

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW010820\
 Data File : VW014559.D
 Acq On : 08 Jan 2020 14:37
 Operator : SY/VA
 Sample : L1056-03
 Misc : 8.36G/5ML/MSVOA_W/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 WP022SB447_200106_35-36

Quant Time: Jan 08 15:40:55 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010320S.M
 Quant Title : SW846 8260
 QLast Update : Fri Jan 03 16:07:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.95	168	604312	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.84	114	942988	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.63	117	673278	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.55	152	216454	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.30	65	251828	47.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.32%	
35) Dibromofluoromethane	7.88	113	258295	47.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.86%	
50) Toluene-d8	10.32	98	978519	43.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.02%	
62) 4-Bromofluorobenzene	12.62	95	246230	32.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	64.56%	
Target Compounds						
16) Acetone	4.12	43	19059	12.775	ug/l	99
20) Methylene Chloride	4.92	84	40438	2.400	ug/l	96
25) 2-Butanone	7.19	43	5494	2.870	ug/l #	83
29) Tetrahydrofuran	7.53	42	3513	2.962	ug/l	94
65) Chlorobenzene	11.65	112	51027	3.699	ug/l	95
74) N-amyl acetate	12.25	43	103886	29.172	ug/l #	64

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

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