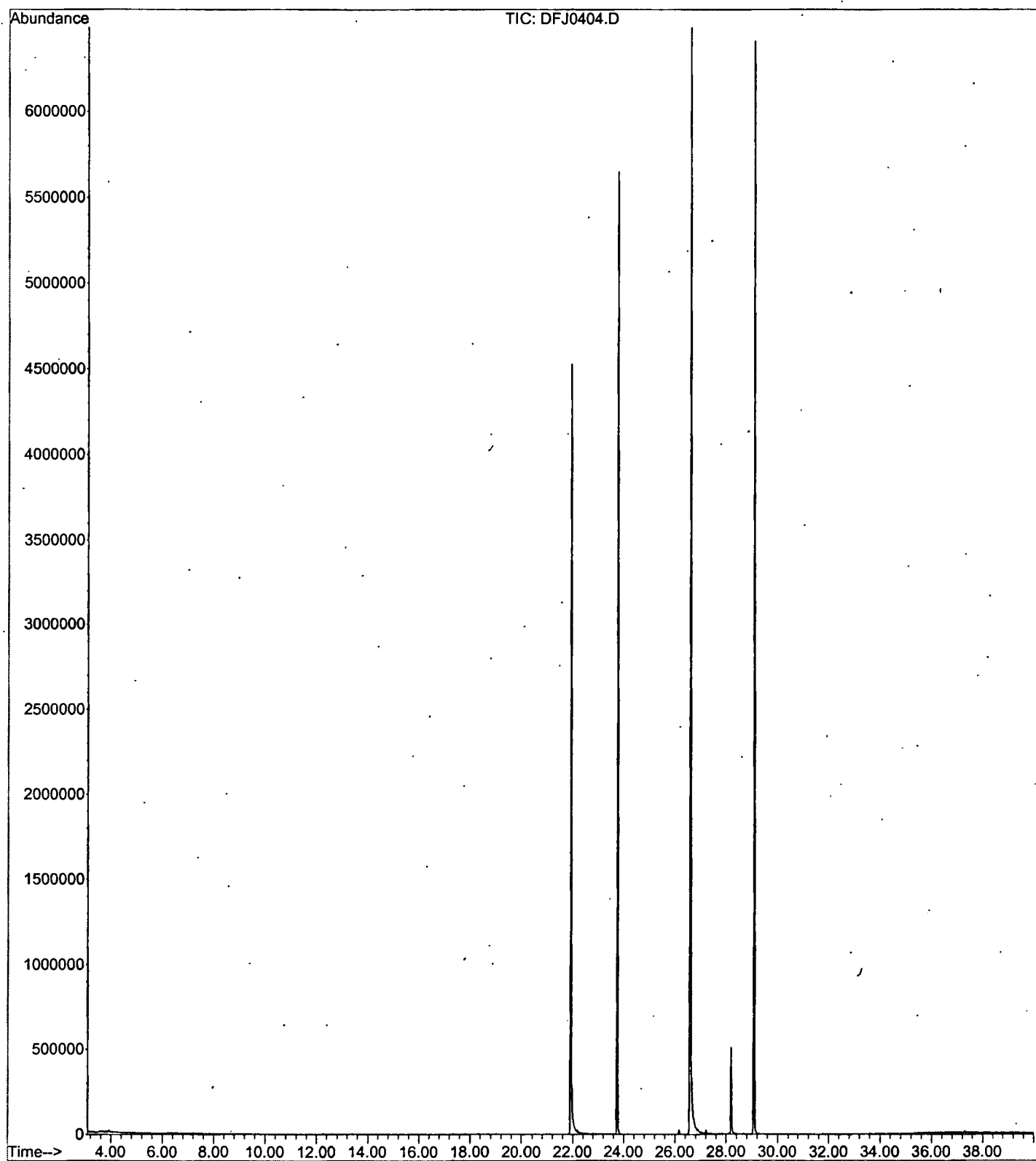


File : C:\MSDCHEM\1\DATA\040402B\DFJ0404.D
 Operator : Bohlk
 Acquired : 4 Apr 2002 7:13 pm using AcqMethod 508SCAN
 Instrument : Instrumen
 Sample Name: DFTPP
 Misc Info :
 Vial Number: 1

Unit, Cali.



DFTPP

Data File : C:\MSDCHEM\1\DATA\040402B\DFJ0404.D
Acq On : 4 Apr 2002 7:13 pm
Sample : DFTPP
Misc :
MS Integration Params: BENZO.P

Vial: 1
Operator: Bohlk
Inst : Instrumen
Multiplr: 1.00

Method : C:\MSDCHEM\1\5973N\5080402.M (RTE Integrator)
Title : Quantitation For Method 508gcscan

Spectrum Information: Scan 3518

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
51	198	30	60	46.1	314624	PASS
68	69	0.00	2	1.6	5291	PASS
69	198	0.00	100	48.6	332032	PASS
70	69	0.00	2	0.6	1910	PASS
127	198	40	60	51.6	352384	PASS
197	198	0.00	1	0.0	0	PASS
198	198	100	100	100.0	683072	PASS
199	198	5	9	6.4	44024	PASS
275	198	10	30	20.4	139200	PASS
365	198	1	100	1.9	13272	PASS
441	443	0.01	100	79.5	58200	PASS
442	198	40	110	54.4	371712	PASS
443	442	17	23	19.7	73200	PASS

DFJ0404.D 5080402.M Fri Apr 05 06:17:53 2002

Response Factor Report Instrumen

Method : C:\MSDCHEM\1\5973N\5080402.M (RTE Integrator)
 Title : Quantitation For Method 508gcscan
 Last Update : Fri Apr 05 05:56:22 2002
 Response via : Initial Calibration

Calibration Files

6 =C100J0404.D 1 =C5J0404.D 2 =C10J0404.D
 3 =C25J0404.D 4 =C40J0404.D 5 =C80J0404.D

Compound	6	1	2	3	4	5	Avg	%RSD

1) I 1,4-Dichlorobenzene-d	-----ISTD-----							
2) T N-nitroso-dimet	0.813	0.640	0.716	0.725	0.870	0.815	0.763	11.03
3) bis(2-chloroeth	0.973	1.263	1.221	1.049	1.142	1.020	1.112	10.42
4) 1,3 - Dichlorob	0.881	1.180	1.099	0.919	1.037	0.918	1.006	11.85
5) 1,4 - Dichlorob	0.865	1.132	1.066	0.910	1.021	0.907	0.983	10.72
6) 1,2 - Dichlorob	0.807	1.058	0.990	0.845	0.949	0.840	0.915	10.86
7) 2,2' - oxybis	1.416	2.089	1.808	1.594	1.817	1.574	1.716	13.85
8) N-Nitroso-di-n-	0.745	0.925	0.908	0.815	0.869	0.816	0.846	7.94
9) I Hexachloroethan	0.297	0.452	0.433	0.372	0.397	0.315	0.378	16.50
10) Naphthalene-d8	-----ISTD-----							
11) Nitrobenzene	0.301	0.338	0.344	0.290	0.333	0.305	0.318	7.06
12) Isophorone	0.573	0.568	0.585	0.513	0.621	0.590	0.575	6.22
13) bis(2-Chloroeth	0.331	0.362	0.369	0.326	0.368	0.336	0.348	5.69
14) 1,2,4-Trichloro	0.155	0.197	0.191	0.158	0.176	0.161	0.173	10.37
15) Naphthalene	0.578	0.829	0.781	0.642	0.698	0.614	0.690	14.21
16) Hexachlorobutad	0.079	0.098	0.092	0.078	0.088	0.080	0.086	9.41
17) Acenaphthene-d10	-----ISTD-----							
18) Hexachlorocyclo	0.123	0.078	0.091	0.099	0.128	0.127	0.107	19.75
19) T 2-chloronaphtha	0.616	0.782	0.746	0.639	0.722	0.635	0.690	9.98
20) I Dimethylphthala	0.744	0.881	0.847	0.729	0.852	0.769	0.804	7.96
21) T 2,6-Dinitrotolu	0.172	0.155	0.174	0.158	0.187	0.174	0.170	6.88
22) Acenaphthylene	0.821	1.022	1.054	0.910	1.008	0.873	0.948	9.85
23) Acenaphthene	0.573	0.808	0.747	0.621	0.683	0.596	0.671	13.72
24) 2,4-Dinitrotolu	0.238	0.200	0.237	0.222	0.263	0.243	0.234	9.04
25) Diethylphthalat	0.562	0.817	0.756	0.636	0.701	0.597	0.678	14.41
26) 4-Chlorophenyl-	0.313	0.407	0.375	0.322	0.362	0.322	0.350	10.70
27) Fluorene	0.630	0.945	0.875	0.713	0.766	0.660	0.765	16.17
28) Azobenzen/ 1,2-	1.077	1.344	1.300	1.108	1.252	1.158	1.207	8.97
29) Phenanthrene-d10	-----ISTD-----							
30) N-Nitrosodiphen	0.315	0.396	0.385	0.319	0.363	0.328	0.351	10.03
31) 4-Bromophenyl-p	0.115	0.128	0.127	0.107	0.126	0.115	0.120	7.13
32) Hexachlorobenze	0.110	0.128	0.119	0.104	0.118	0.111	0.115	7.60
33) Phenanthrene	0.656	0.862	0.813	0.674	0.750	0.670	0.737	11.59
34) Anthracene	0.648	0.815	0.771	0.664	0.745	0.664	0.718	9.58
35) Di-n-butylphtha	0.761	0.905	0.901	0.781	0.876	0.787	0.835	7.88
36) Fluoranthene	0.718	0.810	0.822	0.724	0.803	0.737	0.769	6.16
37) s 2,4-DDD (SURR)	0.187	0.223	0.208	0.200	0.183	0.190	0.199	7.74
38) Chrysene-d12	-----ISTD-----							
39) Pyrene	1.179	1.217	1.193	1.001	1.229	1.193	1.169	7.21
40) Butylbenzylphth	0.521	0.475	0.510	0.455	0.551	0.526	0.506	6.97
41) Benzo (a) antra	0.820	0.811	0.796	0.706	0.852	0.821	0.801	6.26
42) Bis (2-ethylhex	0.655	0.618	0.666	0.607	0.706	0.660	0.652	5.48
43) Chrysene	0.755	0.809	0.800	0.673	0.803	0.765	0.767	6.69
44) Perylene-d12	-----ISTD-----							
45) Di-n-Octylphtha	2.117	1.533	1.855	1.733	2.185	2.078	1.917	13.26
46) Benzo (b) fluor	1.162	1.010	1.034	0.951	1.184	1.147	1.081	8.85
47) Benzo (k) fluor	1.147	1.089	1.229	1.016	1.252	1.159	1.149	7.63

(#) = Out of Range

5080402.M

Fri Apr 05 05:58:38 2002

Page 1

Response Factor Report Instrumen

Method : C:\MSDCHEM\1\5973N\5080402.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Fri Apr 05 05:56:22 2002
Response via : Initial Calibration

Calibration Files

6 =C100J0404.D 1 =C5J0404.D 2 =C10J0404.D
3 =C25J0404.D 4 =C40J0404.D 5 =C80J0404.D

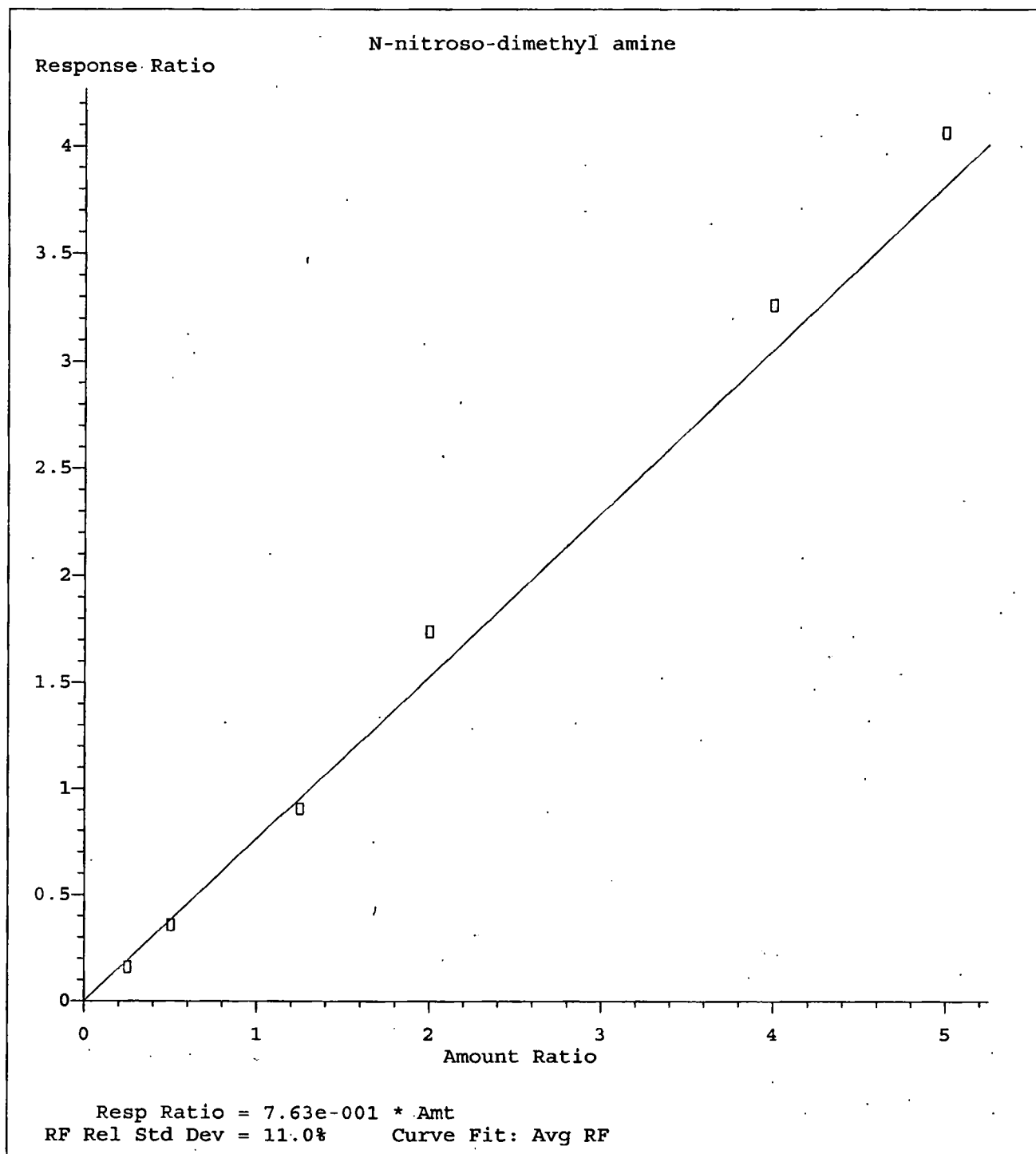
	Compound	6	1	2	3	4	5	Avg	%RSD
48)	Benzo (a) pyren	0.999	0.906	0.927	0.850	1.030	1.007	0.953	7.33
49)	Indeno (1,2,3-c	0.618	0.350	0.423	0.462	0.592	0.600	0.508	21.90
50)	Dibenz (a,h) an	0.577	0.448	0.422	0.452	0.562	0.566	0.505	14.10
51)	Benzo (g,h,i) p	0.585	0.453	0.457	0.461	0.582	0.557	0.516	12.64

