

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110720\
 Data File : VX019348.D
 Acq On : 06 Nov 2020 17:36
 Operator : JC/MD
 Sample : L4662-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-178-DUP-20201104

Quant Time: Nov 07 02:49:23 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102720W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 27 13:55:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	289190	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	511825	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	479371	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	244563	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	203856	54.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.32%	
35) Dibromofluoromethane	5.46	113	159291	49.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.10%	
50) Toluene-d8	8.70	98	616872	50.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.80%	
62) 4-Bromofluorobenzene	11.12	95	242399	52.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.62%	
Target Compounds						
3) Chloromethane	1.31	50	861	0.320	ug/l	89
16) Acetone	2.42	43	5188	4.032	ug/l	96
19) Methyl tert-butyl Ether	3.17	73	7379	0.760	ug/l #	80
30) Chloroform	5.18	83	1216	0.216	ug/l	86
44) Trichloroethene	7.19	130	1710	0.446	ug/l	97
64) Tetrachloroethene	9.32	164	10625	2.909	ug/l	95

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

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