

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY103124\
 Data File : VY020098.D
 Acq On : 31 Oct 2024 16:35
 Operator : SY/MD
 Sample : P4612-03
 Misc : 5.01g/5.0mL/MSVOA_Y/SOIL/A
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 MOO-24-00335

Quant Time: Nov 01 03:49:13 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	167867	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.616	114	346815	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.414	117	302900	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.346	152	94133	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.061	65	126809	60.527	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	= 121.060%	
35) Dibromofluoromethane	7.634	113	119234	54.074	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	= 108.140%	
50) Toluene-d8	10.103	98	422039	48.084	ug/l	0.00
Spiked Amount	50.000	Range	58 - 134	Recovery	= 96.160%	
62) 4-Bromofluorobenzene	12.402	95	121685	42.683	ug/l	0.00
Spiked Amount	50.000	Range	29 - 146	Recovery	= 85.360%	
Target Compounds						
						Qvalue
16) Acetone	3.873	43	65006	133.421	ug/l	# 82
51) 4-Methyl-2-Pentanone	9.999	43	5906	3.670	ug/l	90
52) Toluene	10.164	92	16704	2.726	ug/l	94
59) 2-Hexanone	10.762	43	3921	3.399	ug/l	88

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

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