

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102621\
 Data File : VN069168.D
 Acq On : 25 Oct 2021 22:21
 Operator : JC/MD
 Sample : M4284-08
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 20877

Manual Integrations
 APPROVED

MMDadoda
 10/26/2021 6:15:42 PM

Quant Time: Oct 26 01:37:33 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 13 05:44:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.091	168	372947	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	649367	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	646836	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	295973	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	238528	48.509	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	97.020%		
35) Dibromofluoromethane	8.024	113	190357	50.385	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	100.760%		
50) Toluene-d8	10.443	98	819444	54.208	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	108.420%		
62) 4-Bromofluorobenzene	12.733	95	326902	57.318	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	114.640%		
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	5.369	59	8425	9.377	ug/l #	82
16) Acetone	4.285	43	8202844	4470.225	ug/l	99
17) Carbon Disulfide	4.537	76	13445	1.375	ug/l	95
18) Methyl Acetate	4.859	43	230522	33.120	ug/l	97
20) Methylene Chloride	5.106	84	3809m	0.902	ug/l	
25) 2-Butanone	7.335	43	841088	303.114	ug/l	97
40) Benzene	8.469	78	5170	0.306	ug/l #	87
51) 4-Methyl-2-Pentanone	10.333	43	17385	2.966	ug/l	100
52) Toluene	10.507	92	3938	0.362	ug/l	95
59) 2-Hexanone	11.087	43	202155	48.099	ug/l	99
86) p-Isopropyltoluene	13.619	119	339899	17.130	ug/l	97

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

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