

Data Path : Z:\voasrv\HPCHEM1\MSVOA Y\Data\vv060520\
 Data File : VY002847.D
 Acq On : 05 Jun 2020 14:38
 Operator : SY/MD
 Sample : L2930-01
 Misc : 6.06G/5ML/MSVOA Y/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 S39-HSVE2-IDW01

Quant Time: Jun 18 16:55:18 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\82Y052720S.M
 Quant Title : SW846 8260
 QLast Update : Wed May 27 13:13:33 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.80	168	225261	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.69	114	348211	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.49	117	310008	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.42	152	151220	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.14	65	96738	49.46	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.92%	
35) Dibromofluoromethane	7.72	113	93028	45.59	ug/l	0.00
Spiked Amount	50.000		Recovery	=	91.18%	
50) Toluene-d8	10.18	98	406995	49.33	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.66%	
62) 4-Bromofluorobenzene	12.48	95	127558	45.24	ug/l	0.00
Spiked Amount	50.000		Recovery	=	90.48%	
Target Compounds						
16) Acetone	3.95	43	850220	1853.817	ug/l	97
17) Carbon Disulfide	4.20	76	21683	3.744	ug/l	96
25) 2-Butanone	7.00	43	152238	213.055	ug/l	96
29) Tetrahydrofuran	7.36	42	3048	6.558	ug/l #	92
44) Trichloroethene	8.94	130	6909	2.524	ug/l	98
64) Tetrachloroethene	10.72	164	7551	2.680	ug/l	92

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

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