

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN051821\
 Data File : VN067130.D
 Acq On : 18 May 2021 16:27
 Operator : JC/MD
 Sample : M2387-10
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 30795

Manual Integrations
 APPROVED

MMDadoda
 5/19/2021 4:00:02 PM

Quant Time: May 19 02:50:35 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051421W.M
 Quant Title : SW846 8260
 QLast Update : Tue May 18 04:00:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.597	168	655461	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.525	114	906214	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.357	117	813994	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.296	152	394143	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.959	65	316666	44.72	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	89.44%		
35) Dibromofluoromethane	7.519	113	295482	49.66	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	99.32%		
50) Toluene-d8	10.035	98	1142493	50.66	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	101.32%		
62) 4-Bromofluorobenzene	12.352	95	392960	52.02	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	104.04%		
Target Compounds					Qvalue	
16) Acetone	3.778	43	407003	184.93	ug/l	99
17) Carbon Disulfide	3.998	76	10738	0.65	ug/l #	89
18) Methyl Acetate	4.269	43	118936	18.17	ug/l	99
25) 2-Butanone	6.760	43	61378m	17.88	ug/l	
37) Ethyl Acetate	6.851	43	2526066	345.99	ug/l	98
40) Benzene	7.972	78	8529	0.34	ug/l	94
51) 4-Methyl-2-Pentanone	9.930	43	74675	15.43	ug/l	96
52) Toluene	10.102	92	411780	25.39	ug/l	100
84) 1,2,4-Trimethylbenzene	12.998	105	14463	0.57	ug/l	99
86) p-Isopropyltoluene	13.242	119	92043	3.42	ug/l	98
95) Naphthalene	15.090	128	20019	4.37	ug/l #	96

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

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