

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW081919\
 Data File : VW012002.D
 Acq On : 19 Aug 2019 20:54
 Operator : SY/VA
 Sample : K4391-07
 Misc : 5.07G/5ML/MSVOA W/SOIL
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 S-CORNER-1

Quant Time: Aug 20 05:58:50 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W072719S.M
 Quant Title : SW846 8260
 QLast Update : Sat Jul 27 00:23:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.95	168	265747	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.84	114	448640	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.63	117	371014	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.56	152	165536	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.31	65	180917	59.29	ug/l	0.00
Spiked Amount	50.000		Recovery	=	118.58%	
35) Dibromofluoromethane	7.88	113	136108	54.08	ug/l	0.00
Spiked Amount	50.000		Recovery	=	108.16%	
50) Toluene-d8	10.32	98	590061	57.26	ug/l	0.00
Spiked Amount	50.000		Recovery	=	114.52%	
62) 4-Bromofluorobenzene	12.62	95	218990	59.61	ug/l	0.00
Spiked Amount	50.000		Recovery	=	119.22%	
Target Compounds						
16) Acetone	4.14	43	14889	15.339	ug/l	92
31) Cyclohexane	7.95	56	22275	3.904	ug/l #	56
39) Methylcyclohexane	9.34	83	133312	25.953	ug/l	88
80) 1,3,5-Trimethylbenzene	12.94	105	54391	5.254	ug/l	100
83) tert-Butylbenzene	13.21	119	9643	1.114	ug/l	75
85) sec-Butylbenzene	13.38	105	30870	2.499	ug/l #	66
86) p-Isopropyltoluene	13.50	119	14964	1.335	ug/l	93

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

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