

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062223\
 Data File : VN078357.D
 Acq On : 22 Jun 2023 18:04
 Operator : JC\MD
 Sample : 03360-06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CE-88N-062023

Quant Time: Jun 23 08:03:47 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	465916	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	872501	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	823142	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	262757	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	328869	52.940	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 105.880%	
35) Dibromofluoromethane	8.177	113	303182	54.957	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 109.920%	
50) Toluene-d8	10.571	98	1036848	48.753	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 97.500%	
62) 4-Bromofluorobenzene	12.853	95	341480	51.649	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 103.300%	
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	68682	38.411	ug/l #	79
20) Methylene Chloride	5.289	84	22863	3.869	ug/l #	89
25) 2-Butanone	7.506	43	19164	6.970	ug/l #	70
37) Ethyl Acetate	7.577	43	53902	8.720	ug/l	97
43) Isopropyl Acetate	8.694	43	282754	23.917	ug/l #	82
51) 4-Methyl-2-Pentanone	10.447	43	147007	23.614	ug/l	87

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

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