

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110720\
 Data File : VX019342.D
 Acq On : 06 Nov 2020 15:21
 Operator : JC/MD
 Sample : L4662-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB178-GW-318-320

Quant Time: Nov 07 02:05:36 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	295730	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	522724	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	489479	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	244737	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	209012	54.80	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.60%	
35) Dibromofluoromethane	5.46	113	161631	49.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.46%	
50) Toluene-d8	8.70	98	630341	50.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.86%	
62) 4-Bromofluorobenzene	11.12	95	246532	52.59	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.18%	
Target Compounds						
16) Acetone	2.42	43	5169	3.929	ug/l	90
19) Methyl tert-butyl Ether	3.17	73	6139	0.618	ug/l #	91
24) 1,1-Dichloroethane	3.67	63	1647	0.305	ug/l #	91
30) Chloroform	5.17	83	1672	0.291	ug/l #	83
44) Trichloroethene	7.19	130	1715	0.438	ug/l	85
64) Tetrachloroethene	9.32	164	1335	0.358	ug/l	89

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX110720\
Data File : VX019342.D
Acq On : 06 Nov 2020 15:21
Operator : JC/MD
Sample : L4662-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BP-VPB178-GW-318-320

Quant Time: Nov 07 02:05:36 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X102720W.M
Quant Title : SW846 8260
QLast Update : Tue Oct 27 13:55:07 2020
Response via : Initial Calibration

