

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092921\
 Data File : VY006226.D
 Acq On : 29 Sep 2021 20:20
 Operator : SY/MD
 Sample : M3966-02
 Misc : 7.21G/5ML/MSVOA_Y/SOIL
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 S-831-J-SO-2-2.5-092721

Quant Time: Sep 30 07:35:15 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y092821S.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 28 17:44:36 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.795	168	138928	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.697	114	230215	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.496	117	211868	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.428	152	87138	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.149	65	95589	64.286	ug/l	0.00
Spiked Amount 50.000	Range 50 - 163		Recovery = 128.580%			
35) Dibromofluoromethane	7.722	113	68857	55.984	ug/l	0.00
Spiked Amount 50.000	Range 54 - 147		Recovery = 111.960%			
50) Toluene-d8	10.185	98	289288	56.199	ug/l	0.00
Spiked Amount 50.000	Range 49 - 140		Recovery = 112.400%			
62) 4-Bromofluorobenzene	12.483	95	87699	50.151	ug/l	0.00
Spiked Amount 50.000	Range 25 - 144		Recovery = 100.300%			
Target Compounds					Qvalue	
16) Acetone	3.954	43	35899	86.441	ug/l	97
17) Carbon Disulfide	4.204	76	40723	9.358	ug/l	95
25) 2-Butanone	6.996	43	8038	13.894	ug/l	96
91) 1,2-Dichlorobenzene	13.739	146	4623	1.823	ug/l	97

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

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