

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110720\
 Data File : VX019338.D
 Acq On : 06 Nov 2020 13:51
 Operator : JC/MD
 Sample : L4662-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB178-GW-238-240

Quant Time: Nov 07 01:57:59 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	303682	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	531681	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	499035	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	245155	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	209643	53.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.06%	
35) Dibromofluoromethane	5.46	113	163667	49.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.02%	
50) Toluene-d8	8.70	98	644428	51.19	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.38%	
62) 4-Bromofluorobenzene	11.12	95	251896	52.83	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.66%	
Target Compounds						
16) Acetone	2.41	43	8621	6.381	ug/l	98
19) Methyl tert-butyl Ether	3.17	73	7022	0.689	ug/l #	88
30) Chloroform	5.18	83	3031	0.513	ug/l	99
44) Trichloroethene	7.20	130	1940	0.487	ug/l	81
64) Tetrachloroethene	9.32	164	1091	0.287	ug/l #	87

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

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