

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW102419\
 Data File : VW013718.D
 Acq On : 24 Oct 2019 18:28
 Operator : SY/VA
 Sample : K5501-12RE
 Misc : 5.62G/5ML/MSVOA W/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 184-98-(2.2)RE

Quant Time: Oct 25 07:59:44 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W102319S.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 23 14:11:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.98	168	247	50.00	ug/l	0.03
34) 1,4-Difluorobenzene	8.88	114	326	50.00	ug/l	0.04
63) Chlorobenzene-d5	11.69	117	65	50.00	ug/l	0.07
72) 1,4-Dichlorobenzene-d4	13.55	152	20	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	0.00	65	0	0.00	ug/l	
Spiked Amount	50.000					
			Recovery	=	0.00%	
35) Dibromofluoromethane	7.91	113	82	45.27	ug/l	0.02
Spiked Amount	50.000					
			Recovery	=	90.54%	
50) Toluene-d8	10.34	98	252	34.14	ug/l	0.01
Spiked Amount	50.000					
			Recovery	=	68.28%	
62) 4-Bromofluorobenzene	0.00	95	0	0.00	ug/l	
Spiked Amount	50.000					
			Recovery	=	0.00%	
Target Compounds						
					Qvalue	
18) Methyl Acetate	4.95	43	23	21.356	ug/l #	50
25) 2-Butanone	7.04	43	21	37.179	ug/l #	45
28) Bromochloromethane	7.53	49	25	18.694	ug/l #	7
37) Ethyl Acetate	7.04	43	21	18.069	ug/l #	67
40) Benzene	8.39	78	20	2.328	ug/l #	51
51) 4-Methyl-2-Pentanone	10.29	43	20	18.092	ug/l #	35
74) N-amyl acetate	12.09	43	51	183.563	ug/l #	41
92) 1,2-Dibromo-3-Chloropropan	14.54	75	51	1488.055	ug/l #	1

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

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