

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080823\
 Data File : VN078800.D
 Acq On : 08 Aug 2023 14:24
 Operator : JC\MD
 Sample : 03919-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-36

Quant Time: Aug 09 02:48:47 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080723W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 08 05:03:52 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	183990	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	352036	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	320271	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	83728	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	142816	51.777	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 103.560%		
35) Dibromofluoromethane	8.177	113	115634	48.640	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 97.280%		
50) Toluene-d8	10.571	98	442589	50.590	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 101.180%		
62) 4-Bromofluorobenzene	12.853	95	121297	46.392	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 92.780%		
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	67780	63.866	ug/l #	83
20) Methylene Chloride	5.283	84	8132	3.094	ug/l #	76
25) 2-Butanone	7.500	43	31321	17.428	ug/l #	83
37) Ethyl Acetate	7.577	43	35208	8.978	ug/l #	90
43) Isopropyl Acetate	8.700	43	161661	24.452	ug/l #	77
51) 4-Methyl-2-Pentanone	10.447	43	115310	28.474	ug/l #	80

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

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