

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033022\
 Data File : VY008114.D
 Acq On : 30 Mar 2022 16:51
 Operator : SY/MD
 Sample : N2155-08RE
 Misc : 5.32g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 22L076-HRE

Quant Time: Mar 31 04:34:05 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032522S.M
 Quant Title : SW846 8260
 QLast Update : Sat Mar 26 03:24:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.789	168	130484	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.691	114	235834	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.490	117	224182	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.428	152	101071	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.143	65	73707	56.302	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	= 112.600%	
35) Dibromofluoromethane	7.722	113	36807	24.893	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	= 49.780%#	
50) Toluene-d8	10.179	98	283647	50.918	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	= 101.840%	
62) 4-Bromofluorobenzene	12.483	95	100550	51.161	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	= 102.320%	
Target Compounds						
					Qvalue	
16) Acetone	3.948	43	16037	53.594	ug/l #	88
25) 2-Butanone	6.984	43	11155	24.497	ug/l #	78
68) m/p-Xylenes	11.697	106	12405	3.521	ug/l	96
84) 1,2,4-Trimethylbenzene	13.123	105	20482	3.077	ug/l	99

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

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