

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD091922\
 Data File : VD074351.D
 Acq On : 19 Sep 2022 17:31
 Operator : VA/SY
 Sample : N4655-01
 Misc : 5.08G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SU-3-091522

Quant Time: Sep 19 23:02:09 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D091622S.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 17 04:16:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.881	168	450	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.775	114	745	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.587	117	814	50.000	ug/l	# 0.01
72) 1,4-Dichlorobenzene-d4	0.000	152	0	0.000	ug/l	-13.51
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.234	65	73	16.655	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	33.300%#
35) Dibromofluoromethane	7.799	113	136	27.718	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	55.440%
50) Toluene-d8	10.269	98	748	44.181	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	=	88.360%
62) 4-Bromofluorobenzene	12.563	95	128	18.735	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	=	37.460%
Target Compounds						
6) Chloroethane	2.734	64	1641	721.918	ug/l #	44
16) Acetone	3.999	43	125	126.586	ug/l #	42
43) Isopropyl Acetate	8.446	43	61	11.019	ug/l #	51

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD091922\
Data File : VD074351.D
Acq On : 19 Sep 2022 17:31
Operator : VA/SY
Sample : N4655-01
Misc : 5.08G/5.00ml/MSVOA_D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-3-091522

Quant Time: Sep 19 23:02:09 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D091622S.M
Quant Title : SW846 8260
QLast Update : Sat Sep 17 04:16:54 2022
Response via : Initial Calibration

