

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070924\
 Data File : VD079363.D
 Acq On : 09 Jul 2024 16:02
 Operator : RP/MD
 Sample : P3167-14RE
 Misc : 5.02G/5.0ml/MSVOA_D/SOIL/B
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 BP-VPB-196-GW-820-822RE

Quant Time: Jul 09 16:54:52 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D070824S.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 08 17:06:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.875	168	112369	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.775	114	215916	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.581	117	197455	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.510	152	64578	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.228	65	71950	54.335	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	108.660%
35) Dibromofluoromethane	7.799	113	82004	53.835	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	107.680%
50) Toluene-d8	10.269	98	279744	47.642	ug/l	0.00
Spiked Amount	50.000	Range	58 - 134	Recovery	=	95.280%
62) 4-Bromofluorobenzene	12.569	95	77122	42.258	ug/l	0.00
Spiked Amount	50.000	Range	30 - 143	Recovery	=	84.520%
Target Compounds						
					Qvalue	
16) Acetone	4.022	43	3604	14.310	ug/l	99
20) Methylene Chloride	4.793	84	27633	6.666	ug/l	94

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070924\
Data File : VD079363.D
Acq On : 09 Jul 2024 16:02
Operator : RP/MD
Sample : P3167-14RE
Misc : 5.02G/5.0ml/MSVOA_D/SOIL/B
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BP-VPB-196-GW-820-822RE

Quant Time: Jul 09 16:54:52 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D070824S.M
Quant Title : SW846 8260
QLast Update : Mon Jul 08 17:06:59 2024
Response via : Initial Calibration

