

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062921\
 Data File : VN067555.D
 Acq On : 29 Jun 2021 16:05
 Operator : JC/MD
 Sample : M2849-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 NWB-1613

Manual Integrations
 APPROVED

MMDadoda
 7/5/2021 1:51:44 PM

Quant Time: Jun 29 16:41:28 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062821W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 28 14:31:03 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	167067	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	303690	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	294797	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	114145	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	125453	54.868	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 109.740%			
35) Dibromofluoromethane	8.024	113	92199	48.967	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 97.940%			
50) Toluene-d8	10.443	98	419988	54.910	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 109.820%			
62) 4-Bromofluorobenzene	12.733	95	135300	50.642	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 101.280%			
Target Compounds					Qvalue	
16) Acetone	4.296	43	29884	39.625	ug/l	100
25) 2-Butanone	7.356	43	12609m	10.841	ug/l	
29) Tetrahydrofuran	7.691	42	2977m	3.844	ug/l	
51) 4-Methyl-2-Pentanone	10.338	43	2759	1.063	ug/l	95
52) Toluene	10.510	92	10475	1.931	ug/l	94
68) m/p-Xylenes	11.953	106	14100	3.354	ug/l	83
69) o-Xylene	12.283	106	11605	2.787	ug/l	90
80) 1,3,5-Trimethylbenzene	13.063	105	20437	2.598	ug/l	96
84) 1,2,4-Trimethylbenzene	13.372	105	90962	11.744	ug/l	98
95) Naphthalene	15.507	128	53012	12.001	ug/l	99

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

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