

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX123121\
 Data File : VX026168.D
 Acq On : 30 Dec 2021 15:13
 Operator : JC/MD
 Sample : M5207-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-3D-122021

Quant Time: Dec 31 07:08:54 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X123021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 30 08:02:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	139968	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	239453	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	208095	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	90013	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	103947	51.334	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 102.660%	
35) Dibromofluoromethane	5.391	113	80584	51.764	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 103.520%	
50) Toluene-d8	8.653	98	279926	48.342	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 96.680%	
62) 4-Bromofluorobenzene	11.085	95	96305	46.719	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 93.440%	
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	2889	2.633	ug/l	94
19) Methyl tert-butyl Ether	3.123	73	1241	0.229	ug/l #	85
21) trans-1,2-Dichloroethene	3.093	96	762	0.451	ug/l #	79
27) cis-1,2-Dichloroethene	4.495	96	22544	11.720	ug/l	93
44) Trichloroethene	7.129	130	9640	4.999	ug/l	90
64) Tetrachloroethene	9.275	164	10446	5.948	ug/l	97

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

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