

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110824\
 Data File : VY020230.D
 Acq On : 08 Nov 2024 14:02
 Operator : SY/MD
 Sample : P4722-01
 Misc : 10.39g/5.0mL/MSVOA_Y/SOIL/B
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 WC-1(5.5)

Quant Time: Nov 08 22:18:23 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103024S.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 31 15:23:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.707	168	209744	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.616	114	444878	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.414	117	397795	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.347	152	127098	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.061	65	159975	61.112	ug/l	0.00
Spiked Amount	50.000	Range 50 - 163	Recovery	=	122.220%	
35) Dibromofluoromethane	7.634	113	148294	52.428	ug/l	0.00
Spiked Amount	50.000	Range 54 - 147	Recovery	=	104.860%	
50) Toluene-d8	10.103	98	550182	48.867	ug/l	0.00
Spiked Amount	50.000	Range 58 - 134	Recovery	=	97.740%	
62) 4-Bromofluorobenzene	12.408	95	159777	43.691	ug/l	0.00
Spiked Amount	50.000	Range 29 - 146	Recovery	=	87.380%	
Target Compounds						
						Qvalue
16) Acetone	3.873	43	17678	29.039	ug/l	# 84
20) Methylene Chloride	4.610	84	10151	3.712	ug/l	89
25) 2-Butanone	6.903	43	8035	9.737	ug/l	# 79
30) Chloroform	7.427	83	6195	1.272	ug/l	88
64) Tetrachloroethene	10.646	164	6464	2.461	ug/l	94

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

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