

Method Path : Z:\HPCHEM1\BNA_J\METHOD\
 Method File : BJ072012.M
 Title : CWA degradation products by GCMS
 Last Update : Fri Jul 20 14:09:12 2012
 Response Via : Initial Calibration

Calibration Files

L6 =BJ001035.D L5 =BJ001036.D L4 =BJ001037.D
 L3 =BJ001038.D L2 =BJ001039.D L1 =BJ001040.D

	Compound	L6	L5	L4	L3	L2	L1	Avg	%RSD
1) I	1,4-Dichlorobenzene-d	-----ISTD-----							
2) T	Dimethyl Disulfid	0.649	0.899	0.801	0.680	0.670	0.791	0.748	13.07
3) T	DMMP	0.829	0.929	0.988	0.857			0.901	7.94
4) T	DIMP	1.865	1.857	1.933	1.698	1.189		1.708	17.71
5) T	1,4-Oxathiane	0.871	0.838	0.962	0.932	0.995	0.986	0.931	6.88
6) I	Naphthalene-d8	-----ISTD-----							
7) T	1,4-Dithiane	0.196	0.192	0.179	0.195	0.184	0.181	0.188	3.97
8) T	Thiodiglycol	0.183	0.168	0.157	0.155	0.156	0.138	0.160	9.55
9) T	Benzothiazole	0.585	0.544	0.544	0.514	0.466	0.437	0.515	10.69
10) S	Tripropylphosphat	0.833	0.796	0.746	0.761	0.644	0.610	0.732	11.91
11) T	p-Chlorophenylmet	0.343	0.324	0.312	0.310	0.299	0.306	0.316	4.99
12) S	2-Bromothioanisol	0.169	0.158	0.153	0.159	0.149	0.150	0.156	4.69
13) T	p-Chlorophenylmet	0.243	0.208	0.201	0.191	0.148	0.114	0.184	25.02
14) T	p-Chlorophenylmet	0.344	0.299	0.299	0.331	0.317	0.301	0.315	6.00

(#) = Out of Range