

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022823\
 Data File : VN076835.D
 Acq On : 28 Feb 2023 18:09
 Operator : JC\MD
 Sample : 01703-06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T2-STONE

Quant Time: Mar 01 01:31:32 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021323W.M
 Quant Title : SW846 8260
 QLast Update : Tue Feb 14 01:44:47 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	264925	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.107	114	498461	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.866	117	477946	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.795	152	185415	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	195461	53.431	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 106.860%	
35) Dibromofluoromethane	8.172	113	153139	46.913	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 93.820%	
50) Toluene-d8	10.571	98	637249	52.273	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 104.540%	
62) 4-Bromofluorobenzene	12.848	95	197119	48.820	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 97.640%	
Target Compounds						
						Qvalue
16) Acetone	4.442	43	212868	130.267	ug/l	96
17) Carbon Disulfide	4.731	76	27800	3.850	ug/l #	93
20) Methylene Chloride	5.284	84	29667	8.663	ug/l	95
25) 2-Butanone	7.495	43	35884	14.887	ug/l #	84
37) Ethyl Acetate	7.566	43	52689	10.618	ug/l	97
43) Isopropyl Acetate	8.689	43	483717	59.905	ug/l #	72
51) 4-Methyl-2-Pentanone	10.448	43	27150	5.285	ug/l	88

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

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