

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062222\
 Data File : VN073031.D
 Acq On : 22 Jun 2022 17:42
 Operator : JC\MD
 Sample : N3381-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-7A

Quant Time: Jun 23 07:33:15 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	284793	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	464864	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	414474	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	198406	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	171138	44.448	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	88.900%
35) Dibromofluoromethane	7.951	113	143969	50.106	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	100.220%
50) Toluene-d8	10.375	98	473777	44.313	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	88.620%
62) 4-Bromofluorobenzene	12.669	95	201231	52.342	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	104.680%
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	83750	39.570	ug/l	97
25) 2-Butanone	7.275	43	5306	1.580	ug/l #	78
31) Cyclohexane	8.022	56	26178	5.026	ug/l #	88
39) Methylcyclohexane	9.398	83	2677	0.571	ug/l #	74
40) Benzene	8.392	78	350292	28.669	ug/l	97
52) Toluene	10.445	92	11609	1.548	ug/l	99
67) Ethyl Benzene	11.780	91	3294	0.231	ug/l	94
68) m/p-Xylenes	11.892	106	5322	0.952	ug/l	75
69) o-Xylene	12.216	106	1416	0.254	ug/l	72
73) Isopropylbenzene	12.516	105	8376	0.546	ug/l	96
78) n-propylbenzene	12.857	91	18502	1.096	ug/l	98
95) Naphthalene	15.433	128	5497	0.359	ug/l #	97

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

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