

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX121318\
 Data File : VX006504.D
 Acq On : 13 Dec 2018 19:17
 Operator : JC/MD
 Sample : J6376-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TEST-PIT-1

Quant Time: Dec 14 01:11:50 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112918W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 30 03:55:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	632234	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	1009411	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	948720	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	461199	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	342271	53.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.06%	
35) Dibromofluoromethane	5.50	113	303515	46.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.88%	
50) Toluene-d8	8.71	98	1207775	50.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.62%	
62) 4-Bromofluorobenzene	11.13	95	453058	50.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.14%	
Target Compounds						
16) Acetone	2.45	43	15885	5.245	ug/l	93
18) Methyl Acetate	2.78	43	24591	2.428	ug/l	96
20) Methylene Chloride	2.86	84	27640	4.288	ug/l #	86
43) Isopropyl Acetate	6.46	43	32632	2.056	ug/l	97

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX121318\
Data File : VX006504.D
Acq On : 13 Dec 2018 19:17
Operator : JC/MD
Sample : J6376-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TEST-PIT-1

Quant Time: Dec 14 01:11:50 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112918W.M
Quant Title : SW846 8260
QLast Update : Fri Nov 30 03:55:33 2018
Response via : Initial Calibration