(#) = Out of Range

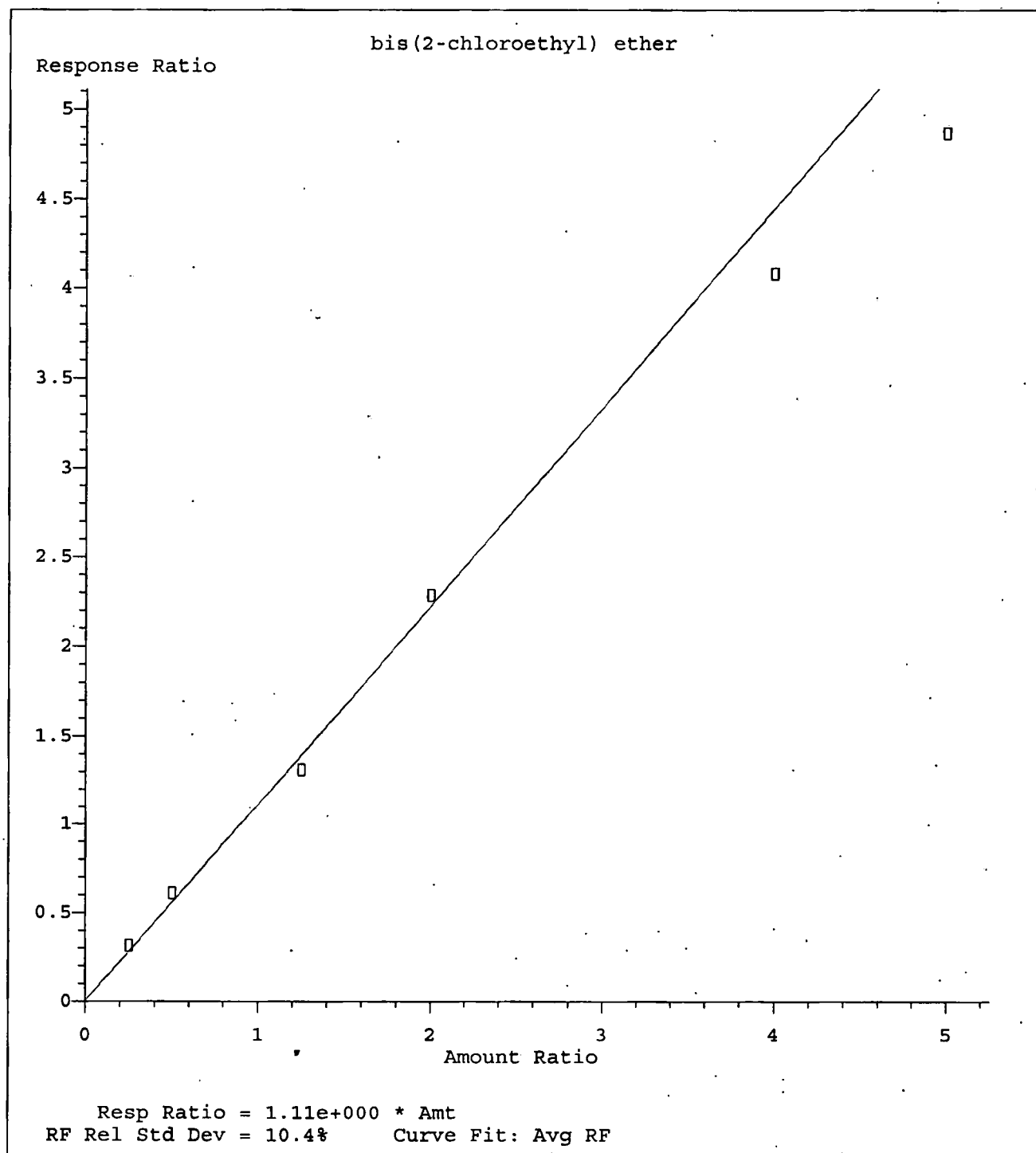
5080402.M

Fri Apr 05 05:58:38 2002

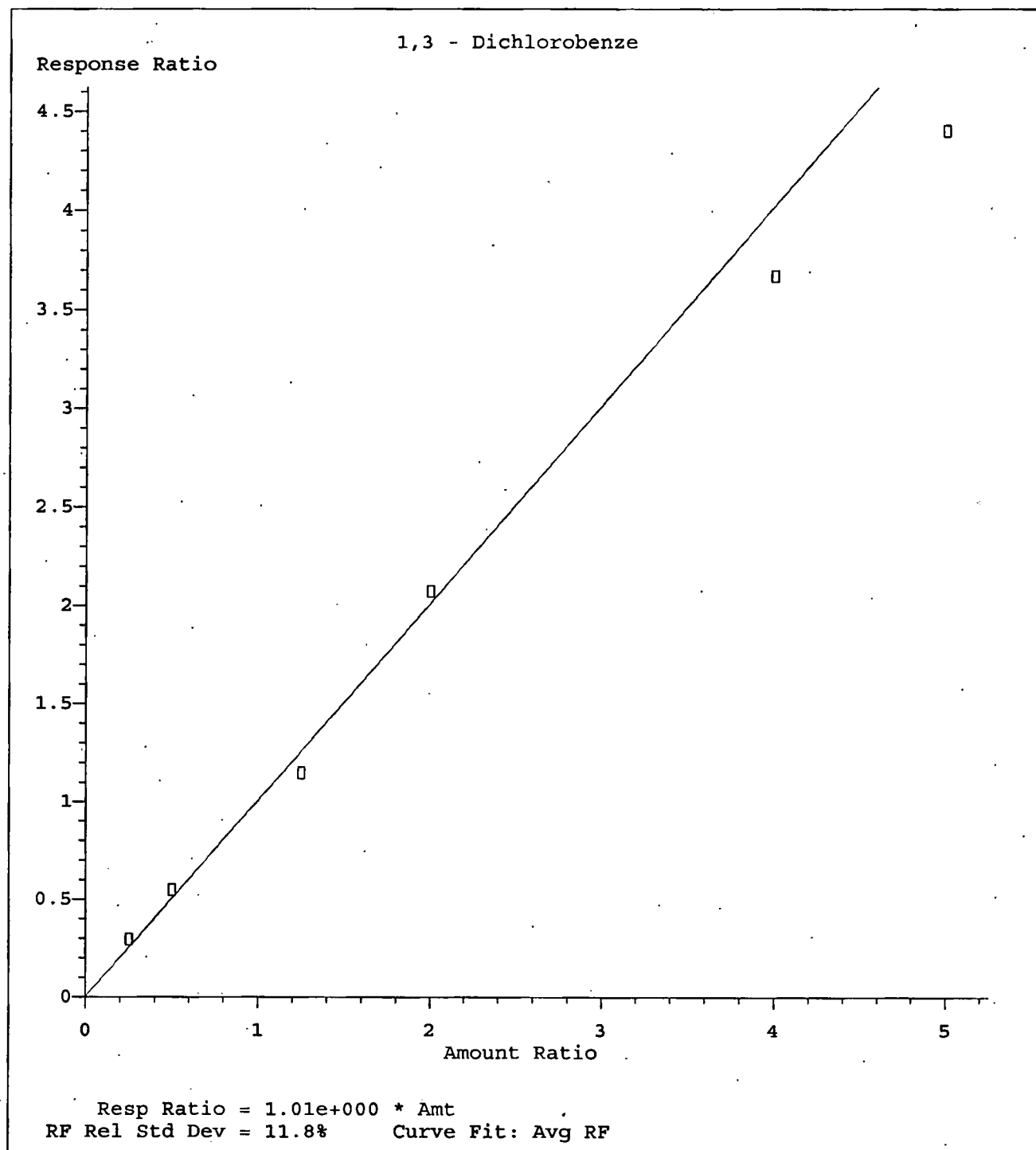
Page 2



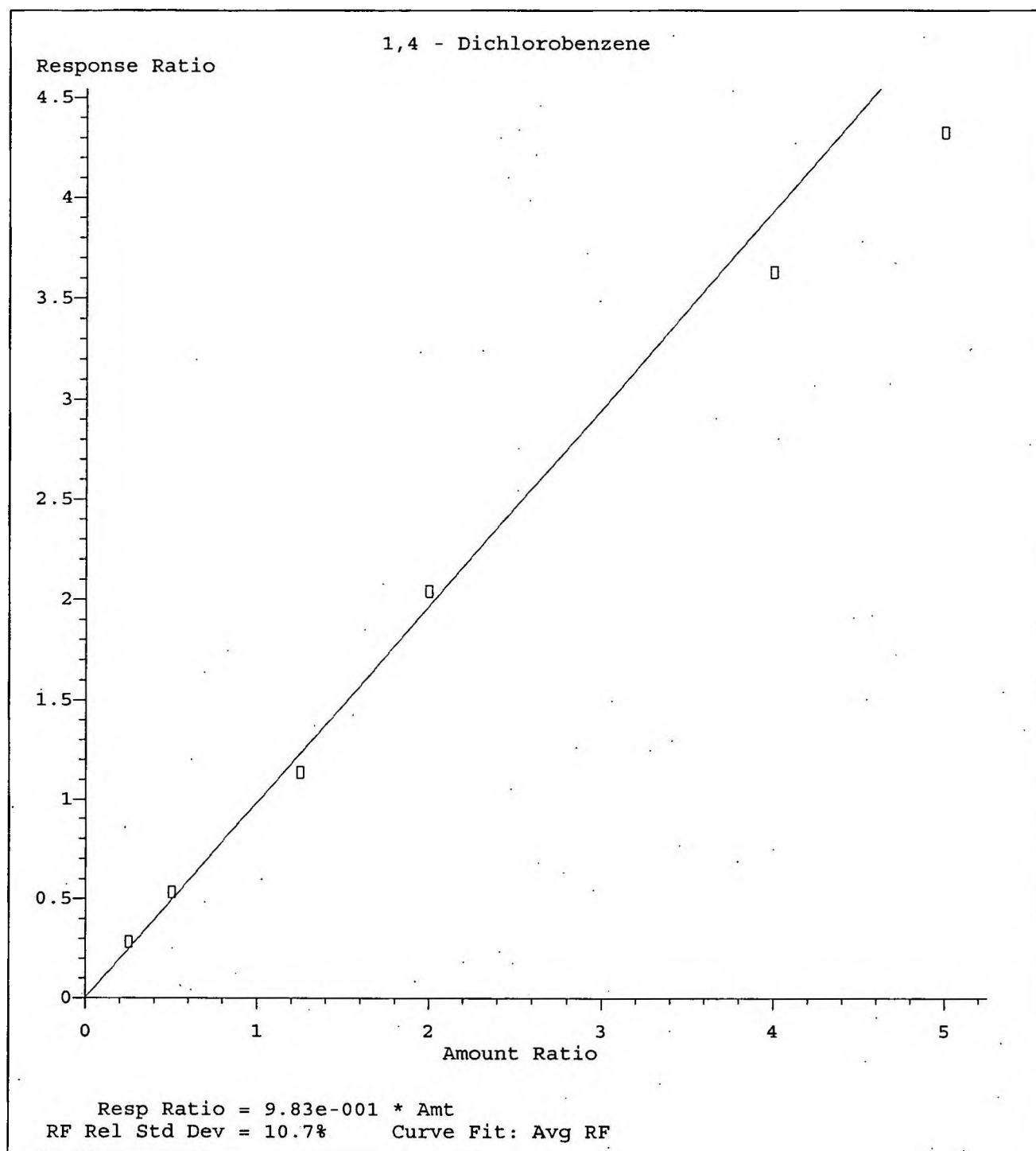
Method Name: C:\MSDCHEM\1\5973N\5080402.M
Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



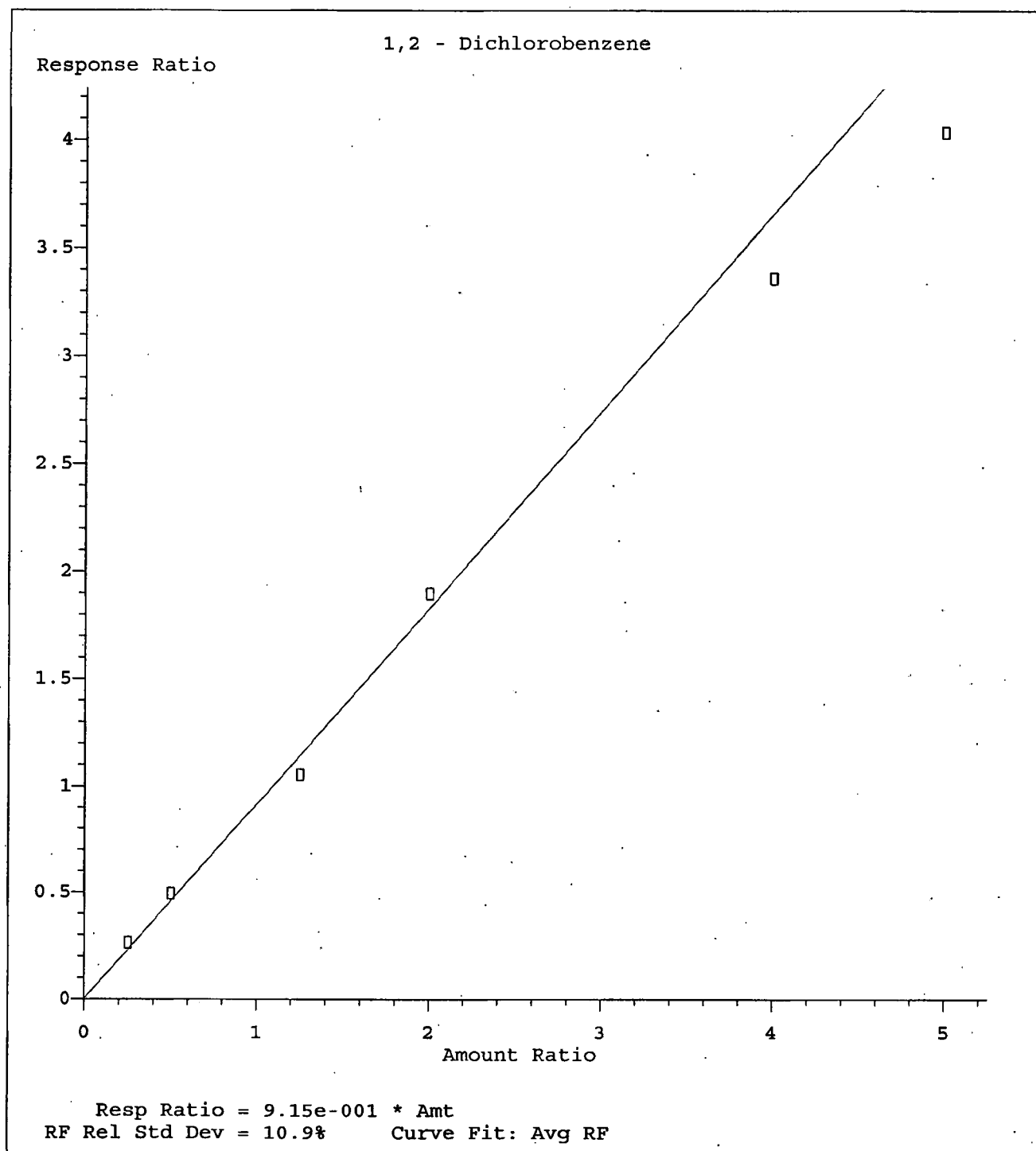
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 Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



