

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD112119\
 Data File : VD064332.D
 Acq On : 21 Nov 2019 18:51
 Operator : VA/SY
 Sample : K5670-09
 Misc : 6.32G/5.00ml/MSVOA D/SOIL
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 OK-01-112119

Quant Time: Nov 22 02:39:09 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110819S.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 12 02:43:24 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.98	168	917896	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.87	114	1319220	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.65	117	1146826	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.58	152	535651	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.34	65	387392	51.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.64%	
35) Dibromofluoromethane	7.92	113	400889	52.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.88%	
50) Toluene-d8	10.34	98	1518700	53.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.32%	
62) 4-Bromofluorobenzene	12.64	95	486449	49.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.34%	
Target Compounds						
16) Acetone	4.17	43	70255	39.284	ug/l	99
18) Methyl Acetate	4.73	43	11707	2.344	ug/l	93
30) Chloroform	7.72	83	143090	9.522	ug/l	94
52) Toluene	10.41	92	26465	1.210	ug/l	96

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

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