

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100218\
 Data File : VU027341.D
 Acq On : 02 Oct 2018 11:51
 Operator : MD/SY
 Sample : PB113364ZHE#02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB113364ZHE#02

Quant Time: Oct 03 02:23:18 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 02 02:04:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	93846	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	163813	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	159686	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	73306	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	98076	62.59	ug/l	0.00
Spiked Amount 50.000			Recovery	=	125.18%	
35) Dibromofluoromethane	4.88	113	67741	60.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	120.68%	
50) Toluene-d8	7.57	98	260760	62.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	124.12%	
62) 4-Bromofluorobenzene	10.31	95	86591	55.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.90%	
Target Compounds						
16) Acetone	2.32	43	10275	11.099	ug/l	97
18) Methyl Acetate	2.62	43	3125	1.480	ug/l	91
20) Methylene Chloride	2.70	84	2611	2.025	ug/l	96
43) Isopropyl Acetate	5.55	43	107426	30.090	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100218\
Data File : VU027341.D
Acq On : 02 Oct 2018 11:51
Operator : MD/SY
Sample : PB113364ZHE#02
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
PB113364ZHE#02

Quant Time: Oct 03 02:23:18 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260
QLast Update : Tue Oct 02 02:04:34 2018
Response via : Initial Calibration

