

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082123\
 Data File : VN078977.D
 Acq On : 21 Aug 2023 16:09
 Operator : JC\MD
 Sample : 04082-10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 362

Quant Time: Aug 22 00:58:43 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	190367	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	359078	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	357338	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	112726	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	153031	51.660	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 103.320%	
35) Dibromofluoromethane	8.171	113	124461	51.491	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.980%	
50) Toluene-d8	10.571	98	457952	51.025	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 102.060%	
62) 4-Bromofluorobenzene	12.853	95	142634	48.621	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 97.240%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	38571	31.557	ug/l	88
20) Methylene Chloride	5.289	84	9514	3.653	ug/l #	84
25) 2-Butanone	7.500	43	10551	5.384	ug/l #	63
43) Isopropyl Acetate	8.694	43	170665	23.735	ug/l #	76
51) 4-Methyl-2-Pentanone	10.453	43	40466	9.489	ug/l #	76

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

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