

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY061722\
 Data File : VY009228.D
 Acq On : 17 Jun 2022 15:41
 Operator : KP/MD
 Sample : N3383-06
 Misc : 6.47g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 RB-04-(3-5)

Quant Time: Jun 17 22:54:22 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y061522S.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 16 02:54:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.783	168	186133	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.685	114	323624	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.483	117	301264	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.422	152	125026	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.136	65	112187	57.699	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	115.400%
35) Dibromofluoromethane	7.710	113	90507	44.607	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	89.220%
50) Toluene-d8	10.173	98	391732	50.718	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	=	101.440%
62) 4-Bromofluorobenzene	12.477	95	134344	48.947	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	=	97.900%
Target Compounds						
					Qvalue	
16) Acetone	3.942	43	24579	50.184	ug/l	91
17) Carbon Disulfide	4.186	76	19929	3.887	ug/l	97
64) Tetrachloroethene	10.715	164	13464	4.824	ug/l	95
65) Chlorobenzene	11.514	112	13423	2.057	ug/l	97
84) 1,2,4-Trimethylbenzene	13.117	105	8352	1.082	ug/l	92

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

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