

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082123\
 Data File : VN078973.D
 Acq On : 21 Aug 2023 14:32
 Operator : JC\MD
 Sample : 04082-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 277

Quant Time: Aug 22 00:57:15 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	8.235	168	190398	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	354995	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	331770	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	98856	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	149536	50.472	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 100.940%	
35) Dibromofluoromethane	8.171	113	125912	52.691	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 105.380%	
50) Toluene-d8	10.571	98	440314	49.624	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 99.240%	
62) 4-Bromofluorobenzene	12.853	95	126730	43.697	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 87.400%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	33672	27.544	ug/l #	86
20) Methylene Chloride	5.289	84	10744	4.124	ug/l	88
25) 2-Butanone	7.500	43	10526	5.371	ug/l #	62
43) Isopropyl Acetate	8.700	43	131277	18.467	ug/l #	73
51) 4-Methyl-2-Pentanone	10.447	43	37040	8.785	ug/l #	83

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

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