

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100318\
 Data File : VX004932.D
 Acq On : 02 Oct 2018 21:15
 Operator : JC/MD
 Sample : PB113401TB
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB113401TB

Quant Time: Oct 03 06:53:50 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 03 04:00:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	192670	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	332122	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	328548	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	173331	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	155127	56.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.72%	
35) Dibromofluoromethane	5.51	113	131980	46.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.78%	
50) Toluene-d8	8.72	98	482267	51.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.08%	
62) 4-Bromofluorobenzene	11.14	95	168130	49.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.74%	
Target Compounds						
16) Acetone	2.45	43	18683	11.627	ug/l #	88
18) Methyl Acetate	2.78	43	12151	2.985	ug/l	100
25) 2-Butanone	4.71	43	3951	1.668	ug/l	93
43) Isopropyl Acetate	6.46	43	198512	26.980	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100318\
Data File : VX004932.D
Acq On : 02 Oct 2018 21:15
Operator : JC/MD
Sample : PB113401TB
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PB113401TB

Quant Time: Oct 03 06:53:50 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 03 04:00:03 2018
Response via : Initial Calibration

