

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080921\
 Data File : VX023617.D
 Acq On : 09 Aug 2021 16:47
 Operator : JC/MD
 Sample : M3278-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-4

Quant Time: Aug 10 04:04:01 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 23 06:05:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	178864	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	301534	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	269250	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	107841	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	119027	53.347	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 106.700%	
35) Dibromofluoromethane	5.397	113	94626	50.191	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 100.380%	
50) Toluene-d8	8.653	98	361805	52.233	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 104.460%	
62) 4-Bromofluorobenzene	11.085	95	121448	51.562	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 103.120%	
Target Compounds						Qvalue
16) Acetone	2.386	43	7789	7.117	ug/l	97
17) Carbon Disulfide	2.514	76	7097	1.638	ug/l	98
30) Chloroform	5.105	83	1852	0.522	ug/l	87

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

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