

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110218\
 Data File : VU027924.D
 Acq On : 02 Nov 2018 14:27
 Operator : MD/SY
 Sample : J5741-30
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 TP-7

Quant Time: Nov 03 05:05:52 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 02 02:28:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	187047	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	332705	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	293610	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	120515	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	146744	52.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.54%	
35) Dibromofluoromethane	4.88	113	107998	51.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.96%	
50) Toluene-d8	7.57	98	428657	52.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.80%	
62) 4-Bromofluorobenzene	10.31	95	138992	46.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.00%	
Target Compounds						
16) Acetone	2.32	43	7921	4.591	ug/l	99
18) Methyl Acetate	2.62	43	5040	1.062	ug/l #	88
20) Methylene Chloride	2.70	84	6640348	2635.916	ug/l	98
43) Isopropyl Acetate	5.55	43	196325	24.544	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU110218\
Data File : VU027924.D
Acq On : 02 Nov 2018 14:27
Operator : MD/SY
Sample : J5741-30
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
TP-7

Quant Time: Nov 03 05:05:52 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260
QLast Update : Fri Nov 02 02:28:20 2018
Response via : Initial Calibration

