

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070623\
 Data File : VX036434.D
 Acq On : 06 Jul 2023 14:41
 Operator : JC/MD
 Sample : 03510-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-WWS-06-G

Quant Time: Jul 07 01:48:02 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062323W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 23 15:03:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	164847	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	283047	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	257604	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	113480	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	110279	46.670	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	93.340%		
35) Dibromofluoromethane	5.391	113	88147	47.230	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	94.460%		
50) Toluene-d8	8.653	98	342571	49.144	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	98.280%		
62) 4-Bromofluorobenzene	11.079	95	117908	48.021	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	96.040%		
Target Compounds						
					Qvalue	
16) Acetone	2.391	43	234725	232.401	ug/l	92
18) Methyl Acetate	2.708	43	4745	1.142	ug/l	98
20) Methylene Chloride	2.788	84	10176	4.679	ug/l	91
25) 2-Butanone	4.568	43	116636	80.374	ug/l	97
37) Ethyl Acetate	4.714	43	16134	5.690	ug/l	97
43) Isopropyl Acetate	6.342	43	69609	15.389	ug/l	97
51) 4-Methyl-2-Pentanone	8.573	43	453908	155.756	ug/l	93

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

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