

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW110618\
 Data File : VW006711.D
 Acq On : 07 Nov 2018 03:45
 Operator : SY/AP
 Sample : J5764-09RE
 Misc : 6.03G/5ML/MSVOA W/SOIL
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 JC-03-110218-ERE

Quant Time: Nov 07 12:30:03 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W110218S.M
 Quant Title : SW846 8260
 QLast Update : Thu Nov 01 03:48:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.95	168	248468	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.85	114	390798	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.63	117	331064	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.57	152	129815	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.30	65	110638	51.80	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.60%	
35) Dibromofluoromethane	7.89	113	259	0.11	ug/l	0.00
Spiked Amount	50.000		Recovery	=	0.22%	
50) Toluene-d8	10.33	98	388912	41.66	ug/l	0.00
Spiked Amount	50.000		Recovery	=	83.32%	
62) 4-Bromofluorobenzene	12.62	95	111181	32.85	ug/l	0.00
Spiked Amount	50.000		Recovery	=	65.70%	
Target Compounds						
					Qvalue	
16) Acetone	4.13	43	15540	45.066	ug/l	100
17) Carbon Disulfide	4.38	76	31489	5.460	ug/l	100
39) Methylcyclohexane	9.34	83	6061	1.390	ug/l	90
40) Benzene	8.33	78	22549	2.310	ug/l	98
52) Toluene	10.39	92	49090	7.498	ug/l	100
68) m/p-Xylenes	11.84	106	5804	1.248	ug/l	96

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

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