

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN092123\
 Data File : VN079322.D
 Acq On : 21 Sep 2023 13:46
 Operator : JC\MD
 Sample : 04501-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-1

Quant Time: Sep 22 02:30:52 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	151682	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	295862	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	269587	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	72643	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	126812	61.933	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 123.860%	
35) Dibromofluoromethane	8.171	113	98450	50.025	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 100.040%	
50) Toluene-d8	10.571	98	356683	49.359	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 98.720%	
62) 4-Bromofluorobenzene	12.853	95	110311	47.857	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 95.720%	
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	23867	27.446	ug/l	91
20) Methylene Chloride	5.283	84	16018	7.569	ug/l #	87
25) 2-Butanone	7.500	43	12651	10.027	ug/l #	76
43) Isopropyl Acetate	8.694	43	142405	28.845	ug/l #	76
51) 4-Methyl-2-Pentanone	10.447	43	29992	10.465	ug/l #	79

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

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