

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022124\
 Data File : VX040380.D
 Acq On : 21 Feb 2024 14:42
 Operator : JC/MD
 Sample : P1521-03 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CG-MW-17-CH2

Quant Time: Feb 22 01:08:16 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
 Quant Title : SW846 8260
 QLast Update : Mon Feb 12 00:21:14 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.555	168	71320	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	127954	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	126575	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	62750	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	55565	52.369	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	104.740%
35) Dibromofluoromethane	5.385	113	42742	47.833	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.660%
50) Toluene-d8	8.646	98	158727	47.487	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	94.980%
62) 4-Bromofluorobenzene	11.079	95	61847	45.368	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.740%
Target Compounds						
					Qvalue	
16) Acetone	2.379	43	2061	4.730	ug/l	96
40) Benzene	6.037	78	43768	12.504	ug/l	97
51) 4-Methyl-2-Pentanone	8.573	43	948	0.639	ug/l	99
67) Ethyl Benzene	10.201	91	1749	0.372	ug/l	91

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

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