

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN092321\
 Data File : VN068624.D
 Acq On : 23 Sep 2021 17:00
 Operator : JC/MD
 Sample : M3826-07
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 AUD-1182

Manual Integrations
 APPROVED

MMDadoda
 9/24/2021 6:58:56 PM

Quant Time: Sep 23 17:25:46 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N092321W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 23 13:52:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	622440	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	1027475	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	945464	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	358709	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	527182	54.384	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 108.760%			
35) Dibromofluoromethane	8.021	113	361886	52.660	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 105.320%			
50) Toluene-d8	10.443	98	1488904	51.095	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 102.180%			
62) 4-Bromofluorobenzene	12.734	95	474563	47.618	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 95.240%			
Target Compounds					Qvalue	
11) Tert butyl alcohol	5.366	59	41914	28.023	ug/l	98
13) Acrolein	4.052	56	10194m	27.950	ug/l	
16) Acetone	4.285	43	1388195	400.816	ug/l	98
25) 2-Butanone	7.337	43	183782	31.024	ug/l	97
37) Ethyl Acetate	7.415	43	151760	12.881	ug/l	98
51) 4-Methyl-2-Pentanone	10.333	43	86302	7.081	ug/l	99

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

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