

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062322\
 Data File : VN073052.D
 Acq On : 23 Jun 2022 16:40
 Operator : JC\MD
 Sample : N3381-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-7A

Quant Time: Jun 24 01:19:36 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 18 06:22:45 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	323887	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	517247	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	473648	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	233467	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	191184	43.661	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	87.320%
35) Dibromofluoromethane	7.951	113	161532	50.525	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.060%
50) Toluene-d8	10.374	98	533091	44.811	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	89.620%
62) 4-Bromofluorobenzene	12.668	95	236734	55.341	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	110.680%
Target Compounds						
					Qvalue	
16) Acetone	4.234	43	90614	37.645	ug/l	91
18) Methyl Acetate	4.792	43	4869	0.760	ug/l #	72
25) 2-Butanone	7.286	43	5576	1.460	ug/l #	72
31) Cyclohexane	8.022	56	27569	4.654	ug/l #	80
40) Benzene	8.392	78	352882	25.956	ug/l	100
52) Toluene	10.439	92	8197	0.982	ug/l	100
73) Isopropylbenzene	12.515	105	10540	0.584	ug/l	99
78) n-propylbenzene	12.857	91	21523	1.083	ug/l	100

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

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