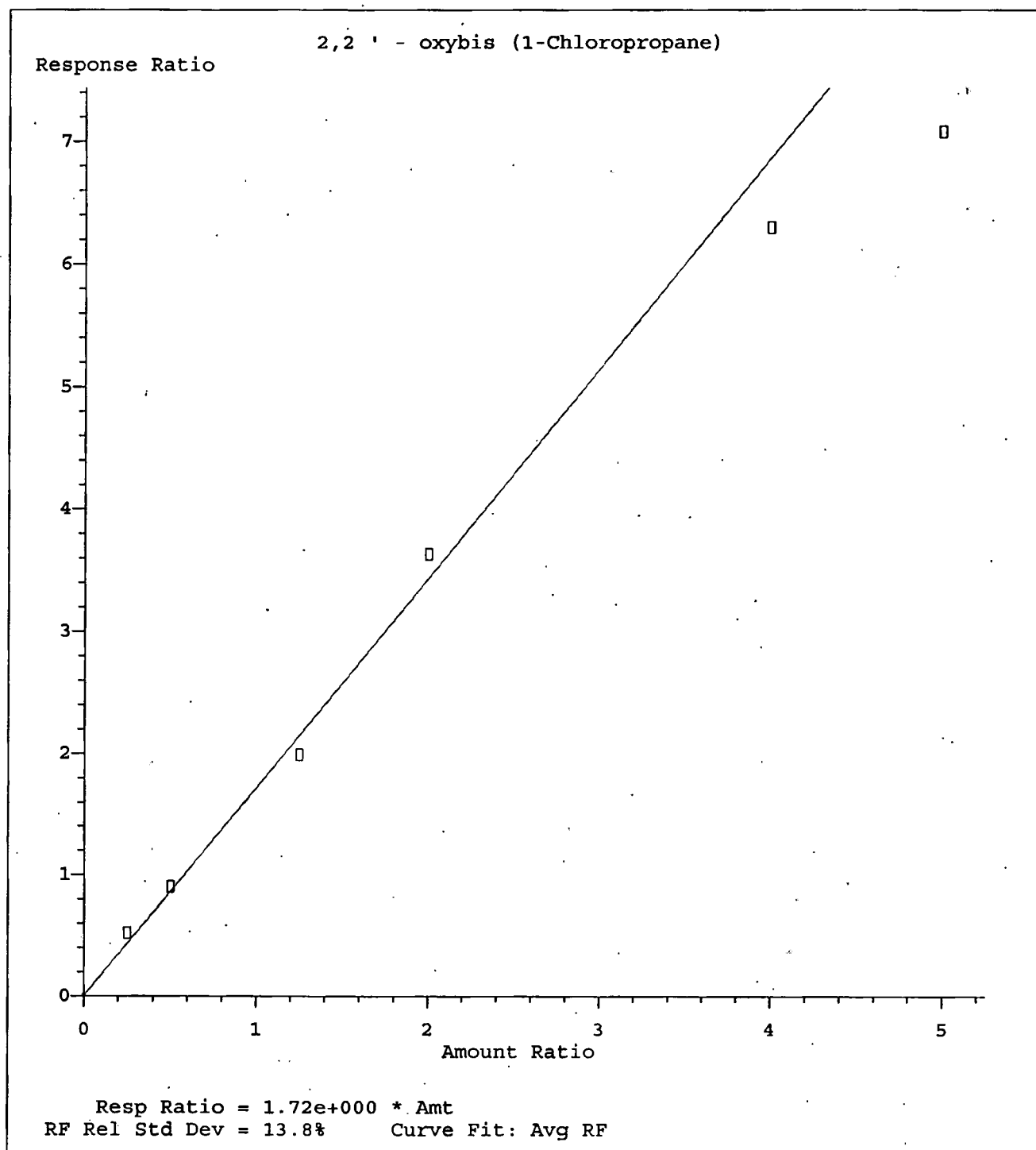
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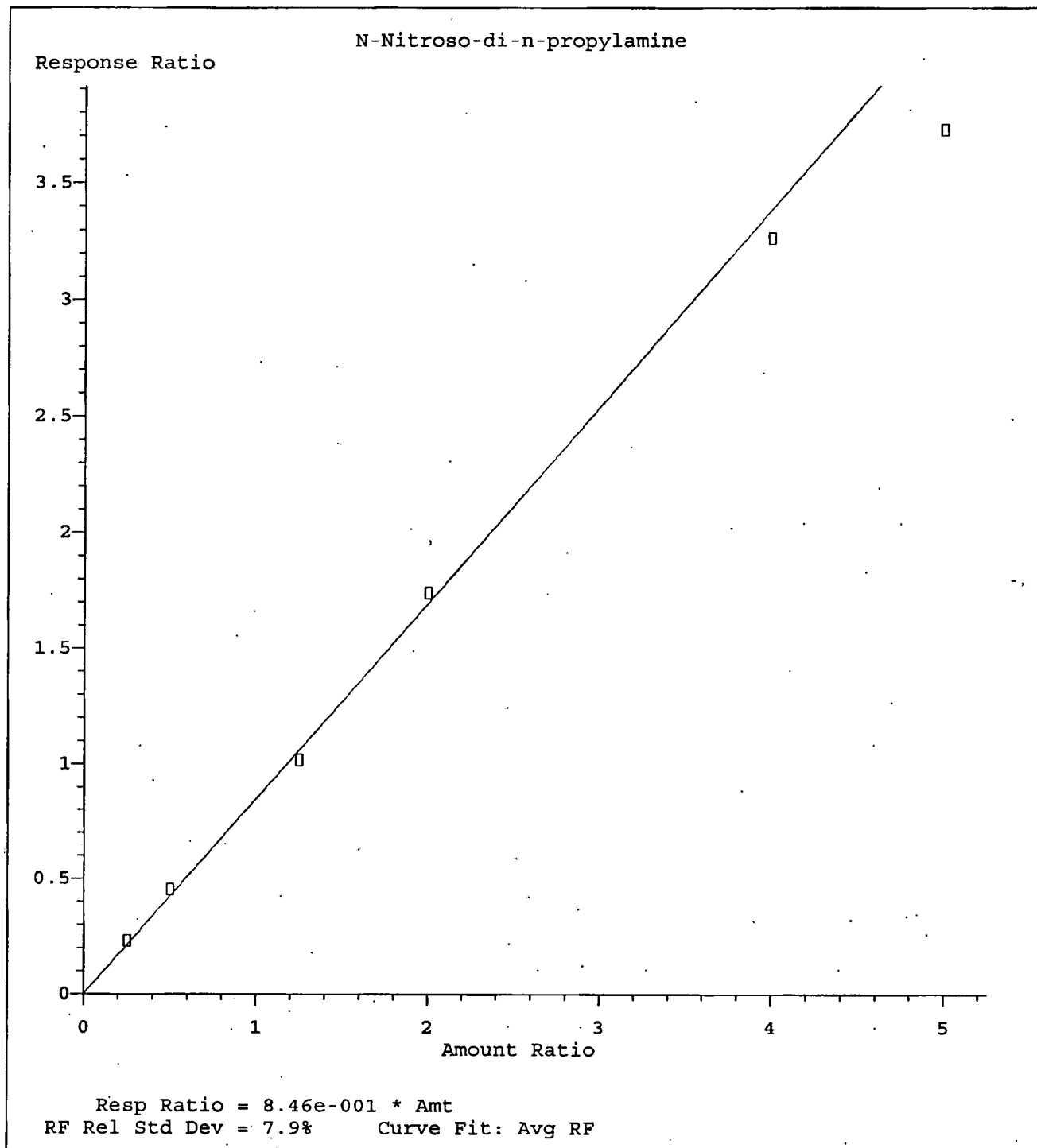
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



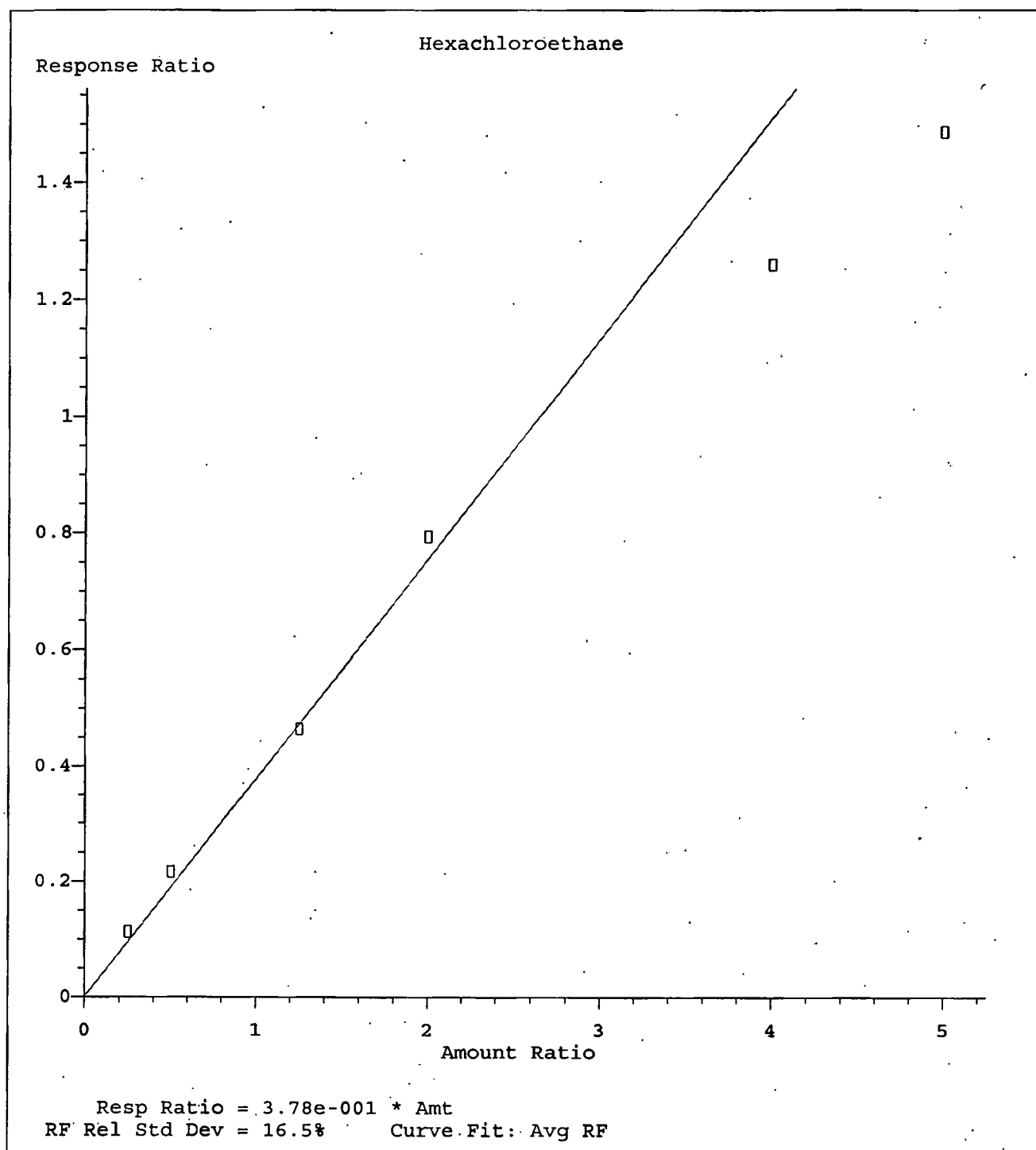
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



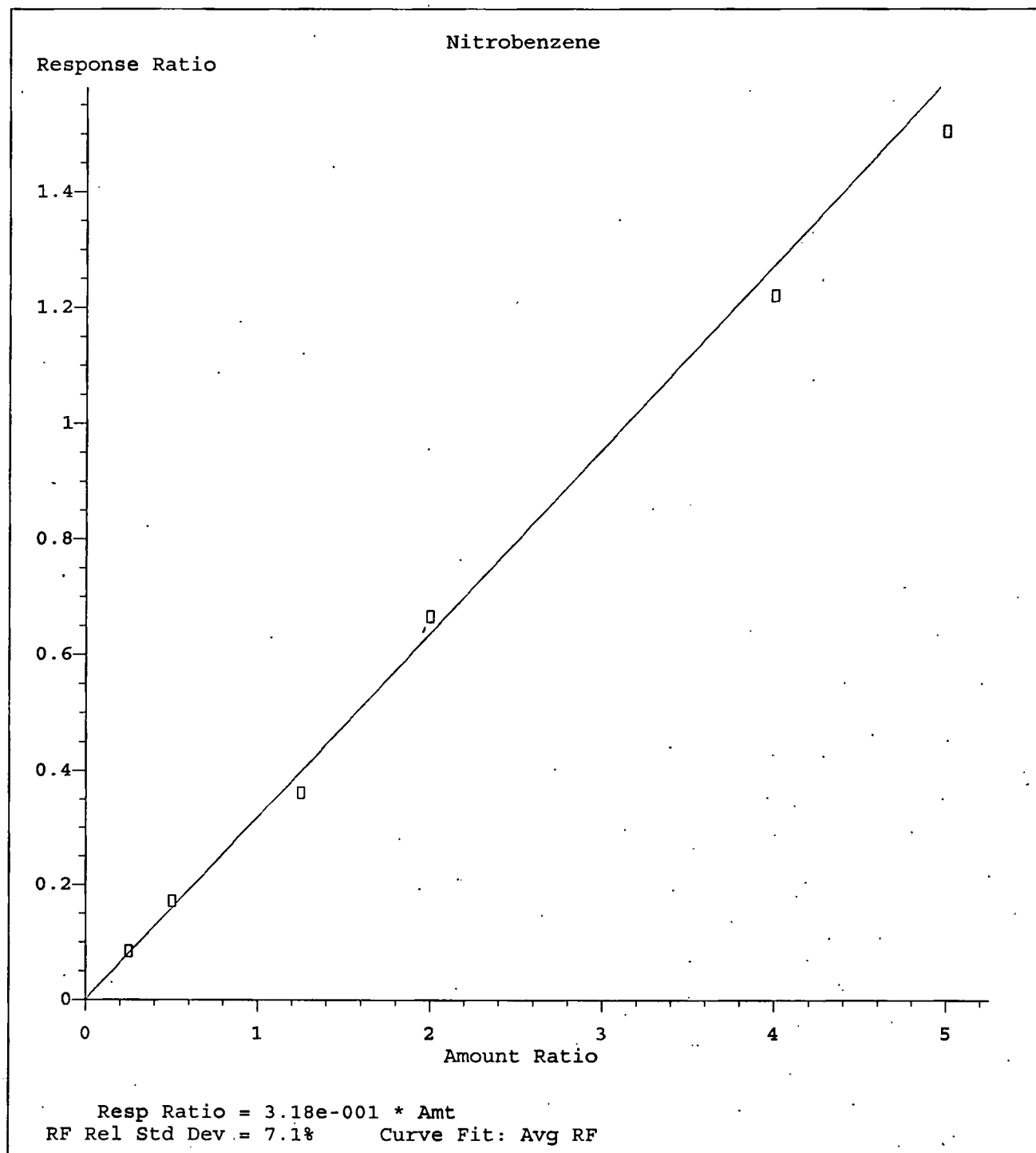
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 Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



