

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY082421\
 Data File : VY005819.D
 Acq On : 24 Aug 2021 14:00
 Operator : SY/MD
 Sample : M3463-15
 Misc : 4.03G/5ML/MSVOA_Y/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 SWCS-16-0204

Quant Time: Aug 25 04:29:45 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y080221S.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 03 04:52:05 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	7.801	168	186485	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.697	114	323789	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.496	117	328867	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.428	152	148534	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.149	65	118518	62.785	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	125.560%
35) Dibromofluoromethane	7.728	113	98583	54.381	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	108.760%
50) Toluene-d8	10.185	98	418497	55.045	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	=	110.080%
62) 4-Bromofluorobenzene	12.483	95	146738	58.283	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	=	116.560%
Target Compounds						Qvalue
16) Acetone	3.960	43	6694	14.065	ug/l	89
17) Carbon Disulfide	4.204	76	23729	4.790	ug/l	97
39) Methylcyclohexane	9.185	83	6182	1.600	ug/l #	45
64) Tetrachloroethene	10.721	164	3142	1.255	ug/l	97
65) Chlorobenzene	11.520	112	9022	1.363	ug/l	99

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

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