

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070921\
 Data File : VN067693.D
 Acq On : 9 Jul 2021 18:17
 Operator : JC/MD
 Sample : M2999-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FRAC-TANK-N43502

Manual Integrations
 APPROVED

MMDadoda
 7/12/2021 5:44:54 PM

Quant Time: Jul 10 05:26:00 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N070721W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jul 07 16:18:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	313902	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	587080	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	543922	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	205483	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	245177	51.681	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 103.360%			
35) Dibromofluoromethane	8.024	113	175755	49.550	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 99.100%			
50) Toluene-d8	10.446	98	749470	50.480	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 100.960%			
62) 4-Bromofluorobenzene	12.733	95	241232	45.995	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 92.000%			
Target Compounds						Qvalue
11) Tert butyl alcohol	5.398	59	4420m	6.484	ug/l	
16) Acetone	4.304	43	5056	3.387	ug/l	96
18) Methyl Acetate	4.862	43	4160m	0.992	ug/l	
30) Chloroform	7.823	83	8348	1.145	ug/l	87

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

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