

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110518\
 Data File : VX005785.D
 Acq On : 05 Nov 2018 18:17
 Operator : JC/MD
 Sample : J5771-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EO-01-110218-C

Quant Time: Nov 06 02:25:49 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103118W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 31 18:23:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	144411	50.00	ug/l	0.01
34) 1,4-Difluorobenzene	6.85	114	251570	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	230996	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	97749	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	114235	54.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.22%	
35) Dibromofluoromethane	5.51	113	88561	50.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.68%	
50) Toluene-d8	8.71	98	302990	49.59	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.18%	
62) 4-Bromofluorobenzene	11.14	95	98364	43.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.24%	
Target Compounds						
16) Acetone	2.44	43	16213	5.833	ug/l	97
18) Methyl Acetate	2.78	43	5307	1.751	ug/l	96
20) Methylene Chloride	2.86	84	5169	2.839	ug/l	93
43) Isopropyl Acetate	6.46	43	109806	22.125	ug/l	97

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX110518\
Data File : VX005785.D
Acq On : 05 Nov 2018 18:17
Operator : JC/MD
Sample : J5771-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
EO-01-110218-C

Quant Time: Nov 06 02:25:49 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X103118W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 31 18:23:06 2018
Response via : Initial Calibration

