

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU112818\
 Data File : VU028376.D
 Acq On : 28 Nov 2018 19:21
 Operator : MD/SY
 Sample : PB115109ZHE#05
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB115109ZHE#05

Quant Time: Nov 29 07:04:19 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U112818W.M
 Quant Title : SW846 8260
 QLast Update : Thu Nov 29 01:04:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	243790	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	465156	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	418120	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	167482	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	210879	51.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.52%	
35) Dibromofluoromethane	4.88	113	155056	50.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.26%	
50) Toluene-d8	7.57	98	608313	49.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.36%	
62) 4-Bromofluorobenzene	10.30	95	203793	45.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.42%	
Target Compounds						
16) Acetone	2.32	43	22796	9.146	ug/l	90
18) Methyl Acetate	2.62	43	7707	1.111	ug/l	96
20) Methylene Chloride	2.70	84	5184	1.403	ug/l	98
43) Isopropyl Acetate	5.55	43	31226	2.841	ug/l	97

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU112818\
Data File : VU028376.D
Acq On : 28 Nov 2018 19:21
Operator : MD/SY
Sample : PB115109ZHE#05
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
PB115109ZHE#05

Quant Time: Nov 29 07:04:19 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U112818W.M
Quant Title : SW846 8260
QLast Update : Thu Nov 29 01:04:49 2018
Response via : Initial Calibration

