

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY120319\
 Data File : VY000842.D
 Acq On : 03 Dec 2019 16:41
 Operator : SY/MD
 Sample : K6120-05
 Misc : 9.46G/5ML/MSVOA Y/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 TH-4

Quant Time: Dec 04 01:24:08 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\82Y112719S.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 27 12:40:47 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.80	168	248045	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.69	114	469337	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.49	117	424456	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.42	152	183216	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.15	65	151858	57.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	115.24%	
35) Dibromofluoromethane	7.72	113	140620	51.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.82%	
50) Toluene-d8	10.18	98	566441	51.80	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.60%	
62) 4-Bromofluorobenzene	12.48	95	187060	45.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.10%	
Target Compounds						
16) Acetone	3.96	43	67928	112.348	ug/l	97
20) Methylene Chloride	4.73	84	21058	6.607	ug/l #	82
67) Ethyl Benzene	11.59	91	86789	5.282	ug/l	97
68) m/p-Xylenes	11.70	106	159510	25.863	ug/l	96
69) o-Xylene	12.03	106	51406	8.939	ug/l	98
84) 1,2,4-Trimethylbenzene	13.12	105	23710	1.962	ug/l	100

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

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