Response Factor Report MSVOA_T

Method Path : Z:\VOASRV\HPCHEM1\MSVOA T\METHODS\
Method File : 82T112818TCLP.M
Title : SW846 8260
Last Update : Thu Nov 29 06:21:27 2018
Response Via : Initial Calibration

Calibration Files
10 =VT019101.D 5 =VT019100.D 20 =VT019102.D
50 =VT019103.D 100 =VT019104.D

		Compound	10	5	20	50	100	Avg	%RSD
1)	I	Pentafluorobenzene	----------ISTD----------						
2)	C	Vinyl Chloride	1.600	1.560	1.589	1.528	1.344	1.524	6.88#
3)	CM	1,1-Dichloroethene	1.150	1.123	1.026	1.019	0.924	1.049	8.61#
4)	T	2-Butanone	0.442	0.441	0.516	0.497	0.452	0.469	7.34
5)	C	Chloroform	1.371	1.486	1.386	1.364	1.259	1.373	5.88#
6)	S	1,2-Dichloroethane-	1.077	0.928	0.975	0.995	0.936	0.982	6.11
7)	I	1,4-Difluorobenzene	----------ISTD----------						
8)	S	Dibromofluoromethan	0.336	0.296	0.312	0.326	0.339	0.322	5.55
9)	T	Carbon Tetrachlorid	0.489	0.533	0.526	0.559	0.574	0.536	6.10
10)	TM	Benzene	1.624	1.743	1.727	1.807	1.820	1.744	4.48
11)	TM	1,2-Dichloroethane	0.610	0.660	0.637	0.649	0.629	0.637	3.00
12)	TM	Trichloroethene	0.311	0.352	0.328	0.353	0.363	0.341	6.20
13)	S	Toluene-d8	1.449	1.271	1.349	1.457	1.520	1.409	7.01
14)	S	4-Bromofluorobenzen	0.522	0.454	0.511	0.568	0.594	0.530	10.18
15)	I	Chlorobenzene-d5	----------ISTD----------						
16)	T	Tetrachloroethene	0.256	0.278	0.272	0.285	0.302	0.279	6.09
17)	PM	Chlorobenzene	1.080	1.173	1.099	1.150	1.168	1.134	3.72
18)	I	1,4-Dichlorobenzene-d	----------ISTD----------						

(#) = Out of Range