

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX121020\
 Data File : VX019850.D
 Acq On : 11 Dec 2020 01:57
 Operator : JC/MD
 Sample : L4983-09
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TR1-C-BOTTOM

Quant Time: Dec 11 03:19:03 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	320580	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	535669	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	495168	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	212418	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	204560	48.55	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	97.10%		
35) Dibromofluoromethane	5.46	113	165204	47.72	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	95.44%		
50) Toluene-d8	8.70	98	657744	49.77	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	99.54%		
62) 4-Bromofluorobenzene	11.12	95	229504	45.62	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	91.24%		

Target Compounds					Qvalue
16) Acetone	2.42	43	25486	16.093 ug/l	96
20) Methylene Chloride	2.83	84	34598	8.386 ug/l	98
43) Isopropyl Acetate	6.41	43	24498	3.083 ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX121020\
Data File : VX019850.D
Acq On : 11 Dec 2020 01:57
Operator : JC/MD
Sample : L4983-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TR1-C-BOTTOM

Quant Time: Dec 11 03:19:03 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

