

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072022\
 Data File : VN073542.D
 Acq On : 20 Jul 2022 23:27
 Operator : JC\MD
 Sample : N3779-10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 070622-H

Quant Time: Jul 21 08:28:28 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 29 01:54:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	277941	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	447532	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	424840	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	229091	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	170186	42.415	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	84.840%
35) Dibromofluoromethane	7.951	113	143668	48.251	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.500%
50) Toluene-d8	10.381	98	480238	45.412	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	90.820%
62) 4-Bromofluorobenzene	12.675	95	213858	52.623	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	105.240%
Target Compounds						
					Qvalue	
16) Acetone	4.222	43	146707	65.139	ug/l	95
25) 2-Butanone	7.269	43	25025	6.979	ug/l #	85
30) Chloroform	7.740	83	17474	2.607	ug/l	92
40) Benzene	8.387	78	93717	6.628	ug/l	95
51) 4-Methyl-2-Pentanone	10.275	43	25683	4.071	ug/l #	85
52) Toluene	10.445	92	119796	13.710	ug/l	98
64) Tetrachloroethene	10.910	164	2055	0.592	ug/l #	85
67) Ethyl Benzene	11.786	91	56959	3.265	ug/l	96
68) m/p-Xylenes	11.892	106	170341	25.875	ug/l	97
69) o-Xylene	12.222	106	88719	13.201	ug/l	96
80) 1,3,5-Trimethylbenzene	12.998	105	12496	0.775	ug/l	97
84) 1,2,4-Trimethylbenzene	13.310	105	53054	3.296	ug/l	97
95) Naphthalene	15.433	128	2068331	117.144	ug/l	100

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

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