

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071423\
 Data File : VN078520.D
 Acq On : 14 Jul 2023 14:22
 Operator : JC\MD
 Sample : 03599-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 IDW-7-11-23

Quant Time: Jul 15 04:01:48 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 10 16:13:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	400285	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	735916	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	699569	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	240779	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	273925	51.046	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 102.100%		
35) Dibromofluoromethane	8.171	113	273217	54.691	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 109.380%		
50) Toluene-d8	10.576	98	891568	47.243	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 94.480%		
62) 4-Bromofluorobenzene	12.853	95	298308	51.318	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 102.640%		
Target Compounds						
						Qvalue
16) Acetone	4.447	43	125226	74.914	ug/l	93
20) Methylene Chloride	5.294	84	27816	4.693	ug/l	98
25) 2-Butanone	7.500	43	78212	29.724	ug/l #	83
37) Ethyl Acetate	7.577	43	18388	3.058	ug/l #	86
43) Isopropyl Acetate	8.700	43	203281	20.534	ug/l #	82
51) 4-Methyl-2-Pentanone	10.453	43	251697	42.744	ug/l	91

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

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