

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD040519\
 Data File : VD061625.D
 Acq On : 5 Apr 2019 17:12
 Operator : VA/AP
 Sample : K2257-25RE
 Misc : 4.88g/5ml/MSVOA D/SOIL
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SP-1-DRE

Quant Time: Apr 06 00:27:38 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D032719S.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 28 09:22:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.03	168	14031	50.00	ug/l	0.01
34) 1,4-Difluorobenzene	8.20	114	32394	50.00	ug/l	0.01
63) Chlorobenzene-d5	12.36	117	28167	50.00	ug/l	0.01
72) 1,4-Dichlorobenzene-d4	14.51	152	10786	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.43	65	11999	100.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	200.22%	
35) Dibromofluoromethane	6.90	113	13980	56.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.94%	
50) Toluene-d8	10.39	98	25590	43.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.20%	
62) 4-Bromofluorobenzene	13.55	95	11107	48.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.56%	
Target Compounds						
					Qvalue	
5) Bromomethane	2.19	94	815	36.629	ug/l #	10
16) Acetone	3.21	43	4854	157.538	ug/l #	89
17) Carbon Disulfide	3.38	76	1317	2.691	ug/l #	75
20) Methylene Chloride	3.80	84	1562	7.414	ug/l #	92
38) Carbon Tetrachloride	7.02	117	487	1.499	ug/l #	10
52) Toluene	10.49	92	1038	2.398	ug/l	93
81) trans-1,4-Dichloro-2-buten	13.55	75	4430	93.252	ug/l #	18

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

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