

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU101018\
 Data File : VU027551.D
 Acq On : 10 Oct 2018 02:58
 Operator : MD/SY
 Sample : J5329-08
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 POND-2

Quant Time: Oct 10 07:40:58 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	143202	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	236868	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	206123	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	84744	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	107712	45.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.08%	
35) Dibromofluoromethane	4.89	113	78973	48.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.30%	
50) Toluene-d8	7.57	98	294218	48.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.86%	
62) 4-Bromofluorobenzene	10.31	95	94210	41.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	83.44%	
Target Compounds						
16) Acetone	2.32	43	12658	8.961	ug/l	95
18) Methyl Acetate	2.62	43	3140	0.974	ug/l #	91
20) Methylene Chloride	2.70	84	2645	1.344	ug/l	95
25) 2-Butanone	4.29	43	6053	3.142	ug/l	95
43) Isopropyl Acetate	5.55	43	128088	24.812	ug/l	100

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU101018\
Data File : VU027551.D
Acq On : 10 Oct 2018 02:58
Operator : MD/SY
Sample : J5329-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
POND-2

Quant Time: Oct 10 07:40:58 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

