

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081221\
 Data File : VY005616.D
 Acq On : 12 Aug 2021 12:46
 Operator : SY/MD
 Sample : M3318-04
 Misc : 5.03G/5ML/MSVOA_Y/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 20210595-AMENDED-COM-9

Quant Time: Aug 13 04:21:10 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y080221S.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 03 04:52:05 2021
 Response via : Initial Calibration

Compound	R.T. QIon		Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	7.801	168	258347	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.697	114	442038	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.496	117	458189	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.428	152	238183	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.149	65	164868	63.044	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	= 126.080%	
35) Dibromofluoromethane	7.722	113	106861	43.179	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	= 86.360%	
50) Toluene-d8	10.185	98	571113	55.024	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	= 110.040%	
62) 4-Bromofluorobenzene	12.483	95	216309	62.933	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	= 125.860%	
Target Compounds						
					Qvalue	
16) Acetone	3.954	43	101693	154.236	ug/l	96
17) Carbon Disulfide	4.204	76	166872	24.314	ug/l	100
20) Methylene Chloride	4.729	84	34712	8.214	ug/l	93
25) 2-Butanone	6.997	43	26589	27.281	ug/l	93
84) 1,2,4-Trimethylbenzene	13.123	105	15230	1.119	ug/l	80

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

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