

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090623\
 Data File : VX037551.D
 Acq On : 06 Sep 2023 13:46
 Operator : JC/MD
 Sample : 04220-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW11

Quant Time: Sep 06 22:45:49 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081723W.M
 Quant Title : SW846 8260
 QLast Update : Fri Aug 18 09:30:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	129922	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	213287	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	188934	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	79998	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	74923	40.000	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	80.000%	
35) Dibromofluoromethane	5.391	113	66265	46.912	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	93.820%	
50) Toluene-d8	8.647	98	248428	48.485	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	96.960%	
62) 4-Bromofluorobenzene	11.079	95	88355	43.879	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	87.760%	
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	2457	3.353	ug/l #	79
19) Methyl tert-butyl Ether	3.117	73	2312	0.456	ug/l #	90
27) cis-1,2-Dichloroethene	4.495	96	1972	1.100	ug/l	77
30) Chloroform	5.099	83	2657	0.889	ug/l	87
44) Trichloroethene	7.129	130	2462	1.509	ug/l	90
64) Tetrachloroethene	9.275	164	2089	1.516	ug/l	93

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

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