

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110918\
 Data File : VX005865.D
 Acq On : 09 Nov 2018 16:52
 Operator : JC/MD
 Sample : J5877-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SSSI-1106-S-TCLP-1

Quant Time: Nov 12 00:43:46 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 07 14:40:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	117912	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	164760	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	140832	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	66090	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	71817	52.98	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.96%	
35) Dibromofluoromethane	5.50	113	58225	51.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.92%	
50) Toluene-d8	8.72	98	199176	53.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.92%	
62) 4-Bromofluorobenzene	11.14	95	62777	46.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.74%	
Target Compounds						
16) Acetone	2.44	43	7513	10.294	ug/l	95
18) Methyl Acetate	2.78	43	13071	5.359	ug/l	100
20) Methylene Chloride	2.86	84	9140	7.624	ug/l	86
43) Isopropyl Acetate	6.46	43	23019	7.606	ug/l	100
95) Naphthalene	13.83	128	17762	4.825	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX110918\
Data File : VX005865.D
Acq On : 09 Nov 2018 16:52
Operator : JC/MD
Sample : J5877-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SSSI-1106-S-TCLP-1

Quant Time: Nov 12 00:43:46 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260
QLast Update : Wed Nov 07 14:40:53 2018
Response via : Initial Calibration

