

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071218\
 Data File : VU025379.D
 Acq On : 13 Jul 2018 08:04
 Operator : MD/SY
 Sample : J3825-02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 CL-02-071118-A

Quant Time: Jul 13 09:11:20 2018
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 10 09:34:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	193884	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	301066	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	277826	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	146700	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	152769	50.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.44%	
35) Dibromofluoromethane	4.89	113	116846	46.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.02%	
50) Toluene-d8	7.57	98	431186	49.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.16%	
62) 4-Bromofluorobenzene	10.31	95	156420	44.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.34%	
Target Compounds						
16) Acetone	2.32	43	27219	22.231	ug/l	94
18) Methyl Acetate	2.62	43	19118	6.199	ug/l	97
20) Methylene Chloride	2.71	84	27285	12.238	ug/l	98
68) m/p-Xylenes	9.38	106	9022	1.972	ug/l	90
84) 1,2,4-Trimethylbenzene	11.15	105	11378	1.177	ug/l	95
95) Naphthalene	13.75	128	74263	8.975	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU071218\
Data File : VU025379.D
Acq On : 13 Jul 2018 08:04
Operator : MD/SY
Sample : J3825-02
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
CL-02-071118-A

Quant Time: Jul 13 09:11:20 2018
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260
QLast Update : Tue Jul 10 09:34:16 2018
Response via : Initial Calibration

