

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071924\
 Data File : VY018877.D
 Acq On : 19 Jul 2024 13:15
 Operator : SY/MD
 Sample : P3246-19
 Misc : 6.92g/5.0mL/MSVOA_Y/SOIL/A
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 WC-CTW2-BL-04-G

Quant Time: Jul 19 16:51:34 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y070924S.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 09 14:28:51 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.783	168	80064	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.685	114	137917	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.483	117	123274	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.422	152	50656	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.130	65	38947	49.435	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	98.860%
35) Dibromofluoromethane	7.703	113	42111	48.751	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	97.500%
50) Toluene-d8	10.173	98	154207	47.121	ug/l	0.00
Spiked Amount	50.000	Range	58 - 134	Recovery	=	94.240%
62) 4-Bromofluorobenzene	12.477	95	49333	44.142	ug/l	0.00
Spiked Amount	50.000	Range	30 - 143	Recovery	=	88.280%
Target Compounds						
					Qvalue	
16) Acetone	3.942	43	3801	26.169	ug/l #	77
17) Carbon Disulfide	4.180	76	5644	2.606	ug/l #	60
20) Methylene Chloride	4.698	84	8153	5.493	ug/l	90
73) Isopropylbenzene	12.331	105	10313	2.782	ug/l #	62
86) p-Isopropyltoluene	13.367	119	8960	2.611	ug/l #	77

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

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