

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051922\
 Data File : VX028839.D
 Acq On : 19 May 2022 20:36
 Operator : JC/MD
 Sample : N2943-11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-30

Quant Time: May 20 01:51:33 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 12 14:58:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	303785	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	544134	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	516097	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	239533	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	227429	54.615	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 109.240%	
35) Dibromofluoromethane	5.391	113	186227	52.236	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 104.480%	
50) Toluene-d8	8.653	98	647125	48.145	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 96.300%	
62) 4-Bromofluorobenzene	11.079	95	254523	50.699	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 101.400%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	4665	2.610	ug/l	98
17) Carbon Disulfide	2.514	76	3880	0.476	ug/l	96
19) Methyl tert-butyl Ether	3.117	73	25877	2.319	ug/l	99
22) Diisopropyl ether	3.764	45	5027	0.466	ug/l #	91
40) Benzene	6.044	78	4635	0.301	ug/l	99

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

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