

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100118\
 Data File : VU027329.D
 Acq On : 01 Oct 2018 19:32
 Operator : MD/SY
 Sample : J5149-01
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WS-2

Quant Time: Oct 02 08:05:43 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	102524	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	173252	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	167890	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	80413	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	103308	60.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	120.68%	
35) Dibromofluoromethane	4.89	113	71579	60.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	120.56%	
50) Toluene-d8	7.57	98	279912	62.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	125.98%	
62) 4-Bromofluorobenzene	10.31	95	93977	56.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.80%	
Target Compounds						
3) Chloromethane	1.33	50	8677	4.336	ug/l	95
16) Acetone	2.32	43	20481	20.251	ug/l	97
25) 2-Butanone	4.28	43	4043	2.932	ug/l	99
51) 4-Methyl-2-Pentanone	7.47	43	12020	4.899	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100118\
Data File : VU027329.D
Acq On : 01 Oct 2018 19:32
Operator : MD/SY
Sample : J5149-01
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WS-2

Quant Time: Oct 02 08:05:43 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

