

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY012420\
 Data File : VY001390.D
 Acq On : 24 Jan 2020 18:08
 Operator : SY/MD
 Sample : L1263-02
 Misc : 6.63G/5ML/MSVOA Y/SOIL
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 CTY-SB-02-6-7

Quant Time: Jan 25 04:46:06 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\82Y012120S.M
 Quant Title : SW846 8260
 QLast Update : Thu Jan 23 01:21:47 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.82	168	162815	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.71	114	295457	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.50	117	267505	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.43	152	110226	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.16	65	96908	54.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.40%	
35) Dibromofluoromethane	7.74	113	87595	50.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.02%	
50) Toluene-d8	10.19	98	352265	52.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.58%	
62) 4-Bromofluorobenzene	12.49	95	122156	45.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.06%	
Target Compounds						
16) Acetone	3.98	43	23166	43.372	ug/l	100
17) Carbon Disulfide	4.22	76	21693	3.990	ug/l	95
25) 2-Butanone	7.02	43	7525	9.759	ug/l	99
95) Naphthalene	15.23	128	37537	7.148	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY012420\
Data File : VY001390.D
Acq On : 24 Jan 2020 18:08
Operator : SY/MD
Sample : L1263-02
Misc : 6.63G/5ML/MSVOA Y/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
CTY-SB-02-6-7

Quant Time: Jan 25 04:46:06 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y012120S.M
Quant Title : SW846 8260
QLast Update : Thu Jan 23 01:21:47 2020
Response via : Initial Calibration

