

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100418\
 Data File : VU027387.D
 Acq On : 04 Oct 2018 03:41
 Operator : MD/SY
 Sample : J5212-08
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 CG-SOIL-2

Quant Time: Oct 04 08:35:44 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	92591	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	167646	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	164622	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	84501	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	98688	63.83	ug/l	0.00
Spiked Amount 50.000			Recovery	=	127.66%	
35) Dibromofluoromethane	4.89	113	69981	60.91	ug/l	0.00
Spiked Amount 50.000			Recovery	=	121.82%	
50) Toluene-d8	7.57	98	260929	60.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	121.36%	
62) 4-Bromofluorobenzene	10.31	95	90754	56.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.58%	
Target Compounds						
16) Acetone	2.32	43	7152	7.830	ug/l	100
18) Methyl Acetate	2.62	43	2537	1.218	ug/l #	87
20) Methylene Chloride	2.70	84	27754	21.814	ug/l	98
43) Isopropyl Acetate	5.55	43	100527	27.514	ug/l	97
95) Naphthalene	13.75	128	49113	9.092	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027387.D
Acq On : 04 Oct 2018 03:41
Operator : MD/SY
Sample : J5212-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
CG-SOIL-2

Quant Time: Oct 04 08:35:44 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

