

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100623\
 Data File : VY015864.D
 Acq On : 06 Oct 2023 15:59
 Operator : SY/MD
 Sample : 04767-01
 Misc : 4.95g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 D-1(0-5)

Quant Time: Oct 07 07:28:24 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100523S.M
 Quant Title : SW846 8260
 QLast Update : Fri Oct 06 05:54:04 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.783	168	224861	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.685	114	392649	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.489	117	362356	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.422	152	165015	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.136	65	147533	66.689	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	= 133.380%	
35) Dibromofluoromethane	7.716	113	126497	56.519	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	= 113.040%	
50) Toluene-d8	10.179	98	461789	63.550	ug/l	0.00
Spiked Amount	50.000	Range	58 - 134	Recovery	= 127.100%	
62) 4-Bromofluorobenzene	12.483	95	158906	53.500	ug/l	0.00
Spiked Amount	50.000	Range	30 - 143	Recovery	= 107.000%	
Target Compounds						
					Qvalue	
16) Acetone	3.972	43	32825	28.639	ug/l #	83
17) Carbon Disulfide	4.180	76	14588	9.025	ug/l	99
27) cis-1,2-Dichloroethene	6.984	96	12079	4.956	ug/l	84
40) Benzene	8.161	78	13435	1.603	ug/l	99
44) Trichloroethene	8.935	130	22616	9.223	ug/l	95
64) Tetrachloroethene	10.715	164	43731	17.170	ug/l	96

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

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