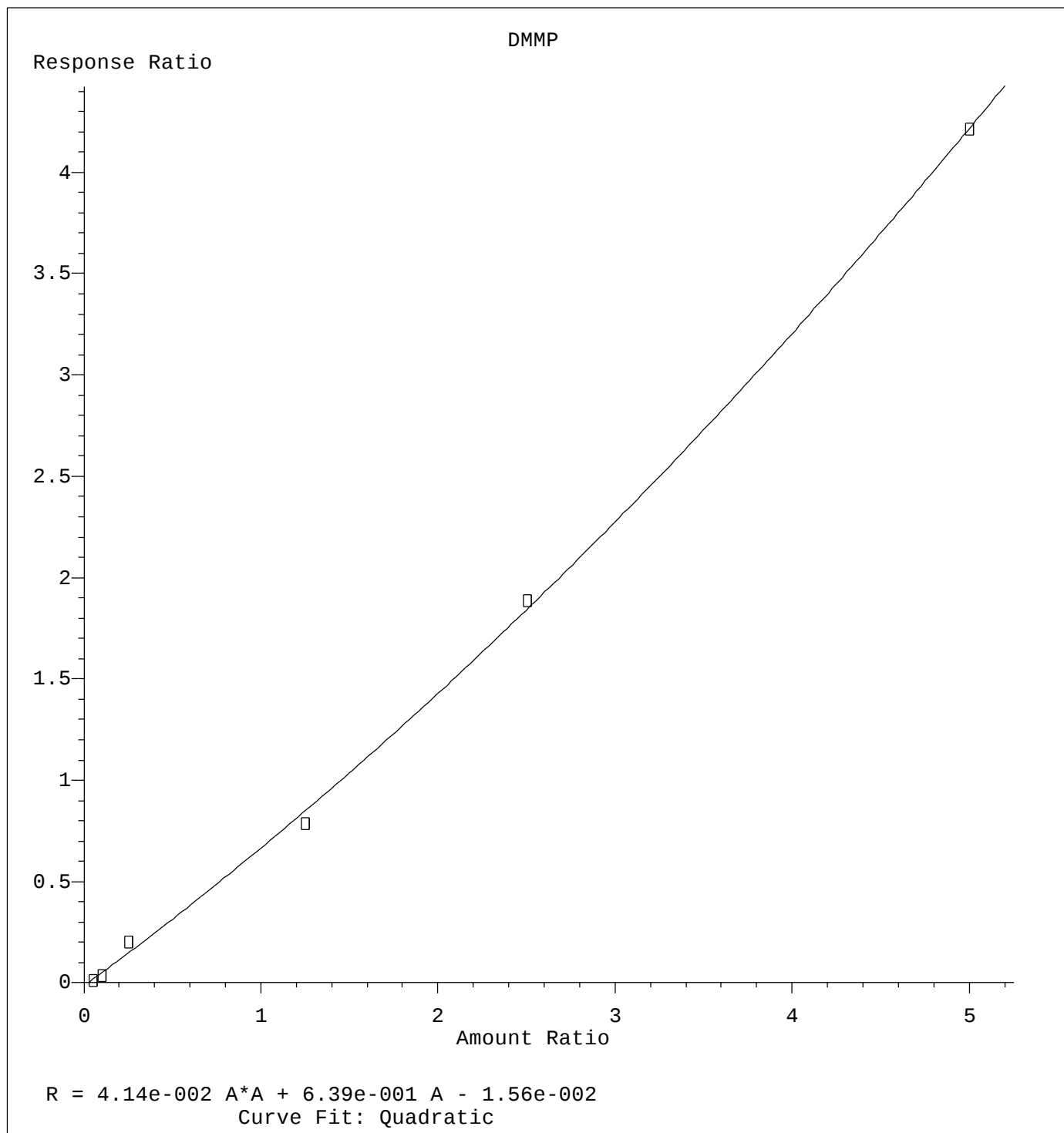


Method : H:\VOL1\GCMSDATA\J\METHODS\BJ091112.M (RTE Integrator)
 Title : CWA degradation products by GCMS
 Last Update : Tue Sep 11 15:38:38 2012
 Response via : Initial Calibration

