

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU011619\
 Data File : VU029211.D
 Acq On : 16 Jan 2019 15:54
 Operator : JC/SP
 Sample : K1010-04
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 OK-01-011519-A

Quant Time: Jan 17 02:26:12 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U010319W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jan 12 01:33:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	110291	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	196134	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	184730	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	74636	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	76805	56.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.66%	
35) Dibromofluoromethane	4.88	113	66928	52.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.28%	
50) Toluene-d8	7.56	98	253682	50.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.08%	
62) 4-Bromofluorobenzene	10.30	95	82182	45.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.12%	
Target Compounds						
16) Acetone	2.32	43	7273	8.468	ug/l	100
20) Methylene Chloride	2.70	84	44773	30.703	ug/l	98
43) Isopropyl Acetate	5.55	43	7225	2.189	ug/l #	97
52) Toluene	7.64	92	85198	22.886	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU011619\
Data File : VU029211.D
Acq On : 16 Jan 2019 15:54
Operator : JC/SP
Sample : K1010-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
OK-01-011519-A

Quant Time: Jan 17 02:26:12 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U010319W.M
Quant Title : SW846 8260
QLast Update : Sat Jan 12 01:33:59 2019
Response via : Initial Calibration

