

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062920\
 Data File : VX017022.D
 Acq On : 29 Jun 2020 21:11
 Operator : JC/SP
 Sample : L3143-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-20200626

Quant Time: Jun 30 06:22:13 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X062920W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 29 17:10:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	254598	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	441467	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	458692	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	238314	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	167271	53.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.58%	
35) Dibromofluoromethane	5.46	113	141616	50.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.42%	
50) Toluene-d8	8.70	98	543863	52.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.34%	
62) 4-Bromofluorobenzene	11.12	95	228729	56.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.32%	
Target Compounds						
6) Chloroethane	1.72	64	2014	1.233	ug/l	91
16) Acetone	2.42	43	23161	19.726	ug/l	99
17) Carbon Disulfide	2.55	76	4259	0.628	ug/l #	89
25) 2-Butanone	4.64	43	7252	3.707	ug/l	99
51) 4-Methyl-2-Pentanone	8.62	43	9900	2.361	ug/l	98
59) 2-Hexanone	9.48	43	953	0.311	ug/l	84

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

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