

Method Path : Z:\HPCHEM1\BNA_J\method\
 Method File : BJ060812.M
 Title : CWA degradation products by GCMS
 Last Update : Fri Jun 08 15:54:08 2012
 Response Via : Initial Calibration

Calibration Files

L6 =BJ000931.D L5 =BJ000932.D L4 =BJ000930.D L3 =BJ000933.D L2 =BJ000934.D
 L1 =BJ000935.D

	Compound	L6	L5	L4	L3	L2	L1	Avg	%RSD
1) I	1,4-Dichlorobenzen...	-----ISTD-----							
2) T	Dimethyl Disul...	0.734	0.800	0.692	0.557	0.640	0.652	0.679	12.30
3) T	DMMP	0.750	0.774	0.756	0.641	0.452	0.411	0.631	25.65
4) T	DIMP	1.933	1.818	1.855	1.729	1.484	1.436	1.709	11.97
5) T	1,4-Oxathiane	0.624	0.585	0.682	0.767	0.700	0.649	0.668	9.52
6) I	Naphthalene-d8	-----ISTD-----							
7) T	1,4-Dithiane	0.155	0.150	0.142	0.142	0.136	0.130	0.142	6.28
8) T	Thiodiglycol	0.164	0.144	0.112	0.106	0.103	0.086	0.119	24.32
9) T	Benzothiazole	0.628	0.589	0.516	0.530	0.505	0.528	0.549	8.77
10) S	Tripropylphosp...	0.903	0.847	0.811	0.759	0.688	0.679	0.781	11.43
11) T	p-Chlorophenyl...	0.381	0.361	0.322	0.342	0.331	0.332	0.345	6.40
12) S	2-Bromothioani...	0.156	0.146	0.133	0.132	0.126	0.127	0.137	8.62
13) T	p-Chlorophenyl...	0.313	0.245	0.214	0.174	0.179	0.176	0.217	25.35
14) T	p-Chlorophenyl...	0.368	0.341	0.266	0.249	0.242	0.280	0.291	17.77

(#) = Out of Range