

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
 Data File : VU027250.D
 Acq On : 28 Sep 2018 01:39
 Operator : MD/SY
 Sample : J5033-15
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BPSI-TT-MW311I-20180919

Quant Time: Sep 28 05:13:54 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 22 01:31:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	90210	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	156613	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	148489	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	68911	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	95650	52.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.36%	
35) Dibromofluoromethane	4.88	113	67760	49.91	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.82%	
50) Toluene-d8	7.57	98	248545	48.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.08%	
62) 4-Bromofluorobenzene	10.31	95	80688	45.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.90%	
Target Compounds						
19) Methyl tert-butyl Ether	3.00	73	13376	2.921	ug/l	99
27) cis-1,2-Dichloroethene	4.23	96	11117	7.400	ug/l	92
30) Chloroform	4.68	83	3310	1.198	ug/l	95
44) Trichloroethene	6.19	130	2504	1.851	ug/l	89

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027250.D
Acq On : 28 Sep 2018 01:39
Operator : MD/SY
Sample : J5033-15
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BPSI-TT-MW311I-20180919

Quant Time: Sep 28 05:13:54 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260
QLast Update : Sat Sep 22 01:31:57 2018
Response via : Initial Calibration

