

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092618\
 Data File : VU027180.D
 Acq On : 26 Sep 2018 16:13
 Operator : MD/SY
 Sample : J5049-07
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-40-SEP18

Quant Time: Sep 27 02:23:51 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	4.98	168	92414	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	154409	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	147982	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	69745	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	96979	52.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.28%	
35) Dibromofluoromethane	4.88	113	66995	50.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.12%	
50) Toluene-d8	7.57	98	244331	47.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.80%	
62) 4-Bromofluorobenzene	10.31	95	81645	46.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.30%	
Target Compounds						
16) Acetone	2.32	43	112691	104.397	ug/l	99
25) 2-Butanone	4.27	43	27938	17.873	ug/l	96
27) cis-1,2-Dichloroethene	4.23	96	5627	3.656	ug/l	95
51) 4-Methyl-2-Pentanone	7.47	43	3159	1.154	ug/l	89
52) Toluene	7.64	92	2253	0.627	ug/l	96

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092618\
Data File : VU027180.D
Acq On : 26 Sep 2018 16:13
Operator : MD/SY
Sample : J5049-07
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-40-SEP18

Quant Time: Sep 27 02:23:51 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

