

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX120920\
 Data File : VX019783.D
 Acq On : 09 Dec 2020 16:57
 Operator : JC/MD
 Sample : L4971-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 5S1-C-TOP

Quant Time: Dec 10 06:14:16 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	332623	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	547380	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	492579	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	212507	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	206459	47.23	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	94.46%		
35) Dibromofluoromethane	5.46	113	166139	46.96	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	93.92%		
50) Toluene-d8	8.70	98	664829	49.23	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	98.46%		
62) 4-Bromofluorobenzene	11.12	95	229732	44.69	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	89.38%		

Target Compounds					Qvalue
16) Acetone	2.42	43	13850	8.429 ug/l	100
20) Methylene Chloride	2.84	84	11843	1.786 ug/l	95
43) Isopropyl Acetate	6.41	43	25877	3.187 ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120920\
Data File : VX019783.D
Acq On : 09 Dec 2020 16:57
Operator : JC/MD
Sample : L4971-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
5S1-C-TOP

Quant Time: Dec 10 06:14:16 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

