

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110818\
 Data File : VU028030.D
 Acq On : 08 Nov 2018 16:47
 Operator : MD/SY
 Sample : J5838-10
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 HC-SOIL-1

Quant Time: Nov 09 09:25:51 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	190316	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	367316	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	338414	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	142088	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	164346	58.08	ug/l	0.00
Spiked Amount 50.000			Recovery	=	116.16%	
35) Dibromofluoromethane	4.89	113	119538	51.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.22%	
50) Toluene-d8	7.57	98	482535	53.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.88%	
62) 4-Bromofluorobenzene	10.31	95	158633	48.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.14%	
Target Compounds						
16) Acetone	2.32	43	16218	9.238	ug/l	99
20) Methylene Chloride	2.71	84	11856	4.625	ug/l	99
30) Chloroform	4.68	83	5691	1.238	ug/l	95
43) Isopropyl Acetate	5.55	43	49622	5.619	ug/l	99
95) Naphthalene	13.75	128	32818	2.644	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU110818\
Data File : VU028030.D
Acq On : 08 Nov 2018 16:47
Operator : MD/SY
Sample : J5838-10
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
HC-SOIL-1

Quant Time: Nov 09 09:25:51 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

