

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU070218\
 Data File : VU025157.D
 Acq On : 02 Jul 2018 20:54
 Operator : MD/SY
 Sample : J3243-06
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 NB-08-062918

Quant Time: Jul 03 05:49:36 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	203817	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	308727	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	283632	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	153819	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	157332	47.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.56%	
35) Dibromofluoromethane	4.89	113	122796	47.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.86%	
50) Toluene-d8	7.57	98	442956	47.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.80%	
62) 4-Bromofluorobenzene	10.31	95	164643	44.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.74%	
Target Compounds						
16) Acetone	2.32	43	16286	9.11	ug/l	97
18) Methyl Acetate	2.62	43	24151	6.30	ug/l	100
20) Methylene Chloride	2.70	84	60610	21.61	ug/l	96
95) Naphthalene	13.75	128	58299	5.76	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU070218\
Data File : VU025157.D
Acq On : 02 Jul 2018 20:54
Operator : MD/SY
Sample : J3243-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
NB-08-062918

Quant Time: Jul 03 05:49:36 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

