

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062822\
 Data File : VN073227.D
 Acq On : 29 Jun 2022 03:31
 Operator : JC\MD
 Sample : N3376-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 220615003-01-VOA

Quant Time: Jun 29 04:28:24 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	324333	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	529957	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	504737	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	252638	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	206003	43.998	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	88.000%
35) Dibromofluoromethane	7.951	113	171834	48.735	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.480%
50) Toluene-d8	10.381	98	567362	45.307	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	90.620%
62) 4-Bromofluorobenzene	12.669	95	254901	52.967	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	105.940%

Target Compounds						Qvalue
11) Tert butyl alcohol	5.281	59	1716039	1512.974	ug/l	100
16) Acetone	4.228	43	33223	12.641	ug/l	98
17) Carbon Disulfide	4.481	76	5030	0.508	ug/l	99
25) 2-Butanone	7.240	43	73040	17.456	ug/l #	68
29) Tetrahydrofuran	7.610	42	252545	88.146	ug/l	89
49) 1,4-Dioxane	9.510	88	4729	45.667	ug/l #	87
51) 4-Methyl-2-Pentanone	10.269	43	13162	1.762	ug/l #	72
52) Toluene	10.439	92	7158	0.692	ug/l	95
68) m/p-Xylenes	11.886	106	3315	0.424	ug/l	94
95) Naphthalene	15.433	128	10376	0.533	ug/l	96

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

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