

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060123\
 Data File : VX035959.D
 Acq On : 01 Jun 2023 15:01
 Operator : JC/MD
 Sample : 02997-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DMA-1103

Quant Time: Jun 02 01:41:29 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 31 12:50:30 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	209742	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	342815	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	297183	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	139610	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	143705	42.193	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	84.380%		
35) Dibromofluoromethane	5.385	113	104253	45.214	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	90.420%		
50) Toluene-d8	8.647	98	396725	47.941	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	95.880%		
62) 4-Bromofluorobenzene	11.079	95	150533	46.081	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	92.160%		
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	33414	23.780	ug/l	96
17) Carbon Disulfide	2.514	76	4433	0.841	ug/l	97
20) Methylene Chloride	2.788	84	2400	0.887	ug/l #	85
25) 2-Butanone	4.574	43	9794	4.870	ug/l	94
30) Chloroform	5.099	83	9270	1.929	ug/l	87
51) 4-Methyl-2-Pentanone	8.574	43	17866	5.199	ug/l	97
52) Toluene	8.720	92	4713	0.812	ug/l	94

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

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