

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061225\
 Data File : VN086976.D
 Acq On : 12 Jun 2025 14:45
 Operator : JC\MD
 Sample : Q2280-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RT-4643

Quant Time: Jun 13 01:28:28 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	187546	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	366227	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	331625	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	155104	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	136177	54.231	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 108.460%	
35) Dibromofluoromethane	8.177	113	111986	51.598	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 103.200%	
50) Toluene-d8	10.571	98	451690	52.572	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 105.140%	
62) 4-Bromofluorobenzene	12.847	95	159349	49.920	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 99.840%	
Target Compounds						
					Qvalue	
16) Acetone	4.459	43	32636	24.508	ug/l	99
18) Methyl Acetate	5.047	43	27184	7.005	ug/l	100
20) Methylene Chloride	5.306	84	9389	3.766	ug/l	99
37) Ethyl Acetate	7.577	43	16833	4.089	ug/l	98
43) Isopropyl Acetate	8.694	43	11144	1.687	ug/l	93

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

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