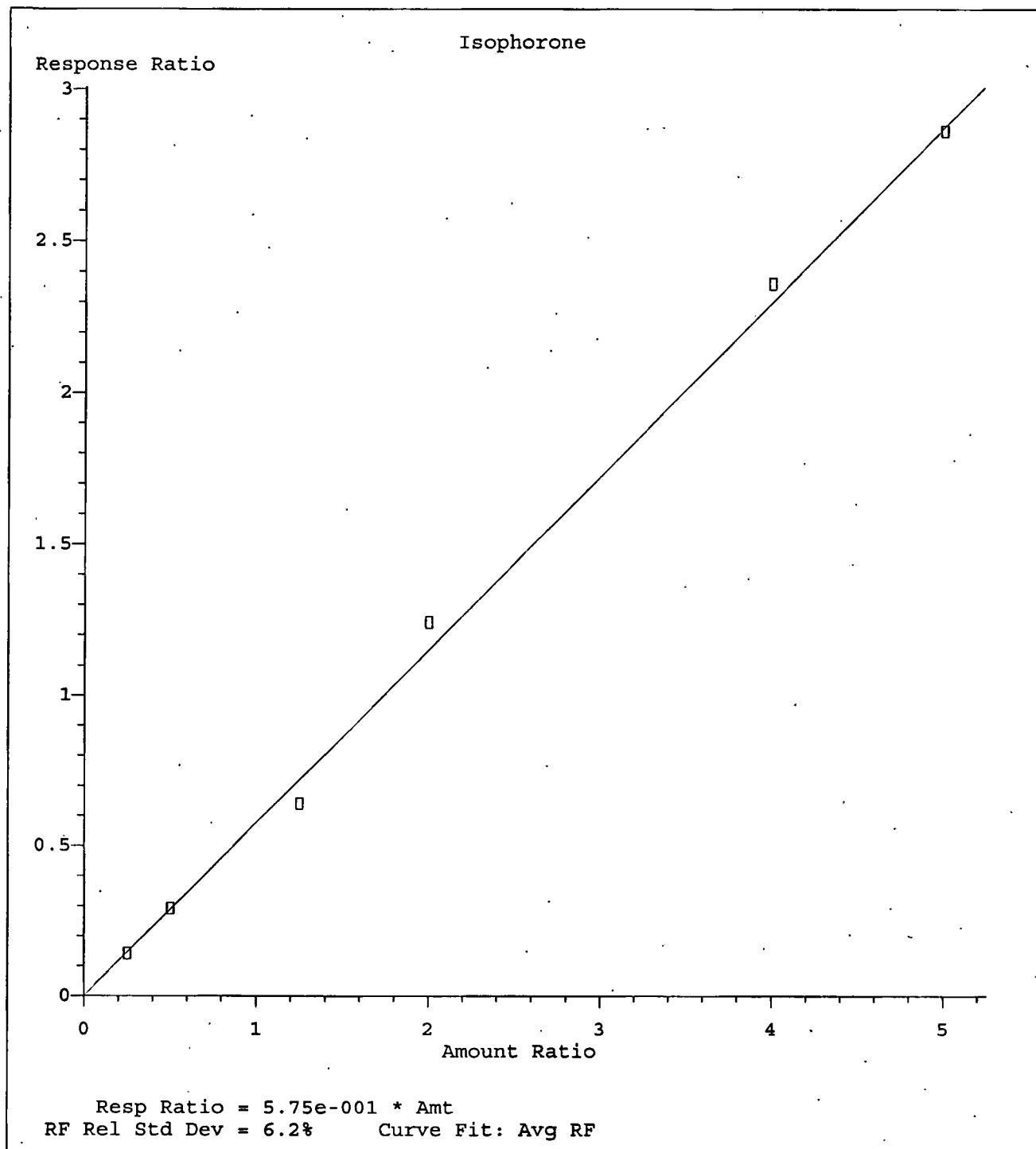
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



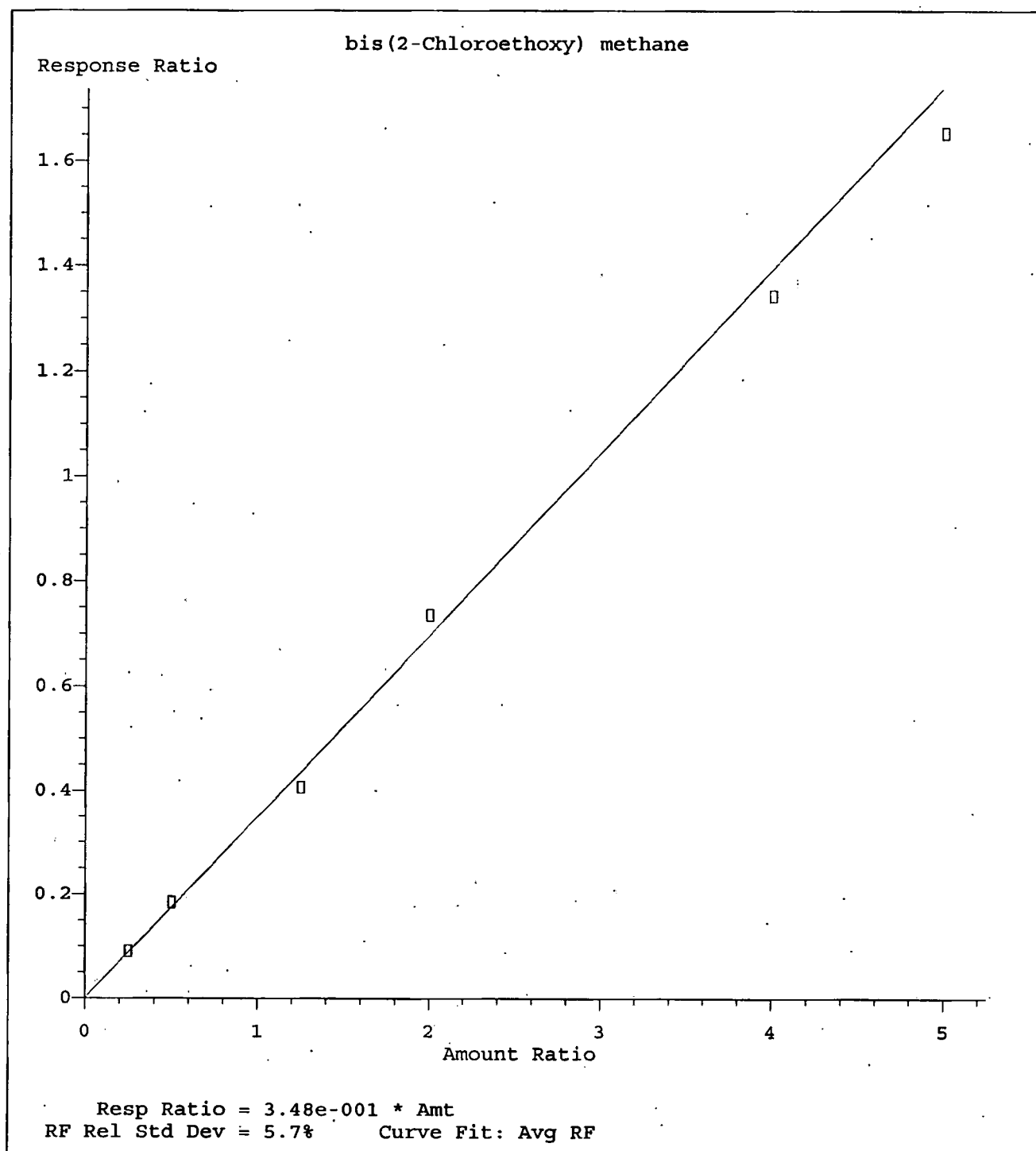
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 Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



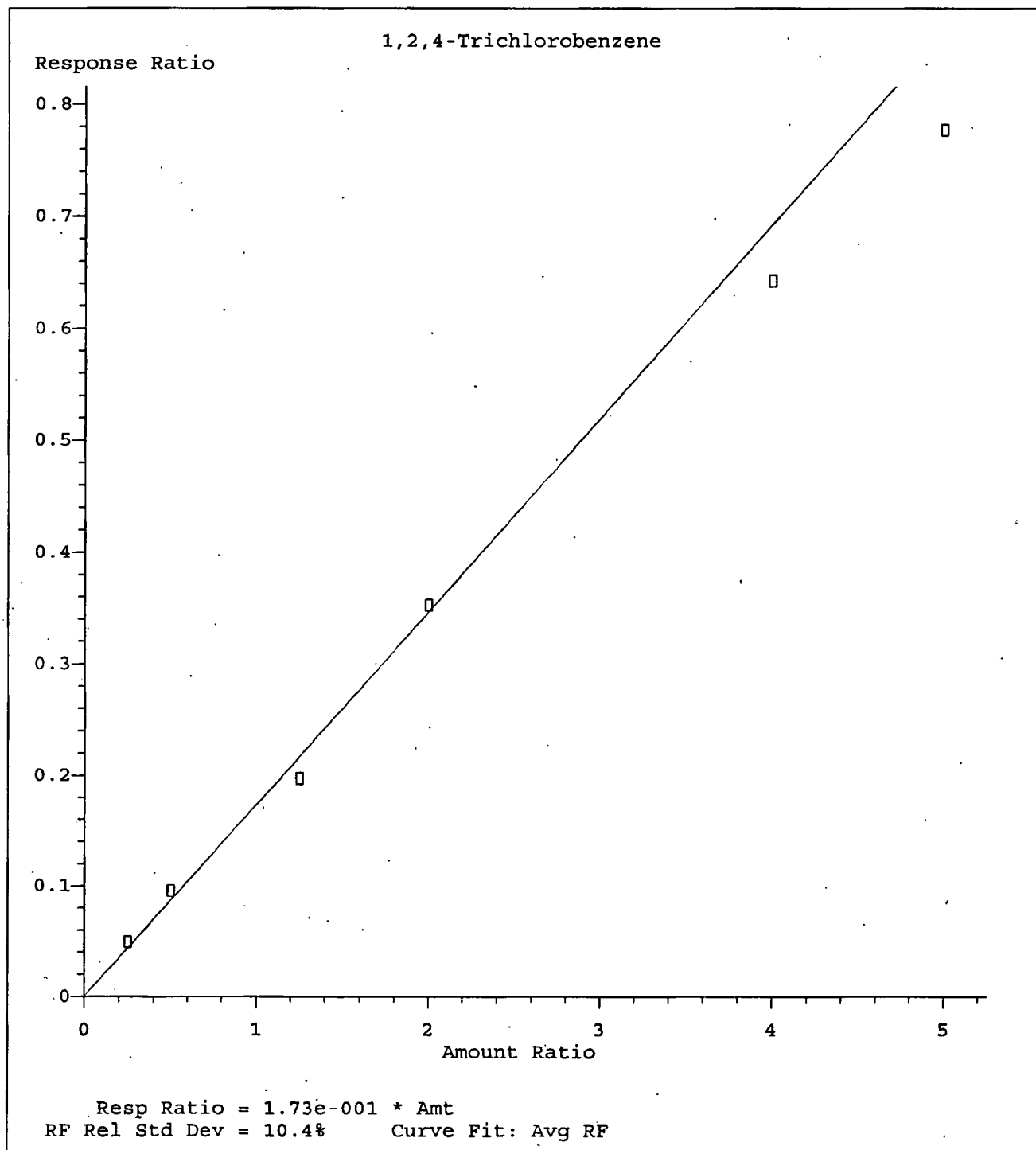
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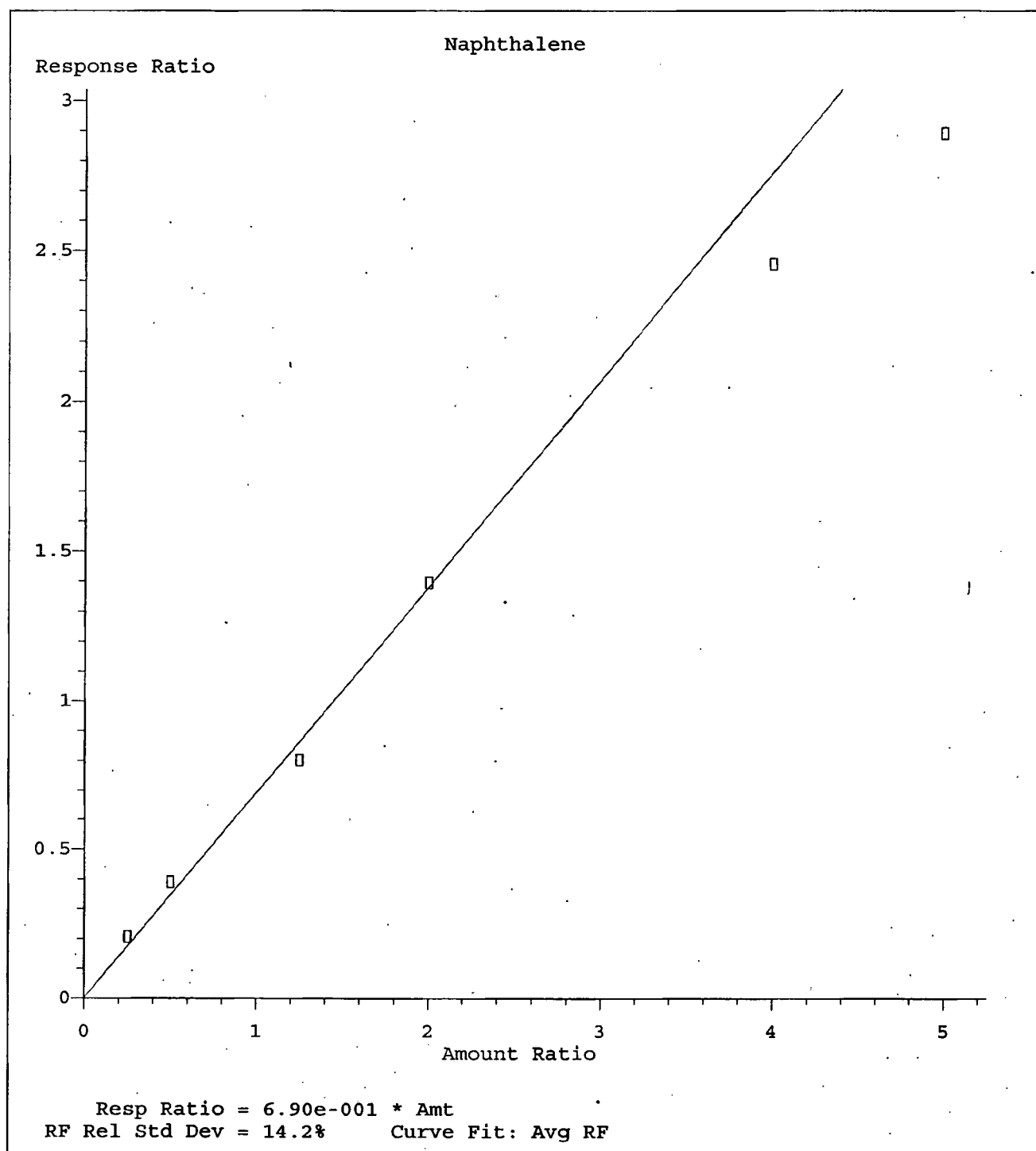
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



