

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082021\
 Data File : VX023851.D
 Acq On : 19 Aug 2021 20:32
 Operator : JC/MD
 Sample : M3483-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BB-TT-VPB-RW7-GW-260-262

Quant Time: Aug 20 03:16:01 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 17 11:57:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	136495	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	236408	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	216536	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	86114	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.970	65	99799	52.387	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 104.780%	
35) Dibromofluoromethane	5.397	113	77570	49.574	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.140%	
50) Toluene-d8	8.653	98	287807	48.327	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 96.660%	
62) 4-Bromofluorobenzene	11.085	95	94417	43.067	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 86.140%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.325	96	1388	1.060	ug/l	93
16) Acetone	2.392	43	3039	3.568	ug/l #	84
19) Methyl tert-butyl Ether	3.135	73	1479	0.324	ug/l	100
24) 1,1-Dichloroethane	3.617	63	4986	1.863	ug/l	99
30) Chloroform	5.111	83	1065	0.376	ug/l #	79
32) 1,1,1-Trichloroethane	5.391	97	2872	1.184	ug/l #	46
44) Trichloroethene	7.135	130	5546	3.101	ug/l	91

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

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