

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN092123\
 Data File : VN079321.D
 Acq On : 21 Sep 2023 13:23
 Operator : JC\MD
 Sample : 04501-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-2

Quant Time: Sep 22 02:30:19 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	162546	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	312469	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	284859	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	77603	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	130323	59.394	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	118.780%
35) Dibromofluoromethane	8.177	113	101361	48.767	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.540%
50) Toluene-d8	10.571	98	367536	48.158	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	96.320%
62) 4-Bromofluorobenzene	12.853	95	113433	46.596	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.200%
Target Compounds						
						Qvalue
16) Acetone	4.442	43	30816	33.068	ug/l	93
20) Methylene Chloride	5.277	84	15515	6.841	ug/l #	81
25) 2-Butanone	7.494	43	15668	11.588	ug/l	88
43) Isopropyl Acetate	8.694	43	132551	25.422	ug/l #	79
51) 4-Methyl-2-Pentanone	10.447	43	29680	9.806	ug/l #	79

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

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