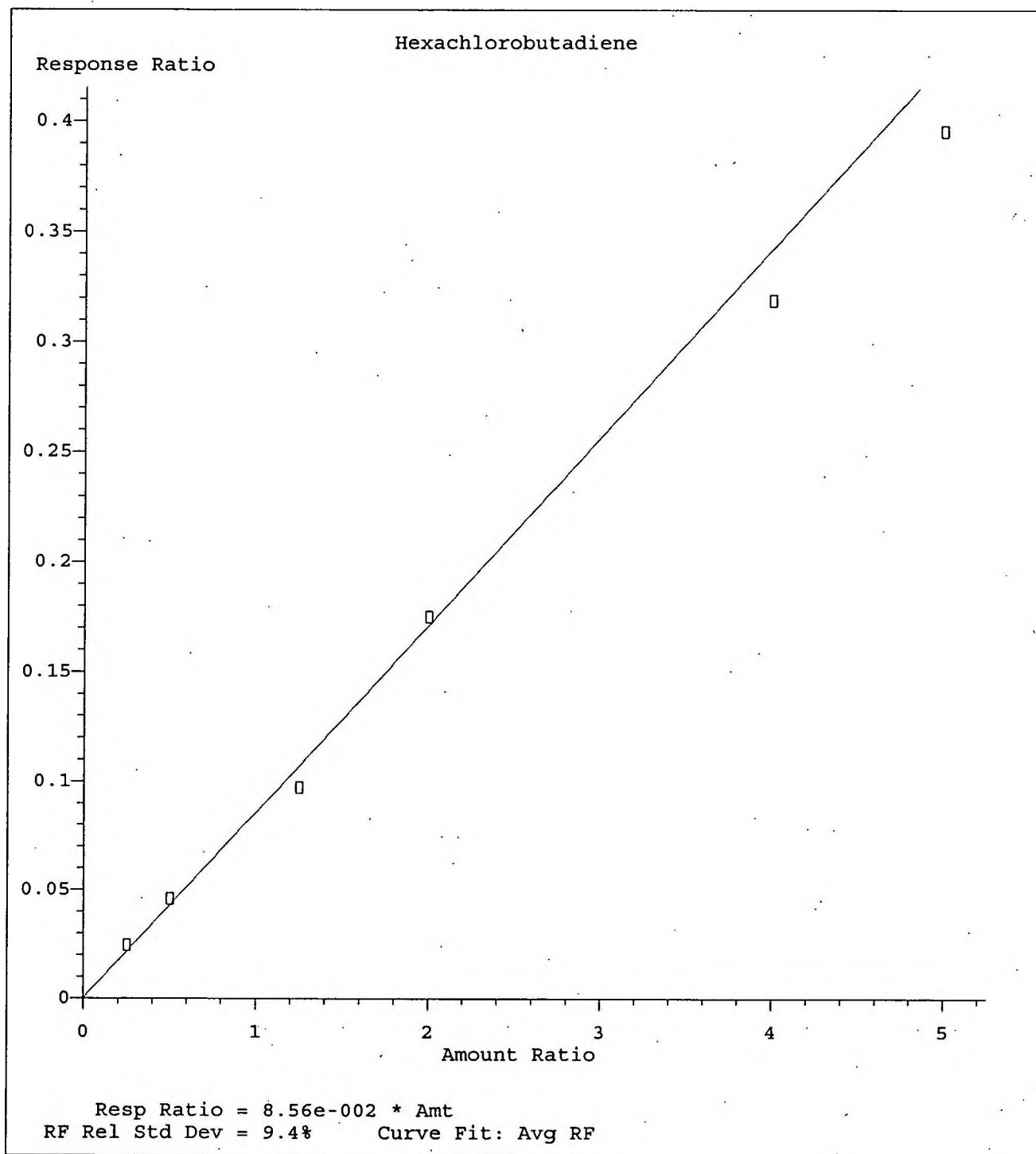
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



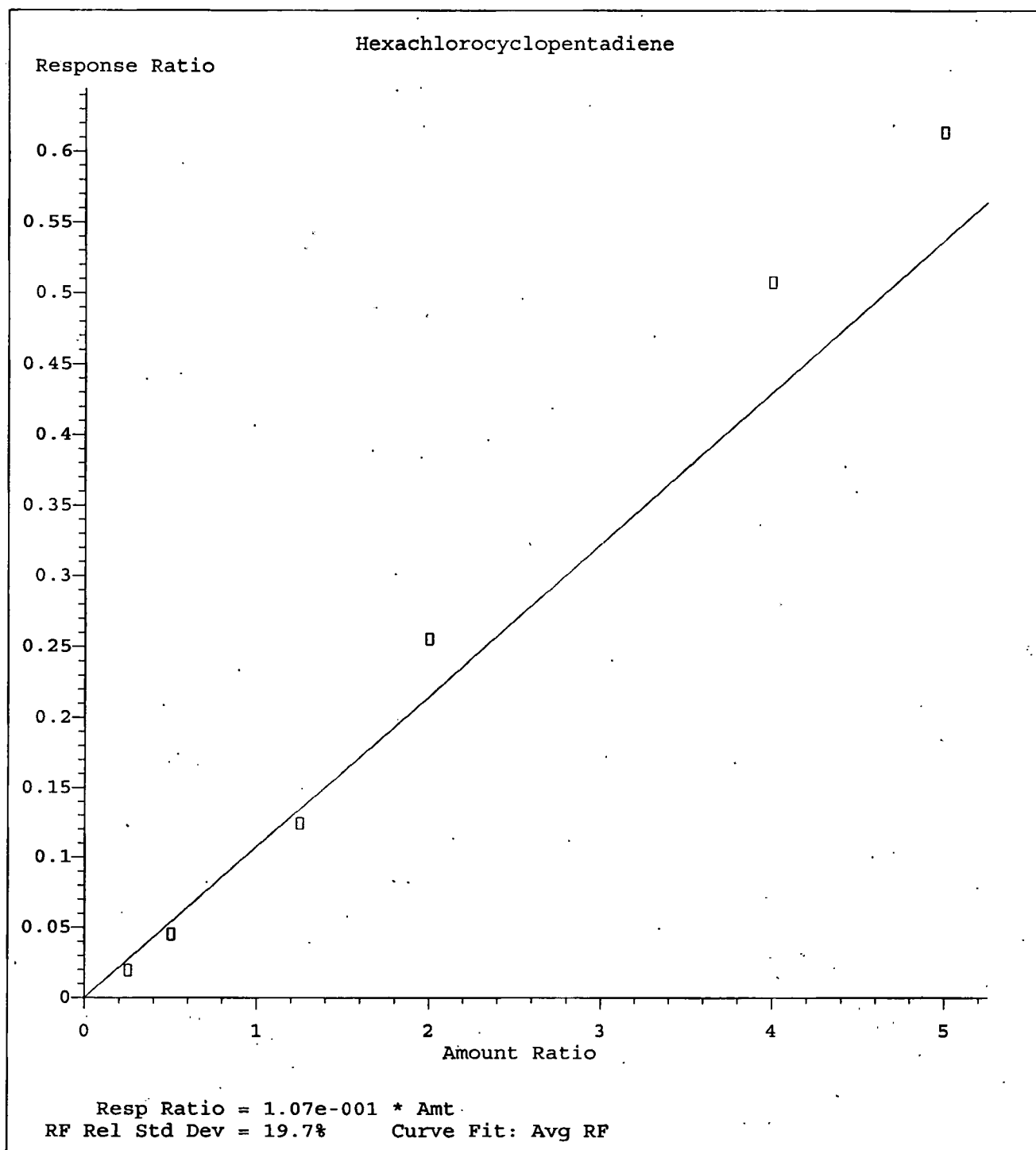
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



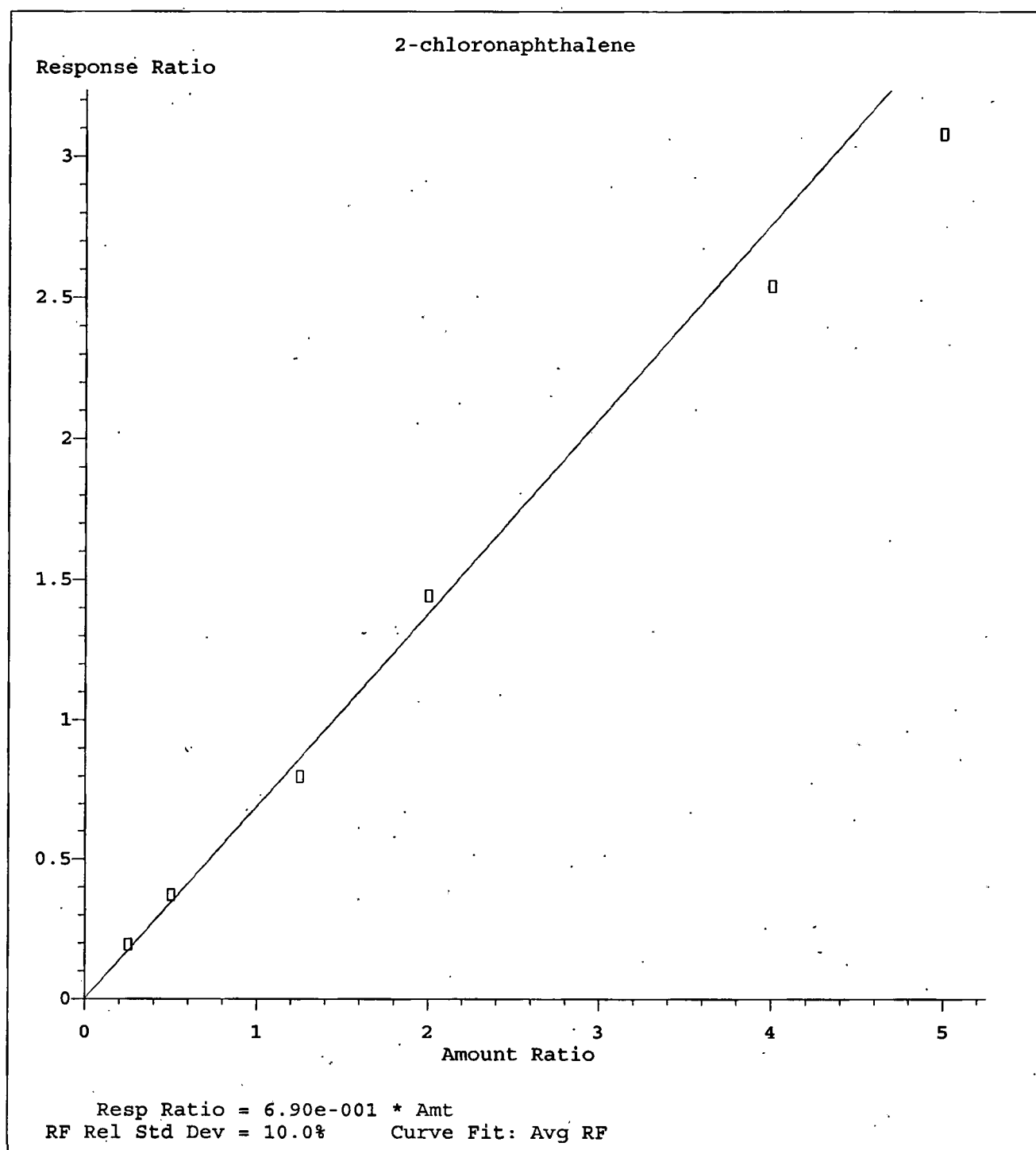
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 Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



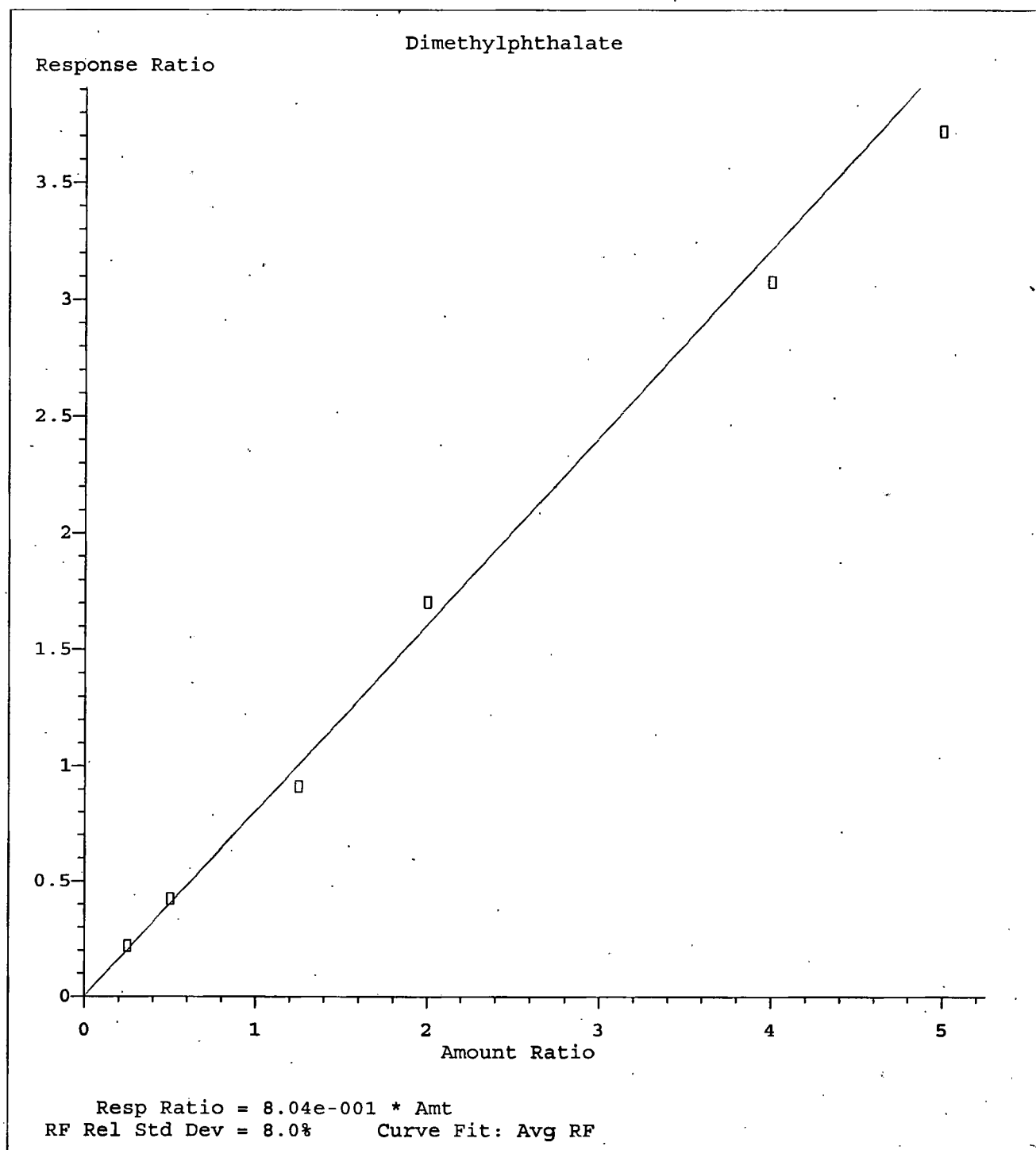
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 Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



