

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN092123\
 Data File : VN079319.D
 Acq On : 21 Sep 2023 12:23
 Operator : JC\MD
 Sample : 04501-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-1

Quant Time: Sep 22 02:29:11 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	158557	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	296309	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	275588	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	75167	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	135703	63.402	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 126.800%#	
35) Dibromofluoromethane	8.171	113	103852	52.690	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 105.380%	
50) Toluene-d8	10.571	98	381988	52.781	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 105.560%	
62) 4-Bromofluorobenzene	12.853	95	115492	50.029	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 100.060%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	23293	25.624	ug/l	91
20) Methylene Chloride	5.277	84	14983	6.773	ug/l #	75
25) 2-Butanone	7.494	43	12916	9.793	ug/l #	73
37) Ethyl Acetate	7.565	43	4467	1.474	ug/l #	75
43) Isopropyl Acetate	8.688	43	114035	23.064	ug/l #	80
51) 4-Methyl-2-Pentanone	10.447	43	27949	9.738	ug/l #	82

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

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