

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN083121\
 Data File : VN068371.D
 Acq On : 31 Aug 2021 15:57
 Operator : JC/MD
 Sample : M3603-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TRE-1242

Quant Time: Sep 01 02:06:54 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N083021W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 31 02:20:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.091	168	394517	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	646060	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	664328	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	339338	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	222338	55.214	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 110.420%			
35) Dibromofluoromethane	8.024	113	204154	49.840	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 99.680%			
50) Toluene-d8	10.446	98	783805	51.116	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 102.240%			
62) 4-Bromofluorobenzene	12.733	95	293506	58.326	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 116.660%			
Target Compounds						Qvalue
16) Acetone	4.299	43	34005	19.091	ug/l	96
29) Tetrahydrofuran	7.691	42	31952	27.008	ug/l	98
40) Benzene	8.464	78	22486	1.417	ug/l	98
52) Toluene	10.507	92	334320	31.318	ug/l	98
67) Ethyl Benzene	11.848	91	240622	11.031	ug/l	99
68) m/p-Xylenes	11.953	106	825102	93.893	ug/l	97
69) o-Xylene	12.283	106	541323	64.032	ug/l	97
78) n-propylbenzene	12.921	91	41533	1.610	ug/l	95
80) 1,3,5-Trimethylbenzene	13.061	105	326877	16.518	ug/l	98
84) 1,2,4-Trimethylbenzene	13.372	105	1657175	85.610	ug/l	99
95) Naphthalene	15.507	128	296174	22.559	ug/l	100

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

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