

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY020520\
 Data File : VY001511.D
 Acq On : 05 Feb 2020 16:49
 Operator : SY/MD
 Sample : L1393-04
 Misc : 5.25G/5ML/MSVOA Y/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 HAM-SB-03-1-2

Quant Time: Feb 06 07:43:47 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\82Y020420S.M
 Quant Title : SW846 8260
 QLast Update : Tue Feb 04 15:10:08 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.82	168	169238	50.00	ug/l	0.01
34) 1,4-Difluorobenzene	8.71	114	319340	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.50	117	281357	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.44	152	104561	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	8.17	65	109428	64.22	ug/l	0.00
Spiked Amount						
			Recovery	=	128.44%	
35) Dibromofluoromethane	7.75	113	94370	51.25	ug/l	0.00
Spiked Amount						
			Recovery	=	102.50%	
50) Toluene-d8	10.19	98	383070	54.98	ug/l	0.00
Spiked Amount						
			Recovery	=	109.96%	
62) 4-Bromofluorobenzene	12.49	95	129788	47.29	ug/l	0.00
Spiked Amount						
			Recovery	=	94.58%	

Target Compounds

						Qvalue
16) Acetone	3.98	43	3403760	6905.647	ug/l	96
17) Carbon Disulfide	4.23	76	20563	3.630	ug/l	100
25) 2-Butanone	7.03	43	25150	39.330	ug/l	93
51) 4-Methyl-2-Pentanone	10.09	43	1644	1.192	ug/l	91
52) Toluene	10.26	92	6167	1.166	ug/l	96
67) Ethyl Benzene	11.61	91	18903	1.851	ug/l	94
68) m/p-Xylenes	11.71	106	42950	11.201	ug/l	99
69) o-Xylene	12.04	106	18115	5.010	ug/l	95

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

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