

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD091120\
 Data File : VD066660.D
 Acq On : 11 Sep 2020 20:36
 Operator : VA/SY
 Sample : L3981-11
 Misc : 11.80G/5.00ml/MSVOA D/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 B20-4-6

Manual Integrations
 APPROVED

MMDadoda
 9/14/2020 12:07:03 PM

Quant Time: Sep 12 02:10:36 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 04 01:57:04 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.98	168	397914	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.86	114	705891	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.64	117	681361	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.57	152	281192	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.33	65	251839	52.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.62%	
35) Dibromofluoromethane	7.91	113	245377	51.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.06%	
50) Toluene-d8	10.34	98	887343	52.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.28%	
62) 4-Bromofluorobenzene	12.63	95	298953	52.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.94%	
Target Compounds						
27) cis-1,2-Dichloroethene	7.21	96	5172m	1.010	ug/l	
44) Trichloroethene	9.11	130	9364	1.774	ug/l	92
64) Tetrachloroethene	10.87	164	773902	168.693	ug/l	93

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD091120\
Data File : VD066660.D
Acq On : 11 Sep 2020 20:36
Operator : VA/SY
Sample : L3981-11
Misc : 11.80G/5.00ml/MSVOA D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B20-4-6

Manual Integrations
APPROVED

MMDadoda
9/14/2020 12:07:03 PM

Quant Time: Sep 12 02:10:36 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260
QLast Update : Fri Sep 04 01:57:04 2020
Response via : Initial Calibration

