

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060123\
 Data File : VX035960.D
 Acq On : 01 Jun 2023 15:24
 Operator : JC/MD
 Sample : 02997-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DMA-1105

Quant Time: Jun 02 01:41:43 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	5.556	168	200625	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	332382	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	305055	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	140372	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	153285	47.051	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	94.100%
35) Dibromofluoromethane	5.385	113	103525	46.308	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.620%
50) Toluene-d8	8.647	98	401355	50.023	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.040%
62) 4-Bromofluorobenzene	11.079	95	152530	48.158	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.320%
Target Compounds						
					Qvalue	
16) Acetone	2.398	43	13520	10.059	ug/l	99
17) Carbon Disulfide	2.508	76	1648	0.327	ug/l	95
25) 2-Butanone	4.574	43	2443	1.270	ug/l	94
30) Chloroform	5.093	83	10474	2.278	ug/l	85
51) 4-Methyl-2-Pentanone	8.574	43	6826	2.049	ug/l	100
52) Toluene	8.720	92	8120	1.444	ug/l	95

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

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