

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112321\
 Data File : VX025283.D
 Acq On : 23 Nov 2021 18:38
 Operator : JC/MD
 Sample : M4739-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-33A

Quant Time: Nov 24 03:15:04 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 17 15:36:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	94960	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	169746	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	164735	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	75619	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	56551	56.498	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 113.000%	
35) Dibromofluoromethane	5.391	113	54621	52.444	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.880%	
50) Toluene-d8	8.653	98	199251	50.882	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 101.760%	
62) 4-Bromofluorobenzene	11.085	95	76478	51.624	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 103.240%	
Target Compounds						Qvalue
16) Acetone	2.380	43	1700	3.057	ug/l	94
19) Methyl tert-butyl Ether	3.117	73	19236	6.273	ug/l	100
27) cis-1,2-Dichloroethene	4.495	96	7071	6.280	ug/l	78

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

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