

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070621\
 Data File : VX023147.D
 Acq On : 06 Jul 2021 15:42
 Operator : JC/MD
 Sample : M2920-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-4

Quant Time: Jul 07 02:52:48 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 02 14:29:42 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	265642	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	437268	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	384748	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	157199	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	181901	50.949	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 101.900%		
35) Dibromofluoromethane	5.397	113	138068	48.896	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 97.800%		
50) Toluene-d8	8.653	98	518279	49.205	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 98.420%		
62) 4-Bromofluorobenzene	11.085	95	184927	45.477	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 90.960%		
Target Compounds						
						Qvalue
8) Diethyl Ether	2.142	74	1711	1.057	ug/l	80
11) Tert butyl alcohol	2.977	59	5033	6.701	ug/l #	94
16) Acetone	2.386	43	3745	2.651	ug/l	98
19) Methyl tert-butyl Ether	3.123	73	4417	0.499	ug/l	97
27) cis-1,2-Dichloroethene	4.501	96	1448	0.467	ug/l	91
29) Tetrahydrofuran	5.031	42	1784	1.303	ug/l #	86
65) Chlorobenzene	10.085	112	2795	0.367	ug/l	95

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

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