

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW010919\
 Data File : VW008102.D
 Acq On : 09 Jan 2019 14:50
 Operator : SY/AP
 Sample : K1071-07
 Misc : 5.10G/5ML/MSVOA W/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 CPSB-1(10-15)

Quant Time: Jan 10 06:28:20 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
 Quant Title : SW846 8260
 QLast Update : Wed Jan 09 04:40:48 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.95	168	76825	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.84	114	128197	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.63	117	107750	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.56	152	36878	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.31	65	41713	48.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.62%	
35) Dibromofluoromethane	7.88	113	39260	52.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.02%	
50) Toluene-d8	10.32	98	158424	50.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.98%	
62) 4-Bromofluorobenzene	12.62	95	46525	41.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	82.72%	
Target Compounds						
16) Acetone	4.12	43	19171	111.832	ug/l	93
17) Carbon Disulfide	4.38	76	6092	2.318	ug/l #	89
25) 2-Butanone	7.17	43	5735	23.872	ug/l #	86
39) Methylcyclohexane	9.34	83	5031	3.005	ug/l	98
65) Chlorobenzene	11.66	112	10092	4.421	ug/l	98

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

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