

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121521\
 Data File : VN070115.D
 Acq On : 15 Dec 2021 16:44
 Operator : JC/MD
 Sample : M4970-13
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SW-2

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/21/2021
 Supervised By : Mahesh Dadoda 12/21/2021

Quant Time: Dec 16 05:40:06 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 03 08:46:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	1152292	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2180634	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	2580180	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	1076449	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	919062	51.658	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 103.320%			
35) Dibromofluoromethane	8.024	113	642968	47.073	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 94.140%			
50) Toluene-d8	10.443	98	2919713	53.235	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 106.480%			
62) 4-Bromofluorobenzene	12.731	95	1293086	65.079	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 130.160%			
Target Compounds						
16) Acetone	4.293	43	526072	49.155	ug/l	97
19) Methyl tert-butyl Ether	5.629	73	10587	0.248	ug/l #	16
25) 2-Butanone	7.345	43	81679	5.617	ug/l	93
40) Benzene	8.464	78	165028	2.601	ug/l	98
51) 4-Methyl-2-Pentanone	10.336	43	38794	1.308	ug/l	94
52) Toluene	10.505	92	31295	0.808	ug/l	96
86) p-Isopropyltoluene	13.616	119	58069m	0.802	ug/l	
95) Naphthalene	15.512	128	20509	4.344	ug/l #	95

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

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