

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY091824\
 Data File : VY019618.D
 Acq On : 18 Sep 2024 20:02
 Operator : SY/MD
 Sample : P4088-15
 Misc : 5.93g/5.0mL/MSVOA_Y/SOIL/A
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 NB-614-VOC-09

Quant Time: Sep 19 01:28:24 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y090924S.M
 Quant Title : SW846 8260
 QLast Update : Mon Sep 09 16:00:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.707	168	304405	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.616	114	588356	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.414	117	519151	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.347	152	189932	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.061	65	187039	66.356	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	= 132.720%	
35) Dibromofluoromethane	7.634	113	186876	56.020	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	= 112.040%	
50) Toluene-d8	10.109	98	702687	56.950	ug/l	0.00
Spiked Amount	50.000	Range	58 - 134	Recovery	= 113.900%	
62) 4-Bromofluorobenzene	12.408	95	220754	46.530	ug/l	0.00
Spiked Amount	50.000	Range	29 - 146	Recovery	= 93.060%	
Target Compounds						
					Qvalue	
16) Acetone	3.873	43	56282	83.813	ug/l	88
25) 2-Butanone	6.896	43	22242	23.288	ug/l	91

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY091824\
Data File : VY019618.D
Acq On : 18 Sep 2024 20:02
Operator : SY/MD
Sample : P4088-15
Misc : 5.93g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NB-614-VOC-09

Quant Time: Sep 19 01:28:24 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y090924S.M
Quant Title : SW846 8260
QLast Update : Mon Sep 09 16:00:30 2024
Response via : Initial Calibration

