

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031623\
 Data File : VX034568.D
 Acq On : 17 Mar 2023 06:48
 Operator : JC/MD
 Sample : 01907-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 33957

Quant Time: Mar 17 07:55:12 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Mar 13 10:29:32 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	237675	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	412392	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	376031	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	176578	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	168907	48.850	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.700%
35) Dibromofluoromethane	5.385	113	129972	47.735	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.460%
50) Toluene-d8	8.647	98	504392	49.623	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.240%
62) 4-Bromofluorobenzene	11.079	95	198128	48.257	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.520%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.953	59	6734	9.352	ug/l #	91
16) Acetone	2.373	43	89638	55.035	ug/l	98
17) Carbon Disulfide	2.514	76	7789	1.064	ug/l	96
18) Methyl Acetate	2.703	43	6545	1.230	ug/l	100
20) Methylene Chloride	2.788	84	11651	3.686	ug/l	95
37) Ethyl Acetate	4.714	43	65498	14.336	ug/l	98
43) Isopropyl Acetate	6.336	43	267588	34.573	ug/l	99

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

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