

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN092123\
 Data File : VN079320.D
 Acq On : 21 Sep 2023 12:59
 Operator : JC\MD
 Sample : 04501-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-3

Quant Time: Sep 22 02:29:48 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091123W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 12 05:05:24 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	175575	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	327343	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	301337	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	82057	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	142087	59.950	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	119.900%
35) Dibromofluoromethane	8.171	113	110429	50.715	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.440%
50) Toluene-d8	10.571	98	396277	49.565	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.120%
62) 4-Bromofluorobenzene	12.853	95	121294	47.561	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	95.120%
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	21442	21.302	ug/l #	85
20) Methylene Chloride	5.277	84	15832	6.463	ug/l #	82
25) 2-Butanone	7.494	43	12388	8.482	ug/l #	77
43) Isopropyl Acetate	8.694	43	108585	19.879	ug/l #	81
51) 4-Methyl-2-Pentanone	10.447	43	27159	8.565	ug/l #	78

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

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