

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU101018\
 Data File : VU027573.D
 Acq On : 10 Oct 2018 16:23
 Operator : MD/SY
 Sample : J5326-08
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 FL-PIT-2

Quant Time: Oct 11 05:24:35 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 02 02:04:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	148393	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	250743	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	220236	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	94526	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	107455	43.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.74%	
35) Dibromofluoromethane	4.88	113	80194	46.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.34%	
50) Toluene-d8	7.57	98	312734	48.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.26%	
62) 4-Bromofluorobenzene	10.31	95	101431	42.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	84.88%	
Target Compounds						
16) Acetone	2.32	43	8643	5.904	ug/l	92
18) Methyl Acetate	2.62	43	3448	1.033	ug/l	91
20) Methylene Chloride	2.70	84	916523	449.481	ug/l	94
43) Isopropyl Acetate	5.55	43	122598	22.435	ug/l	99
95) Naphthalene	13.75	128	28150	5.468	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU101018\
Data File : VU027573.D
Acq On : 10 Oct 2018 16:23
Operator : MD/SY
Sample : J5326-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-PIT-2

Quant Time: Oct 11 05:24:35 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260
QLast Update : Tue Oct 02 02:04:34 2018
Response via : Initial Calibration

