

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU070218\
 Data File : VU025148.D
 Acq On : 02 Jul 2018 17:14
 Operator : MD/SY
 Sample : J3771-02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 REACTOR-PAD

Quant Time: Jul 03 05:20:04 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	205182	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	312820	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	284946	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	151670	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	158593	47.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.70%	
35) Dibromofluoromethane	4.89	113	124133	47.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.62%	
50) Toluene-d8	7.57	98	448341	47.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.68%	
62) 4-Bromofluorobenzene	10.31	95	165457	44.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.02%	
Target Compounds						
16) Acetone	2.32	43	15762	8.76	ug/l	89
18) Methyl Acetate	2.62	43	8835	2.29	ug/l	100
20) Methylene Chloride	2.71	84	1255253	444.50	ug/l	97
84) 1,2,4-Trimethylbenzene	11.15	105	21621	2.10	ug/l	98
95) Naphthalene	13.75	128	103694	9.85	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU070218\
Data File : VU025148.D
Acq On : 02 Jul 2018 17:14
Operator : MD/SY
Sample : J3771-02
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
REACTOR-PAD

Quant Time: Jul 03 05:20:04 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

