

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX121020\
 Data File : VX019852.D
 Acq On : 11 Dec 2020 02:44
 Operator : JC/MD
 Sample : L4983-11
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TR1-D-BOTTOM

Quant Time: Dec 11 05:55:42 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X120420W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 04 15:32:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	324632	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	546671	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	508463	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	221062	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	6.03	65	209407	49.08	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	98.16%		
35) Dibromofluoromethane	5.46	113	167987	47.55	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	95.10%		
50) Toluene-d8	8.70	98	670238	49.69	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	99.38%		
62) 4-Bromofluorobenzene	11.12	95	232161	45.22	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	90.44%		

Target Compounds

					Qvalue
16) Acetone	2.42	43	25502	15.902	ug/l 100
20) Methylene Chloride	2.83	84	33351	7.912	ug/l 98
43) Isopropyl Acetate	6.41	43	24090	2.971	ug/l 99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX121020\
Data File : VX019852.D
Acq On : 11 Dec 2020 02:44
Operator : JC/MD
Sample : L4983-11
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TR1-D-BOTTOM

Quant Time: Dec 11 05:55:42 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X120420W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 04 15:32:53 2020
Response via : Initial Calibration