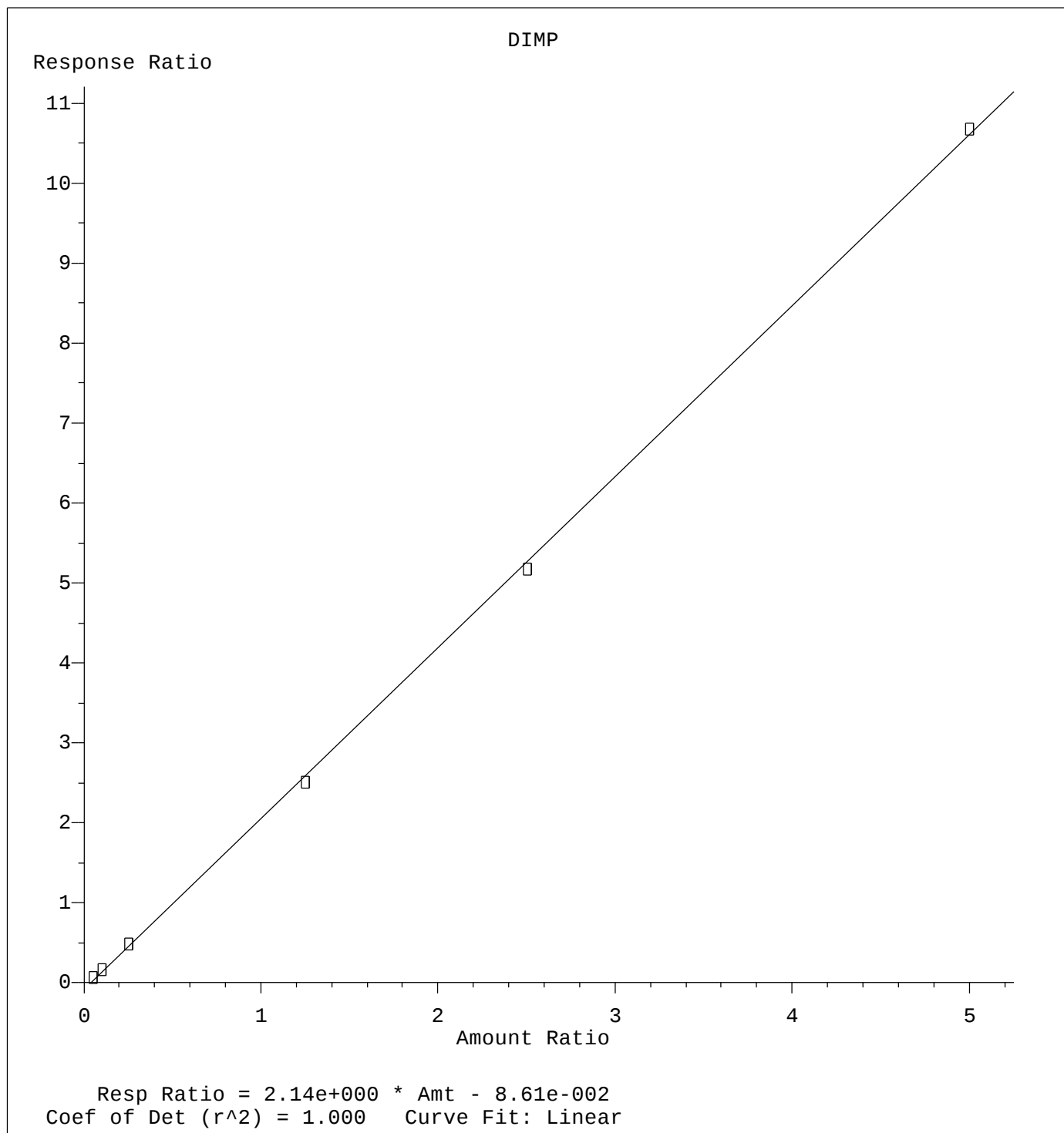
Calibration Files

L6 =BJ001103.D L5 =BJ001104.D L4 =BJ001105.D
 L3 =BJ001106.D L2 =BJ001107.D L1 =BJ001108.D

	Compound	L6	L5	L4	L3	L2	L1	Avg	%RSD
1) I	1,4-Dichlorobenzene-d	-----ISTD-----							
2) T	Dimethyl Disulfide		0.846	0.745	1.023	0.845	0.942	0.880	12.04
3) T	DMMP	0.842	0.753	0.627	0.796	0.343	0.191	0.592	44.95
4) T	DIMP	2.134	2.066	2.006	1.924	1.581	1.328	1.840	17.20
5) T	1,4-Oxathiane	0.756	0.877	0.916	0.859	0.837	0.798	0.840	6.81
6) I	Naphthalene-d8	-----ISTD-----							
7) T	1,4-Dithiane	0.171	0.166	0.161	0.182	0.173	0.170	0.171	4.13
8) T	Thiodiqlvcol	0.180	0.159	0.149	0.167	0.151	0.167	0.162	7.02
9) T	Benzothiazole	0.607	0.555	0.525	0.485	0.556	0.531	0.543	7.47
10) S	Tripropylphosphate	1.004	0.851	0.831	0.796	0.864	0.836	0.864	8.40
11) T	p-Chlorophenylmethy	0.345	0.321	0.303	0.316	0.313	0.329	0.321	4.46
12) S	2-Bromothioanisole	0.164	0.154	0.147	0.154	0.151	0.156	0.154	3.72
13) T	p-Chlorophenylmethy	0.259	0.206	0.180	0.190	0.193	0.206	0.206	13.53
14) T	p-Chlorophenylmethy	0.395	0.300	0.268	0.332	0.260	0.285	0.307	16.45



