

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU070218\
 Data File : VU025177.D
 Acq On : 03 Jul 2018 06:16
 Operator : MD/SY
 Sample : J3809-14
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DUP-1-062818

Quant Time: Jul 03 08:08:16 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 13 13:55:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	199676	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	295566	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	276725	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	143953	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	152477	46.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.56%	
35) Dibromofluoromethane	4.89	113	115904	47.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.50%	
50) Toluene-d8	7.57	98	426642	47.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.36%	
62) 4-Bromofluorobenzene	10.31	95	157770	44.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.82%	
Target Compounds						
11) Tert butyl alcohol	2.82	59	384756	485.31	ug/l	Qvalue 100
16) Acetone	2.32	43	8380	4.78	ug/l	100
19) Methyl tert-butyl Ether	3.00	73	79471	9.30	ug/l	98
22) Diisopropyl ether	3.57	45	61521	7.35	ug/l	97
40) Benzene	5.40	78	5136	0.52	ug/l	94

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU070218\
Data File : VU025177.D
Acq On : 03 Jul 2018 06:16
Operator : MD/SY
Sample : J3809-14
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DUP-1-062818

Quant Time: Jul 03 08:08:16 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260
QLast Update : Wed Jun 13 13:55:26 2018
Response via : Initial Calibration