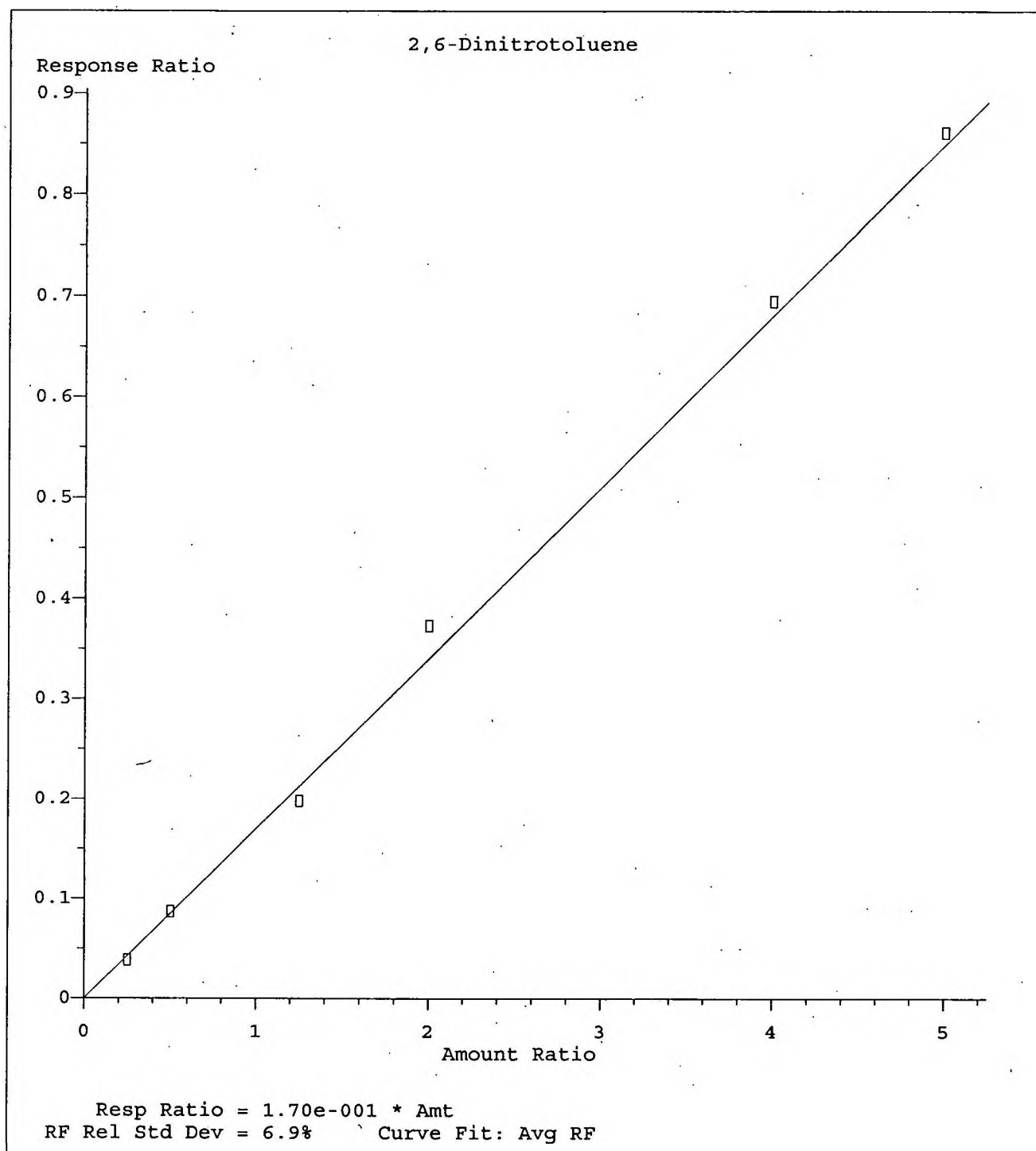
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



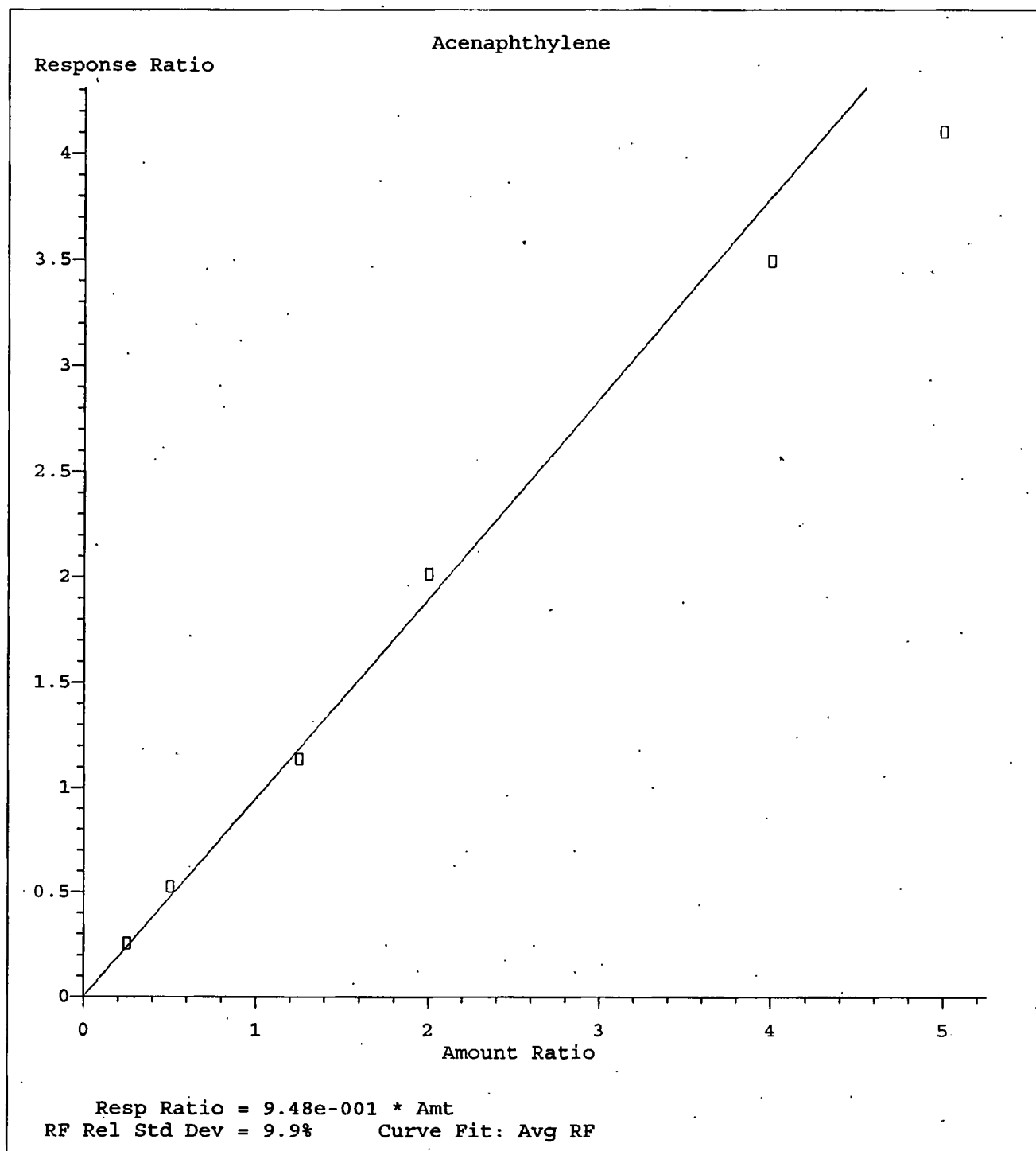
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 Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



