

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010522\
 Data File : VX026353.D
 Acq On : 06 Jan 2022 03:57
 Operator : JC/MD
 Sample : N1019-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-9

Quant Time: Jan 06 05:06:06 2022
 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	5.556	168	156983	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	265108	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	227306	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	99322	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	114334	50.343	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 100.680%		
35) Dibromofluoromethane	5.391	113	84663	49.121	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 98.240%		
50) Toluene-d8	8.653	98	304664	47.522	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 95.040%		
62) 4-Bromofluorobenzene	11.079	95	112959	49.495	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 99.000%		
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	12533	10.185	ug/l	91
19) Methyl tert-butyl Ether	3.117	73	7959	1.311	ug/l	96
37) Ethyl Acetate	4.721	43	6395m	1.984	ug/l	
44) Trichloroethene	7.129	130	3942	1.846	ug/l	95
64) Tetrachloroethene	9.275	164	1932	1.007	ug/l	95
68) m/p-Xylenes	10.299	106	3913	1.174	ug/l	91

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

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