

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY121422\
 Data File : VY011858.D
 Acq On : 14 Dec 2022 20:05
 Operator : KP/MD
 Sample : N6038-11
 Misc : 3.93g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 RB-34-(1.5-2)

Quant Time: Dec 15 01:24:14 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y120922S.M
 Quant Title : SW846 8260
 QLast Update : Sat Dec 10 04:19:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.789	168	261837	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.691	114	440482	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.489	117	362087	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.428	152	141938	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.142	65	88201	49.297	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	98.600%
35) Dibromofluoromethane	7.722	113	32552	13.819	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	27.640%#
50) Toluene-d8	10.179	98	259186	38.924	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	=	77.840%
62) 4-Bromofluorobenzene	12.483	95	95478	36.118	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	=	72.240%
Target Compounds						
					Qvalue	
16) Acetone	3.948	43	75977	106.268	ug/l	97
20) Methylene Chloride	4.716	84	33152	8.103	ug/l	97
25) 2-Butanone	6.990	43	9985	10.857	ug/l	98
84) 1,2,4-Trimethylbenzene	13.123	105	12701	1.660	ug/l	91

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY121422\
Data File : VY011858.D
Acq On : 14 Dec 2022 20:05
Operator : KP/MD
Sample : N6038-11
Misc : 3.93g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
RB-34-(1.5-2)

Quant Time: Dec 15 01:24:14 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y120922S.M
Quant Title : SW846 8260
QLast Update : Sat Dec 10 04:19:59 2022
Response via : Initial Calibration

