

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120221\
 Data File : VX025456.D
 Acq On : 03 Dec 2021 02:53
 Operator : JC/MD
 Sample : M4917-20
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-36

Quant Time: Dec 03 06:15:10 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	113776	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	191022	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	168741	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	75798	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	65802	54.868	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 109.740%	
35) Dibromofluoromethane	5.391	113	58518	49.928	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.860%	
50) Toluene-d8	8.653	98	220071	49.940	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 99.880%	
62) 4-Bromofluorobenzene	11.085	95	75093	45.044	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 90.080%	
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	1076	1.615	ug/l #	85
19) Methyl tert-butyl Ether	3.117	73	819	0.223	ug/l #	86
40) Benzene	6.044	78	1453	0.288	ug/l	98

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

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