

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100418\
 Data File : VX004987.D
 Acq On : 03 Oct 2018 23:21
 Operator : JC/MD
 Sample : PB113446ZHE#13
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB113446ZHE#13

Quant Time: Oct 04 03:32:57 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	5.66	168	290713	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	471144	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	441594	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	248106	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	201555	48.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.06%	
35) Dibromofluoromethane	5.51	113	177198	44.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.76%	
50) Toluene-d8	8.72	98	647302	48.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.52%	
62) 4-Bromofluorobenzene	11.14	95	241018	50.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.80%	
Target Compounds						
16) Acetone	2.45	43	21875	9.022	ug/l	95
18) Methyl Acetate	2.78	43	14348	2.336	ug/l	97
25) 2-Butanone	4.71	43	5457	1.527	ug/l	97
43) Isopropyl Acetate	6.46	43	247886	23.749	ug/l	95

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100418\
Data File : VX004987.D
Acq On : 03 Oct 2018 23:21
Operator : JC/MD
Sample : PB113446ZHE#13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PB113446ZHE#13

Quant Time: Oct 04 03:32:57 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