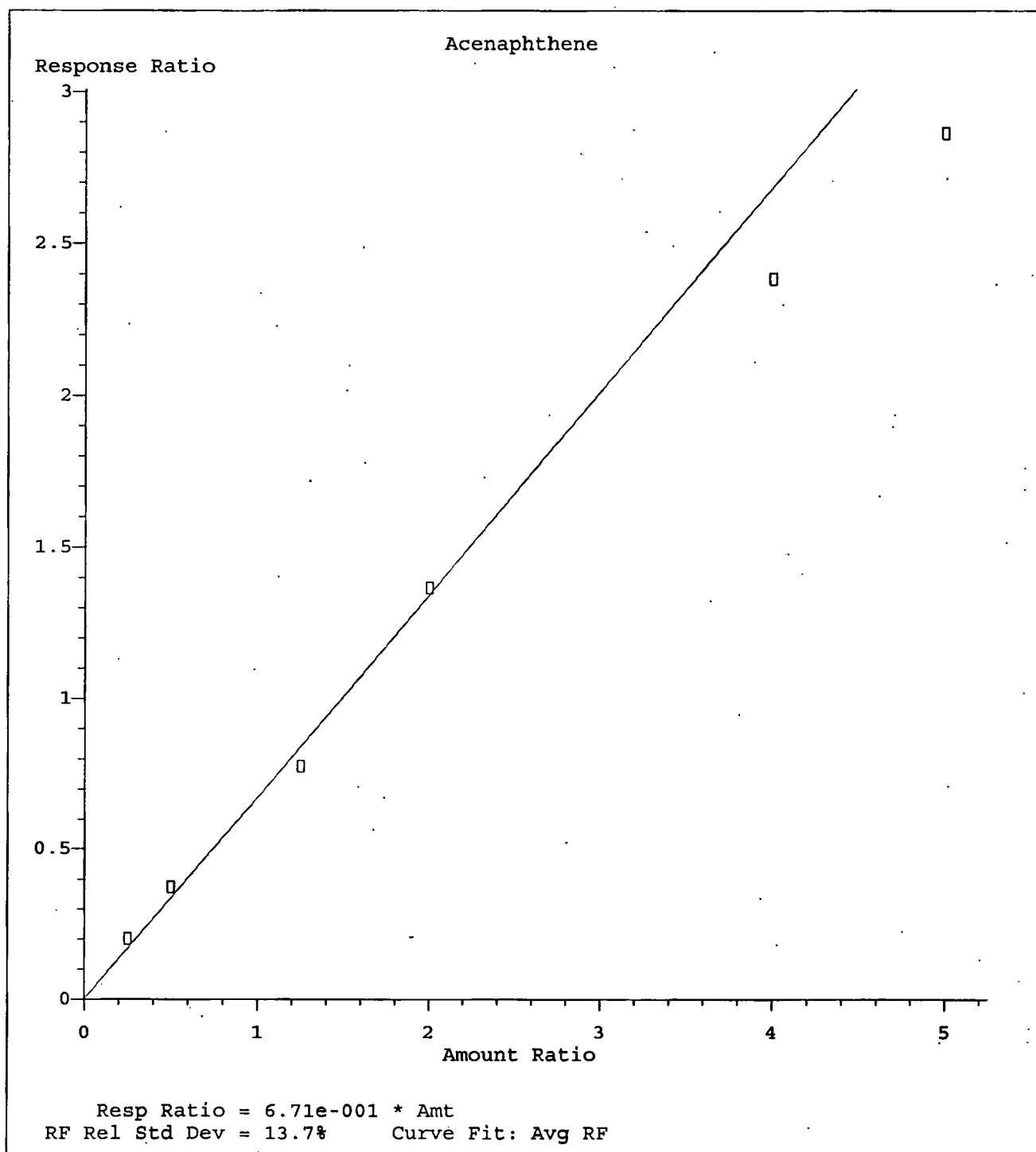
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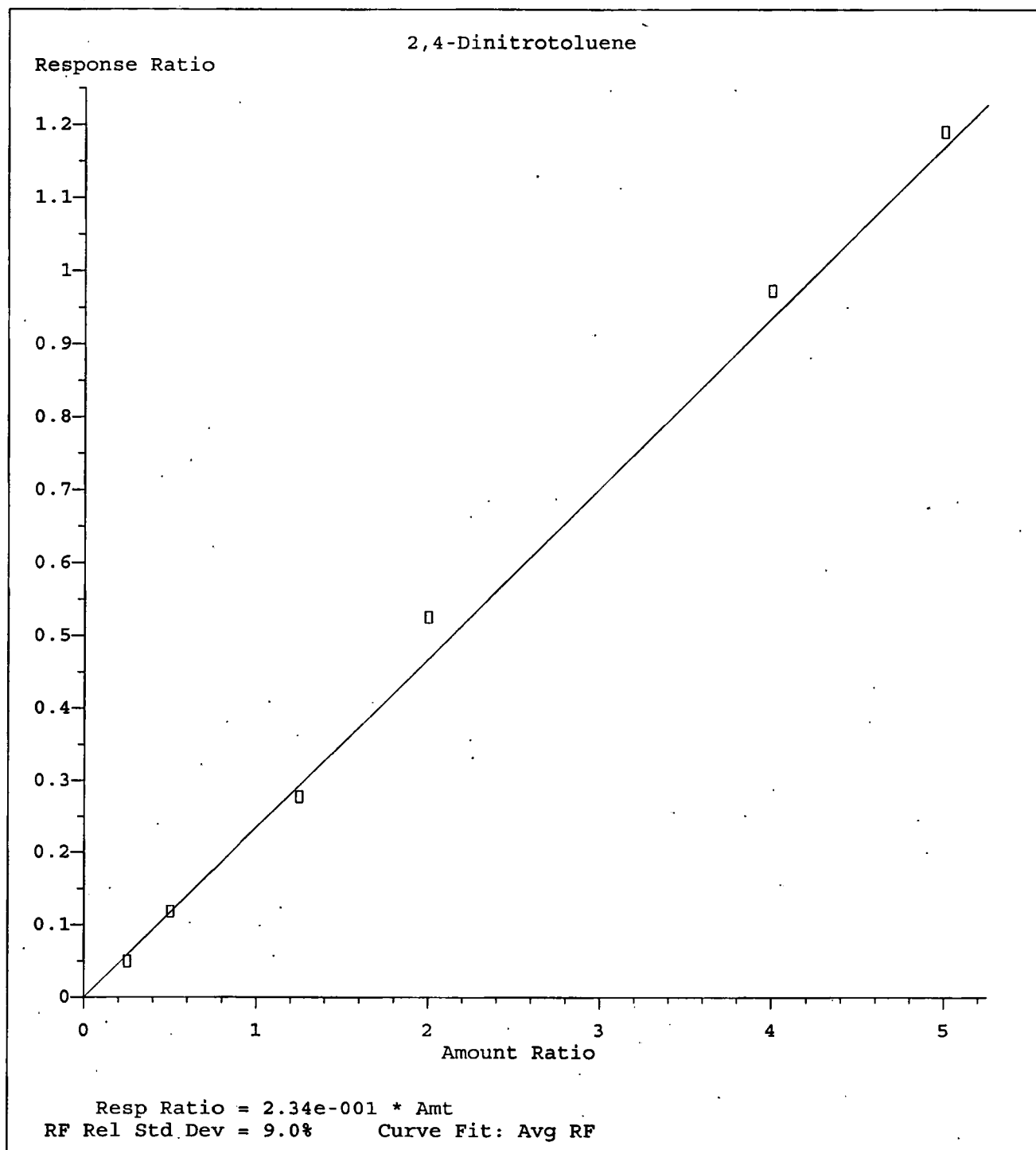
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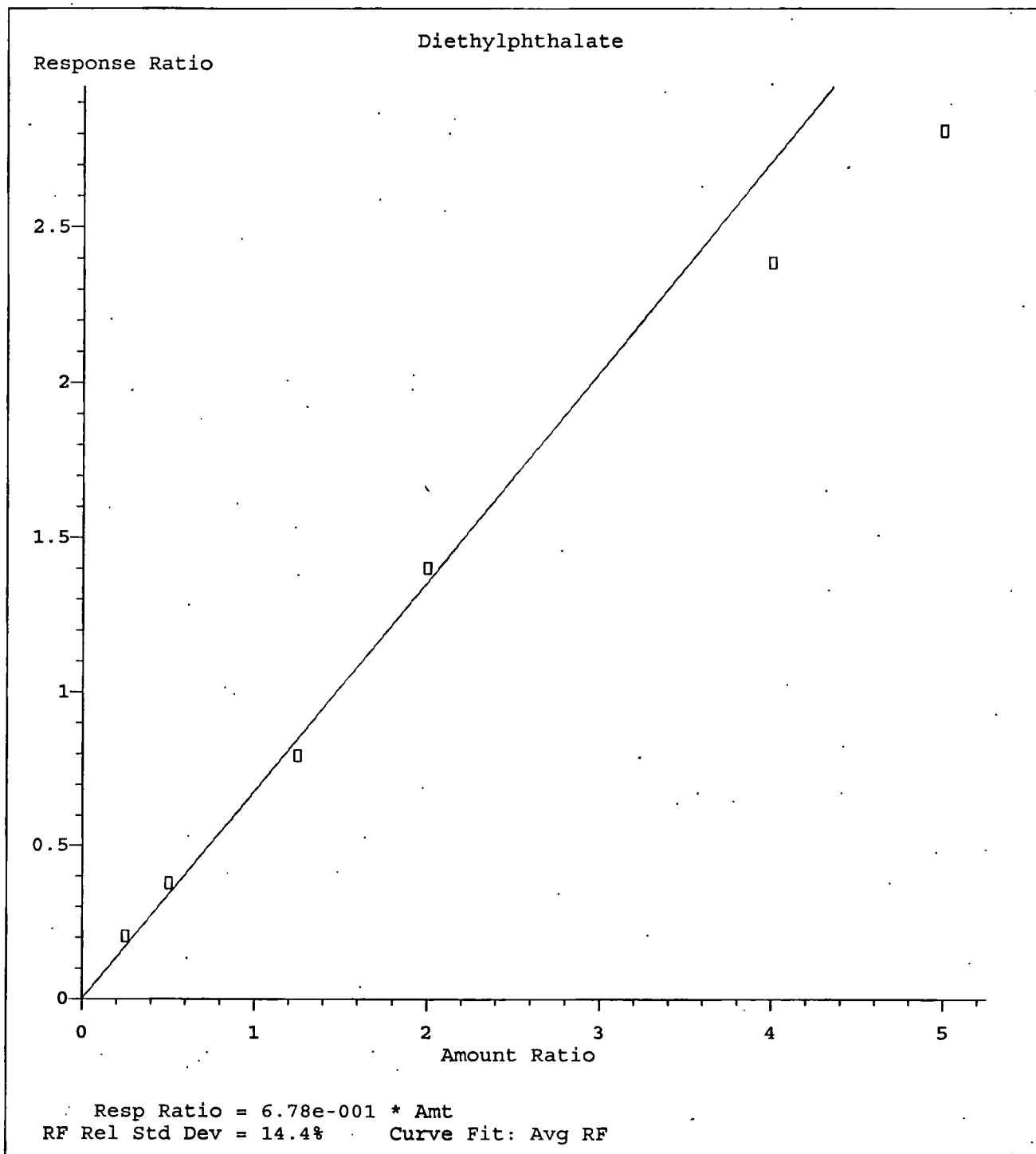
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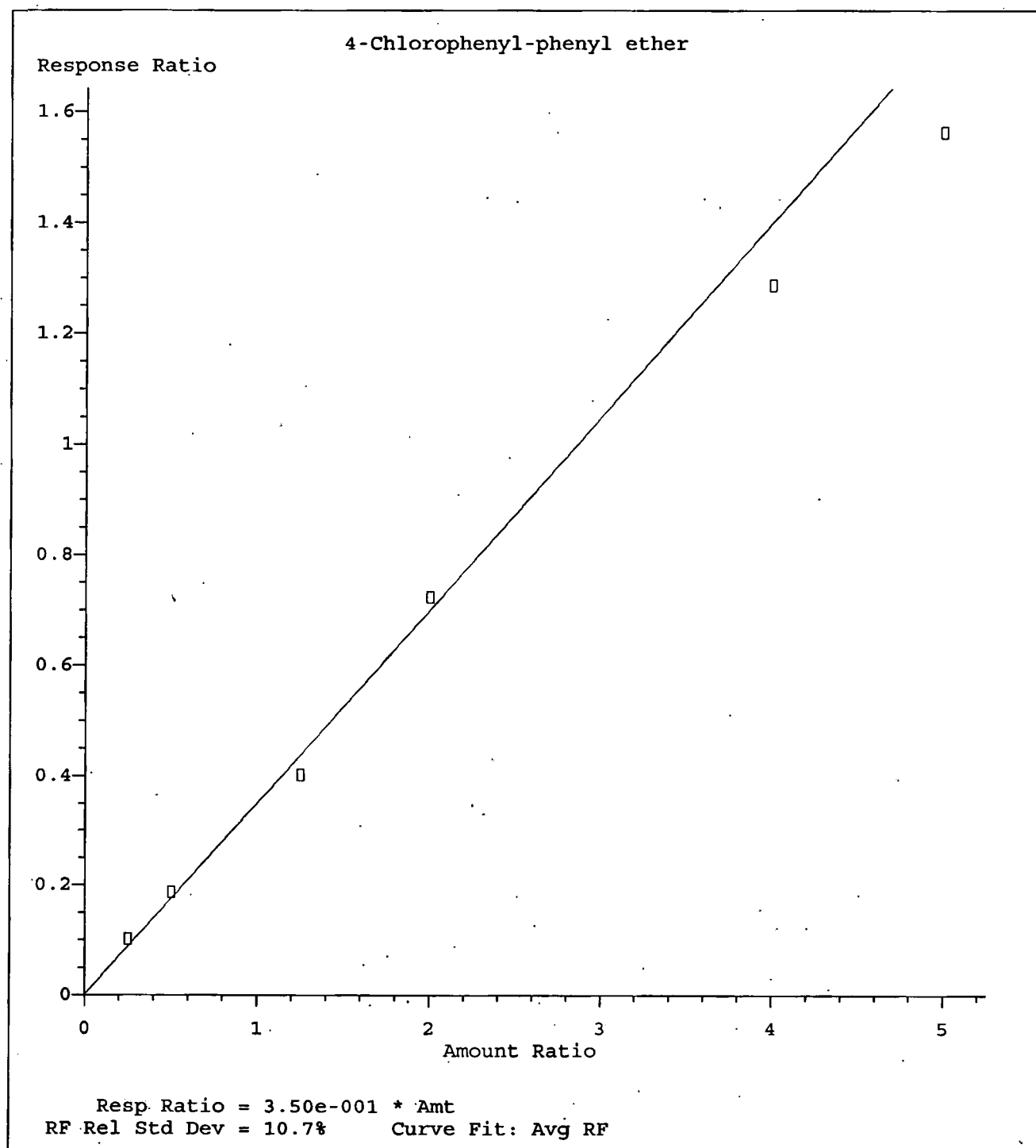
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 Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



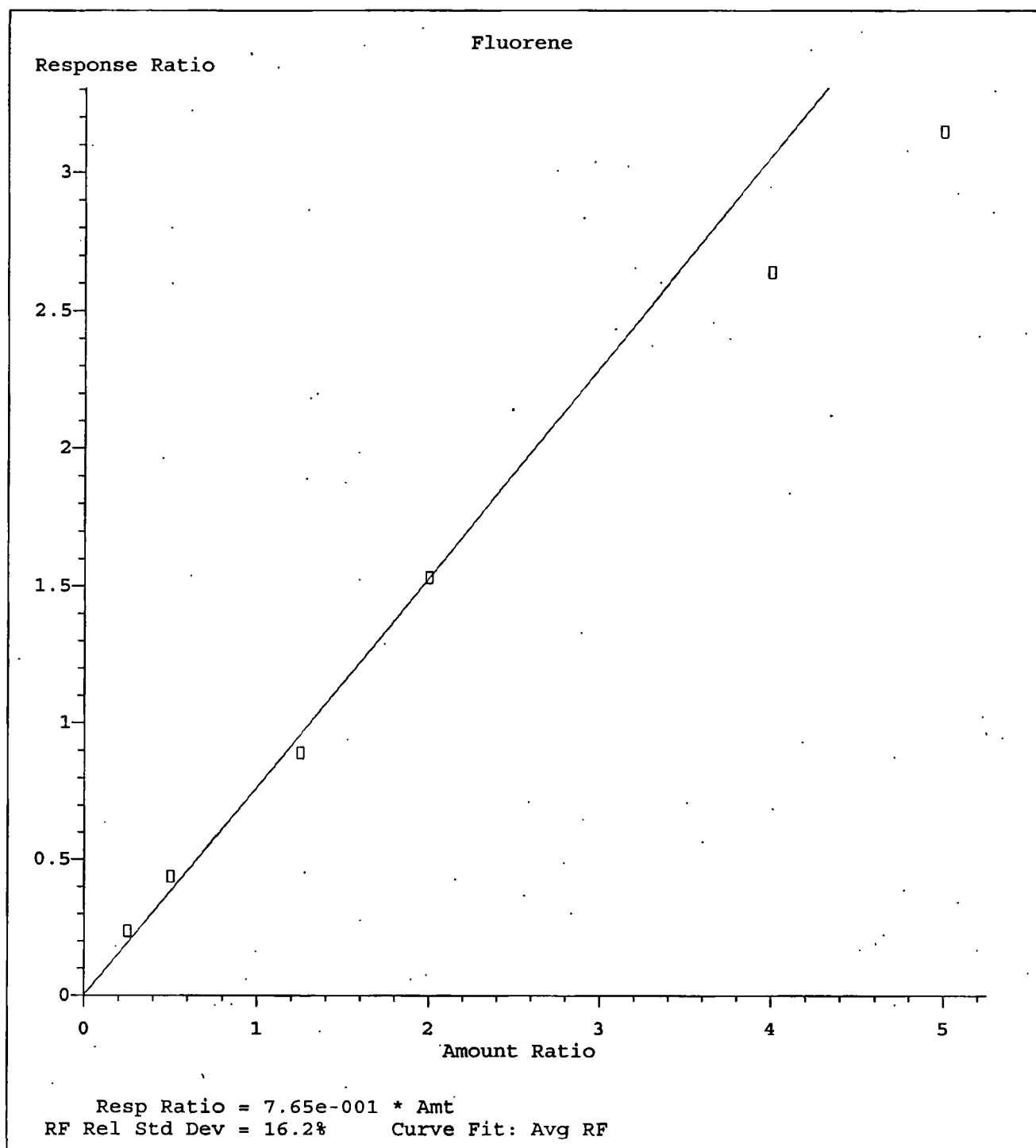
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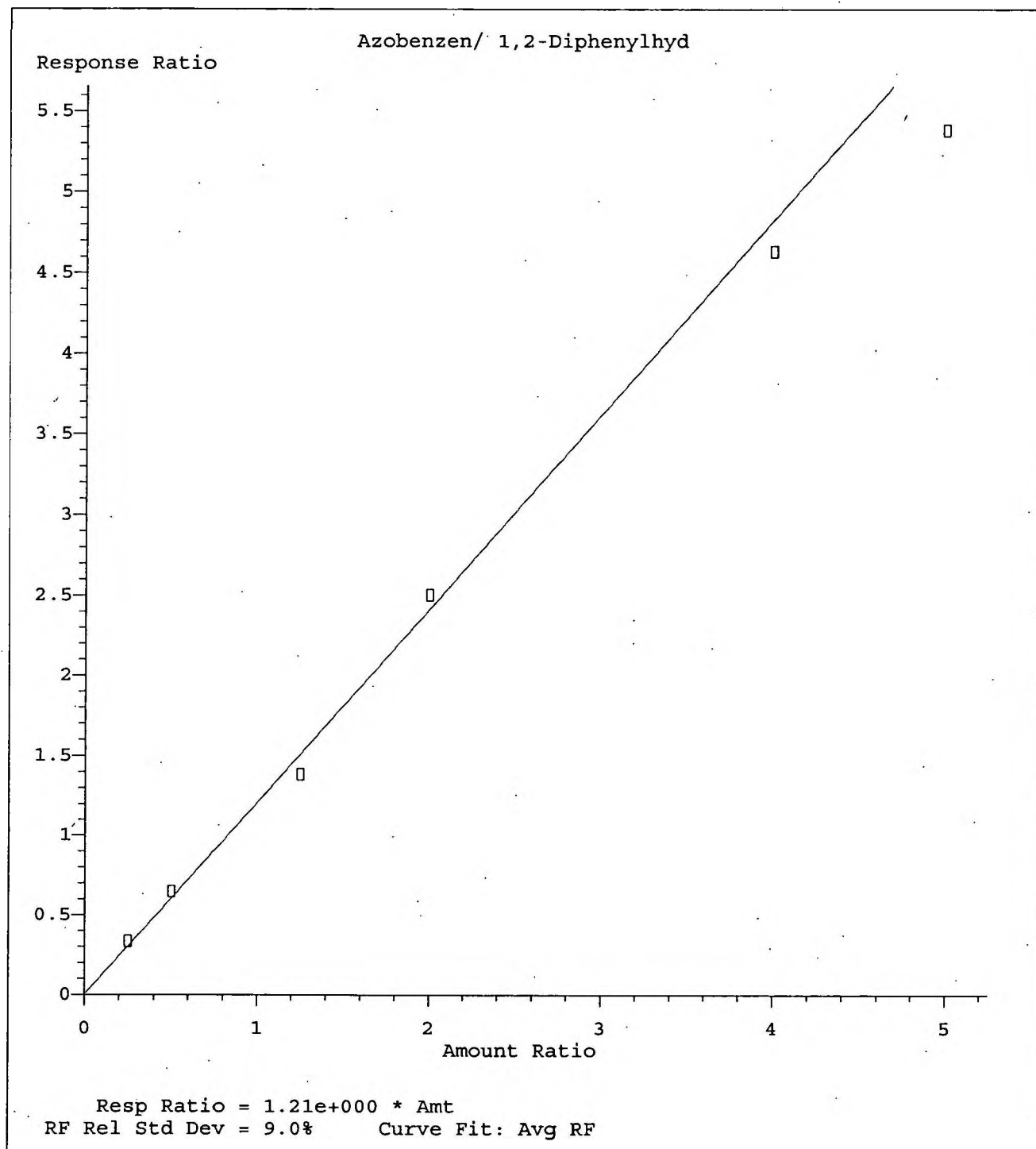
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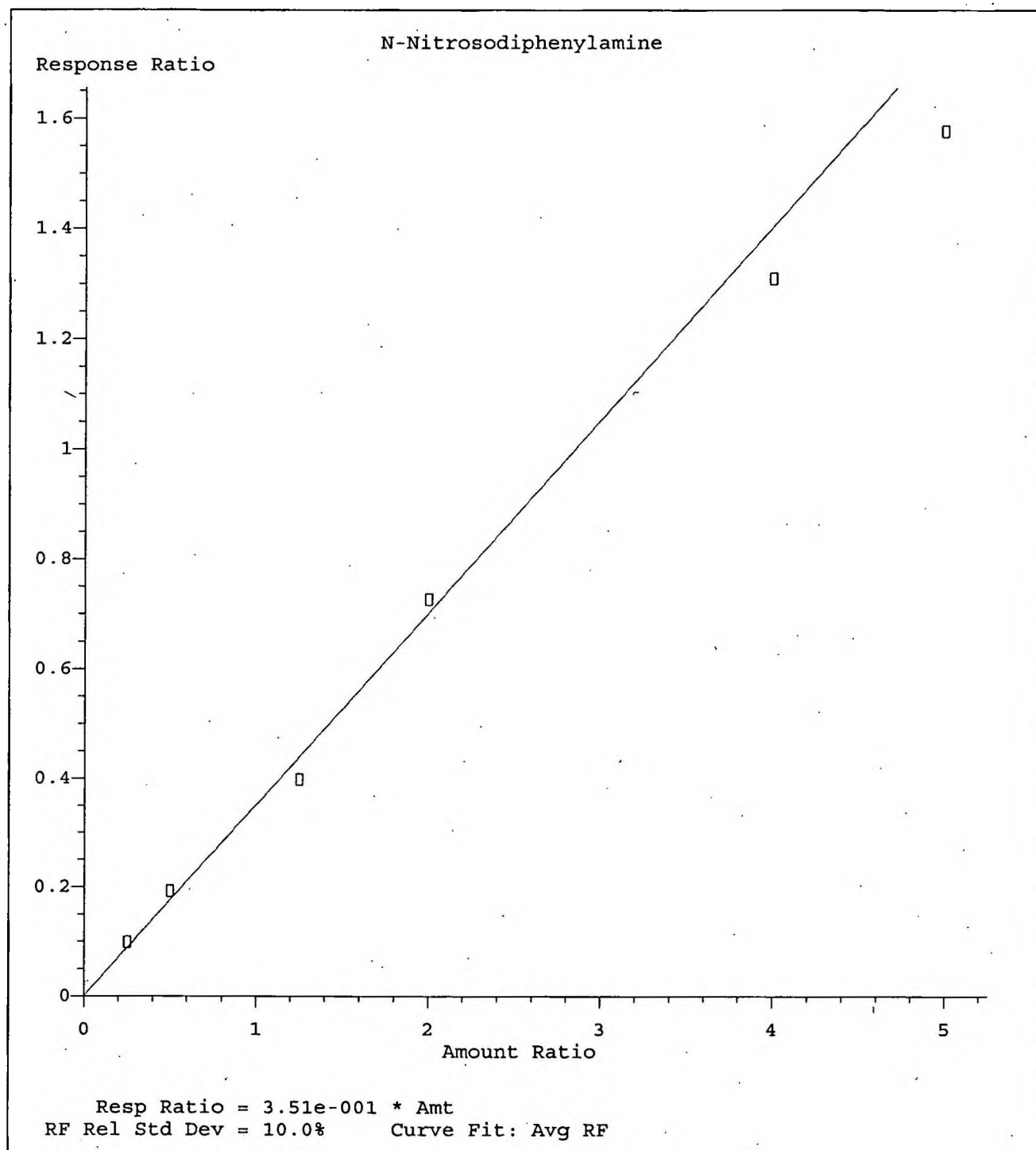
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



