

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD091120\
 Data File : VD066650.D
 Acq On : 11 Sep 2020 16:02
 Operator : VA/SY
 Sample : L3981-01
 Misc : 15.68G/5.00ml/MSVOA D/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 B17-0-2

Manual Integrations
 APPROVED

MMDadoda
 9/14/2020 12:06:59 PM

Quant Time: Sep 21 00:53:06 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	7.97	168	364239	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.86	114	620192	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.64	117	533727	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.58	152	229538	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.33	65	186092	42.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	84.46%	
35) Dibromofluoromethane	7.91	113	196159	46.88	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.76%	
50) Toluene-d8	10.34	98	740231	49.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.00%	
62) 4-Bromofluorobenzene	12.63	95	229270	46.24	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.48%	
Target Compounds						
20) Methylene Chloride	4.96	84	63913	9.041	ug/l	97
27) cis-1,2-Dichloroethene	7.21	96	532096	113.544	ug/l	89
44) Trichloroethene	9.11	130	990952	213.713	ug/l	98
64) Tetrachloroethene	10.87	164	43899433m	12215.993	ug/l	
91) 1,2-Dichlorobenzene	13.89	146	13698	2.083	ug/l #	85

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

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