

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090623\
 Data File : VX037548.D
 Acq On : 06 Sep 2023 12:32
 Operator : JC/MD
 Sample : 04220-01 20X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW1R

Quant Time: Sep 06 22:45:09 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	130338	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	216446	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	190707	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	86227	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	73867	39.311	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	78.620%
35) Dibromofluoromethane	5.385	113	65186	45.474	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	90.940%
50) Toluene-d8	8.647	98	251642	48.395	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.800%
62) 4-Bromofluorobenzene	11.079	95	90735	44.403	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	88.800%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.977	59	88999	201.911	ug/l #	94
16) Acetone	2.392	43	1993	2.711	ug/l #	86
17) Carbon Disulfide	2.507	76	855	0.251	ug/l	100
19) Methyl tert-butyl Ether	3.117	73	40038	7.868	ug/l	97
22) Diisopropyl ether	3.763	45	10205	2.395	ug/l #	69
29) Tetrahydrofuran	5.019	42	5126	7.262	ug/l #	83
49) 1,4-Dioxane	7.683	88	473	10.904	ug/l #	67

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

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