

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121021\
 Data File : VN069988.D
 Acq On : 11 Dec 2021 5:34
 Operator : JC/MD
 Sample : M4990-12
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 R8-6

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/11/2021
 Supervised By : Amit Patel 12/14/2021

Quant Time: Dec 11 08:31:34 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	1058314	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2098125	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	2454960	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	1022973	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	885417	54.186	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 108.380%			
35) Dibromofluoromethane	8.024	113	644017	49.004	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 98.000%			
50) Toluene-d8	10.443	98	2770505	52.501	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 105.000%			
62) 4-Bromofluorobenzene	12.731	95	1184760	61.972	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 123.940%			
Target Compounds						
16) Acetone	4.296	43	66626	6.778	ug/l	94
18) Methyl Acetate	4.856	43	48344m	1.629	ug/l	
20) Methylene Chloride	5.095	84	53926	4.127	ug/l	86
43) Isopropyl Acetate	8.566	43	155589	3.537	ug/l #	80

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

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