

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071818\
 Data File : VX003454.D
 Acq On : 18 Jul 2018 18:14
 Operator : JC/MD
 Sample : PB111206TB
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB111206TB

Quant Time: Jul 19 03:11:32 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071618W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 16 13:15:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	267879	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	405178	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	385590	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	211560	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	175259	49.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.92%	
35) Dibromofluoromethane	5.51	113	153154	46.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.44%	
50) Toluene-d8	8.72	98	581355	48.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.10%	
62) 4-Bromofluorobenzene	11.14	95	199086	46.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.18%	
Target Compounds						
16) Acetone	2.45	43	19485	13.35	ug/l #	89
18) Methyl Acetate	2.78	43	7315	1.54	ug/l	92
20) Methylene Chloride	2.85	84	7379	Below	Cal	91
43) Isopropyl Acetate	6.47	43	145190	20.81	ug/l	95
74) N-amyl acetate	6.47	43	145190	22.29	ug/l #	95

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071818\
Data File : VX003454.D
Acq On : 18 Jul 2018 18:14
Operator : JC/MD
Sample : PB111206TB
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PB111206TB

Quant Time: Jul 19 03:11:32 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071618W.M
Quant Title : SW846 8260
QLast Update : Mon Jul 16 13:15:28 2018
Response via : Initial Calibration

