

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082123\
 Data File : VN078975.D
 Acq On : 21 Aug 2023 15:20
 Operator : JC\MD
 Sample : 04082-06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 72-11929

Quant Time: Aug 22 00:57:58 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	8.236	168	195088	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	374703	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	359467	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	113616	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	152996	50.398	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	100.800%
35) Dibromofluoromethane	8.177	113	125658	49.819	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.640%
50) Toluene-d8	10.571	98	466472	49.807	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.620%
62) 4-Bromofluorobenzene	12.853	95	143316	46.817	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.640%
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	30375	24.250	ug/l	89
20) Methylene Chloride	5.283	84	11827	4.431	ug/l #	84
25) 2-Butanone	7.500	43	11554	5.753	ug/l #	82
43) Isopropyl Acetate	8.694	43	153358	20.439	ug/l #	73
51) 4-Methyl-2-Pentanone	10.453	43	40294	9.054	ug/l #	79

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

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