

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071118\
 Data File : VX003335.D
 Acq On : 11 Jul 2018 22:09
 Operator : JC/MD
 Sample : PB110983ZHE#09
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB110983ZHE#09

Quant Time: Jul 12 04:36:46 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X070918W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 09 15:24:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	250692	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	378961	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	354651	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	190782	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	171930	43.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.16%	
35) Dibromofluoromethane	5.51	113	146140	35.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	71.80%	
50) Toluene-d8	8.72	98	545436	37.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	75.44%	
62) 4-Bromofluorobenzene	11.14	95	184392	34.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	69.02%	
Target Compounds						
10) Methyl Iodide	2.52	142	1506	6.042	ug/l	95
16) Acetone	2.45	43	25911	18.499	ug/l	92
18) Methyl Acetate	2.78	43	29049	7.352	ug/l	94
20) Methylene Chloride	2.86	84	15309	4.694	ug/l	97

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071118\
Data File : VX003335.D
Acq On : 11 Jul 2018 22:09
Operator : JC/MD
Sample : PB110983ZHE#09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PB110983ZHE#09

Quant Time: Jul 12 04:36:46 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X070918W.M
Quant Title : SW846 8260
QLast Update : Mon Jul 09 15:24:16 2018
Response via : Initial Calibration

