

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082323\
 Data File : VN079019.D
 Acq On : 23 Aug 2023 16:24
 Operator : JC\MD
 Sample : 04091-17
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-G

Quant Time: Aug 24 01:17:27 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 16 02:42:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	181131	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	353116	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	353014	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	111528	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	148728	52.768	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 105.540%	
35) Dibromofluoromethane	8.177	113	124266	52.279	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.560%	
50) Toluene-d8	10.571	98	451198	51.122	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 102.240%	
62) 4-Bromofluorobenzene	12.853	95	140562	48.724	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 97.440%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	25208	21.675	ug/l #	86
20) Methylene Chloride	5.283	84	12599	5.084	ug/l #	83
25) 2-Butanone	7.494	43	9856	5.286	ug/l #	56
43) Isopropyl Acetate	8.694	43	207657	29.367	ug/l #	71
51) 4-Methyl-2-Pentanone	10.447	43	39593	9.441	ug/l #	74

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

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