

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082323\
 Data File : VN079017.D
 Acq On : 23 Aug 2023 15:36
 Operator : JC\MD
 Sample : 04091-11
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-E

Quant Time: Aug 24 01:12:32 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	8.229	168	184340	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	358750	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	357916	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	107478	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	149993	52.290	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 104.580%	
35) Dibromofluoromethane	8.171	113	123431	51.112	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.220%	
50) Toluene-d8	10.570	98	450778	50.272	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 100.540%	
62) 4-Bromofluorobenzene	12.853	95	143319	48.900	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 97.800%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	27674	23.382	ug/l	92
20) Methylene Chloride	5.283	84	13245	5.251	ug/l #	93
25) 2-Butanone	7.494	43	9831	5.181	ug/l #	63
43) Isopropyl Acetate	8.694	43	168570	23.465	ug/l #	74
51) 4-Methyl-2-Pentanone	10.447	43	36871	8.654	ug/l #	81

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

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