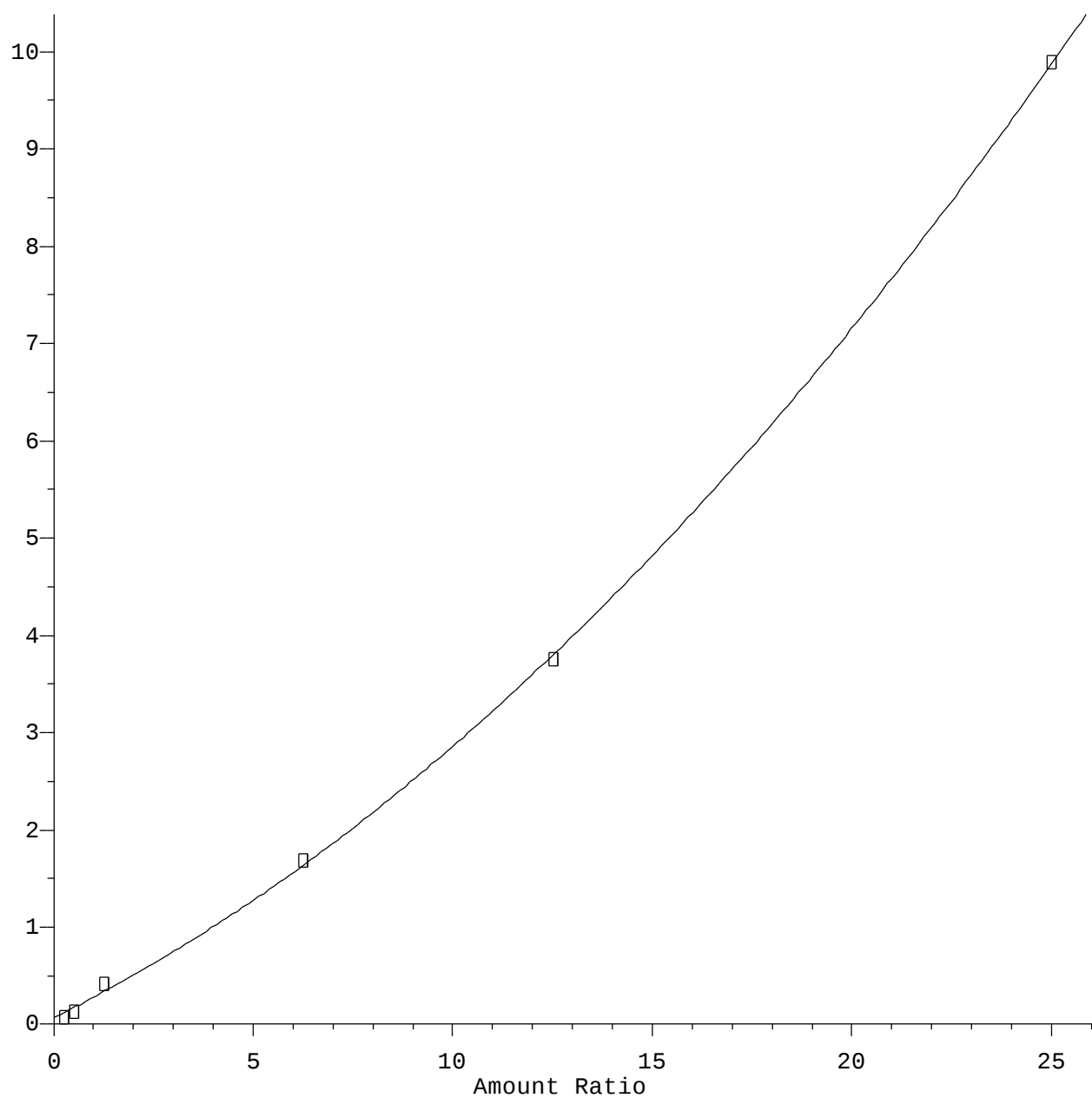
Method Name: H:\VOL1\GCMSDATA\J\METHODS\BJ091112.M
Calibration Table Last Updated: Tue Sep 11 15:38:38 2012



Method Name: H:\VOL1\GCMSDATA\J\METHODS\BJ091112.M
Calibration Table Last Updated: Tue Sep 11 15:38:38 2012

p-Chlorophenylmethylsulfone

Response Ratio



$R = 7.56e-003 A^2 + 2.03e-001 A + 6.76e-002$
Curve Fit: Quadratic

Method Name: H:\VOL1\GCMSDATA\J\METHODS\BJ091112.M
Calibration Table Last Updated: Tue Sep 11 15:38:38 2012