

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110720\
 Data File : VX019341.D
 Acq On : 06 Nov 2020 14:59
 Operator : JC/MD
 Sample : L4662-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB178-GW-298-300

Quant Time: Nov 07 02:03:58 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	294465	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	513031	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	487526	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	245474	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	206273	54.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.64%	
35) Dibromofluoromethane	5.46	113	159999	49.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.30%	
50) Toluene-d8	8.70	98	624605	51.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.84%	
62) 4-Bromofluorobenzene	11.12	95	244101	53.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.12%	
Target Compounds						
16) Acetone	2.42	43	4462	3.406	ug/l	91
19) Methyl tert-butyl Ether	3.17	73	7623	0.771	ug/l #	86
30) Chloroform	5.19	83	1182	0.206	ug/l	91
44) Trichloroethene	7.19	130	781	0.203	ug/l	81
64) Tetrachloroethene	9.32	164	998	0.269	ug/l	88

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX110720\
Data File : VX019341.D
Acq On : 06 Nov 2020 14:59
Operator : JC/MD
Sample : L4662-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BP-VPB178-GW-298-300

Quant Time: Nov 07 02:03:58 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

