #	93
84) 1,2,4-Trimethylbenzene	13.476	105	24829	2.965	ug/l	99

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

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