

Method Path : Z:\HPCHEM1\BNA_J\METHOD\
 Method File : BJ112712.M
 Title : CWA degradation products by GCMS
 Last Update : Wed Nov 28 14:20:09 2012
 Response Via : Initial Calibration

Calibration Files

L6 =BJ001290.D L5 =BJ001291.D L4 =BJ001292.D
 L3 =BJ001293.D L2 =BJ001294.D L1 =BJ001295.D

	Compound	L6	L5	L4	L3	L2	L1	Avg	%RSD
1) I	1,4-Dichlorobenzene-d	-----ISTD-----							
2) T	Dimethyl Disulfid	0.737	1.031	0.892	1.311	0.774	0.792	0.923	23.59
3) T	DMMP	1.014	1.100	1.034	0.898	0.616	0.520	0.864	27.80
4) T	DIMP	2.296	2.433	2.327	2.057	1.655	1.533	2.050	18.35
5) T	1,4-Oxathiane	1.045	1.201	1.261	1.345	1.345	1.047	1.207	11.29
6) I	Naphthalene-d8	-----ISTD-----							
7) T	1,4-Dithiane	0.267	0.252	0.249	0.243	0.230	0.212	0.242	7.84
8) T	Thiodiglycol	0.237	0.226	0.182	0.189	0.186	0.211	0.205	11.25
9) T	Benzothiazole	0.581	0.575	0.546	0.564	0.521	0.565	0.559	3.96
10) S	Tripropylphosphat	1.182	1.079	1.046	0.920	0.868	0.893	0.998	12.40
11) T	p-Chlorophenylmet	0.373	0.353	0.338	0.350	0.350	0.336	0.350	3.79
12) S	2-Bromothioanisol	0.185	0.170	0.162	0.167	0.154	0.160	0.166	6.48
13) T	p-Chlorophenylmet	0.278	0.268	0.240	0.215	0.196	0.223	0.236	13.42
14) T	p-Chlorophenylmet	0.409	0.378	0.329	0.317	0.308	0.315	0.343	11.98

(#) = Out of Range