

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062223\
 Data File : VN078359.D
 Acq On : 22 Jun 2023 18:52
 Operator : JC\MD
 Sample : 03357-20
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-SCAN-04-G

Quant Time: Jun 23 08:04:07 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060623W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 07 05:05:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	441923	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	850091	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	770370	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	233386	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	323704	54.938	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 109.880%		
35) Dibromofluoromethane	8.177	113	299310	55.685	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 111.380%		
50) Toluene-d8	10.571	98	1016563	49.059	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 98.120%		
62) 4-Bromofluorobenzene	12.853	95	307305	47.705	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 95.420%		
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	270323	159.390	ug/l	98
20) Methylene Chloride	5.300	84	13894	2.479	ug/l	93
25) 2-Butanone	7.500	43	55016	21.095	ug/l #	84
37) Ethyl Acetate	7.577	43	18855	3.131	ug/l #	94
43) Isopropyl Acetate	8.700	43	124942	10.847	ug/l #	76
51) 4-Methyl-2-Pentanone	10.453	43	327022	53.915	ug/l	90

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

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