

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY071020\
 Data File : VY003170.D
 Acq On : 10 Jul 2020 17:54
 Operator : SY/MD
 Sample : L3255-07RE
 Misc : 6.80G/5ML/MSVOA Y/SOIL
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 RBR-07RE

Quant Time: Jul 11 06:36:44 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\82Y071020S.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 10 13:18:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.80	168	388816	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.69	114	582734	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.49	117	362680	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.42	152	86383	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.14	65	185683	56.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.16%	
35) Dibromofluoromethane	7.72	113	185644	55.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.72%	
50) Toluene-d8	10.18	98	594319	47.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.34%	
62) 4-Bromofluorobenzene	12.48	95	117071	25.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	51.98%	
Target Compounds						
11) Tert butyl alcohol	4.96	59	61172	172.383	ug/l	99
16) Acetone	3.95	43	15877	21.659	ug/l #	79
18) Methyl Acetate	4.48	43	7771	3.817	ug/l	95
25) 2-Butanone	6.99	43	6224	5.171	ug/l	98
73) Isopropylbenzene	12.33	105	11556	2.099	ug/l	98
80) 1,3,5-Trimethylbenzene	12.81	105	8222	1.792	ug/l	93
86) p-Isopropyltoluene	13.37	119	13269	2.554	ug/l #	85

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

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