

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW091119\
 Data File : VW012918.D
 Acq On : 11 Sep 2019 15:35
 Operator : SY/VA
 Sample : K4683-02
 Misc : 4.17G/5ML/MSVOA W/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 20190664-LOCATION-2

Quant Time: Sep 12 08:13:42 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W090719S.M
 Quant Title : SW846 8260
 QLast Update : Sun Sep 08 05:04:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.95	168	263349	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.84	114	392992	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.63	117	312243	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.55	152	111009	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.30	65	136198	46.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.58%	
35) Dibromofluoromethane	7.88	113	115210	49.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.62%	
50) Toluene-d8	10.32	98	461198	50.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.76%	
62) 4-Bromofluorobenzene	12.61	95	133106	40.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	80.12%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	5.21	59	7599	33.754	ug/l #	90
16) Acetone	4.13	43	236563	294.901	ug/l	98
17) Carbon Disulfide	4.38	76	36235	7.849	ug/l #	93
18) Methyl Acetate	4.68	43	6988	3.713	ug/l #	91
25) 2-Butanone	7.17	43	63643	69.321	ug/l	99
29) Tetrahydrofuran	7.54	42	4287	6.618	ug/l	92

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

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