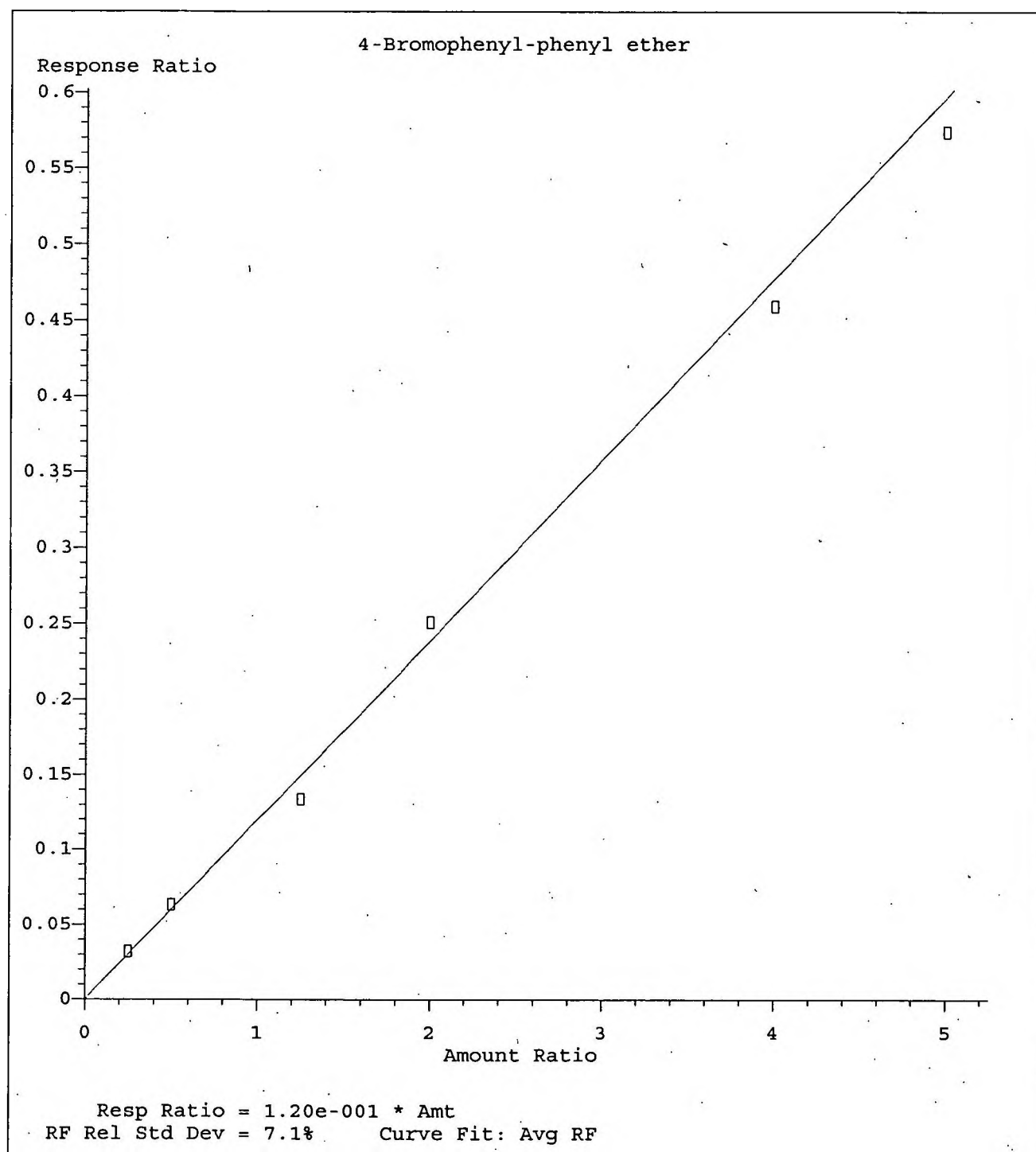
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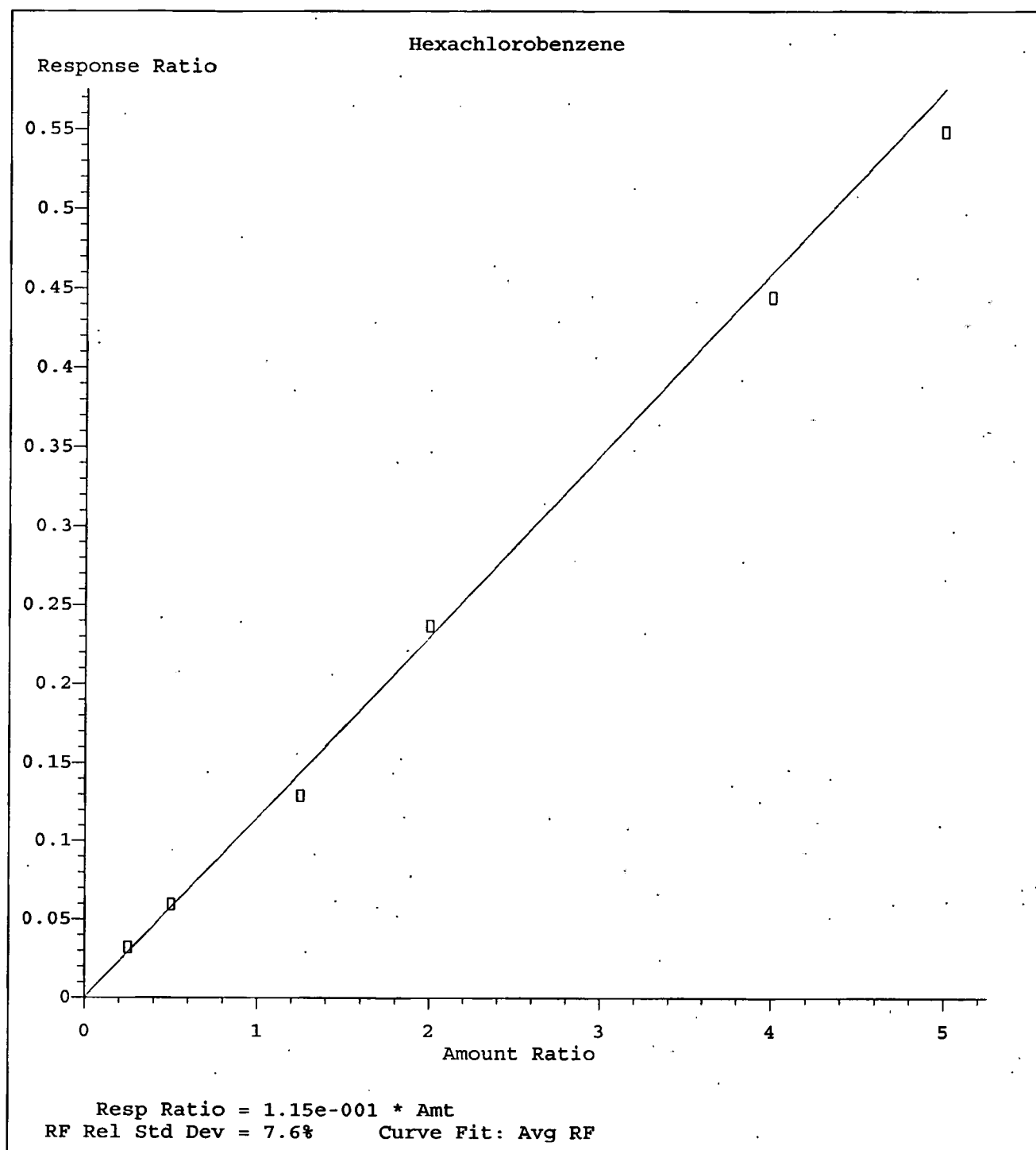
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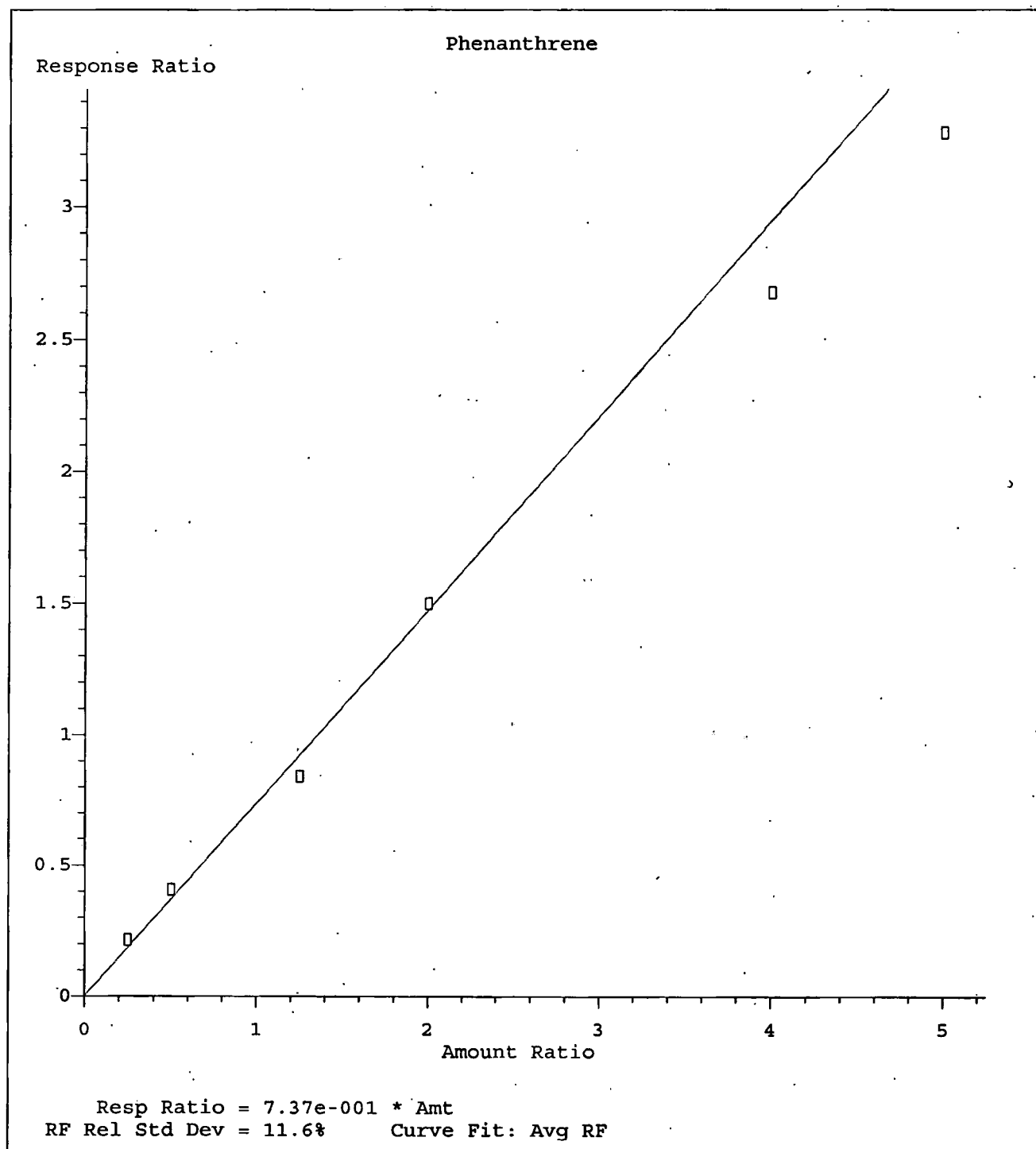
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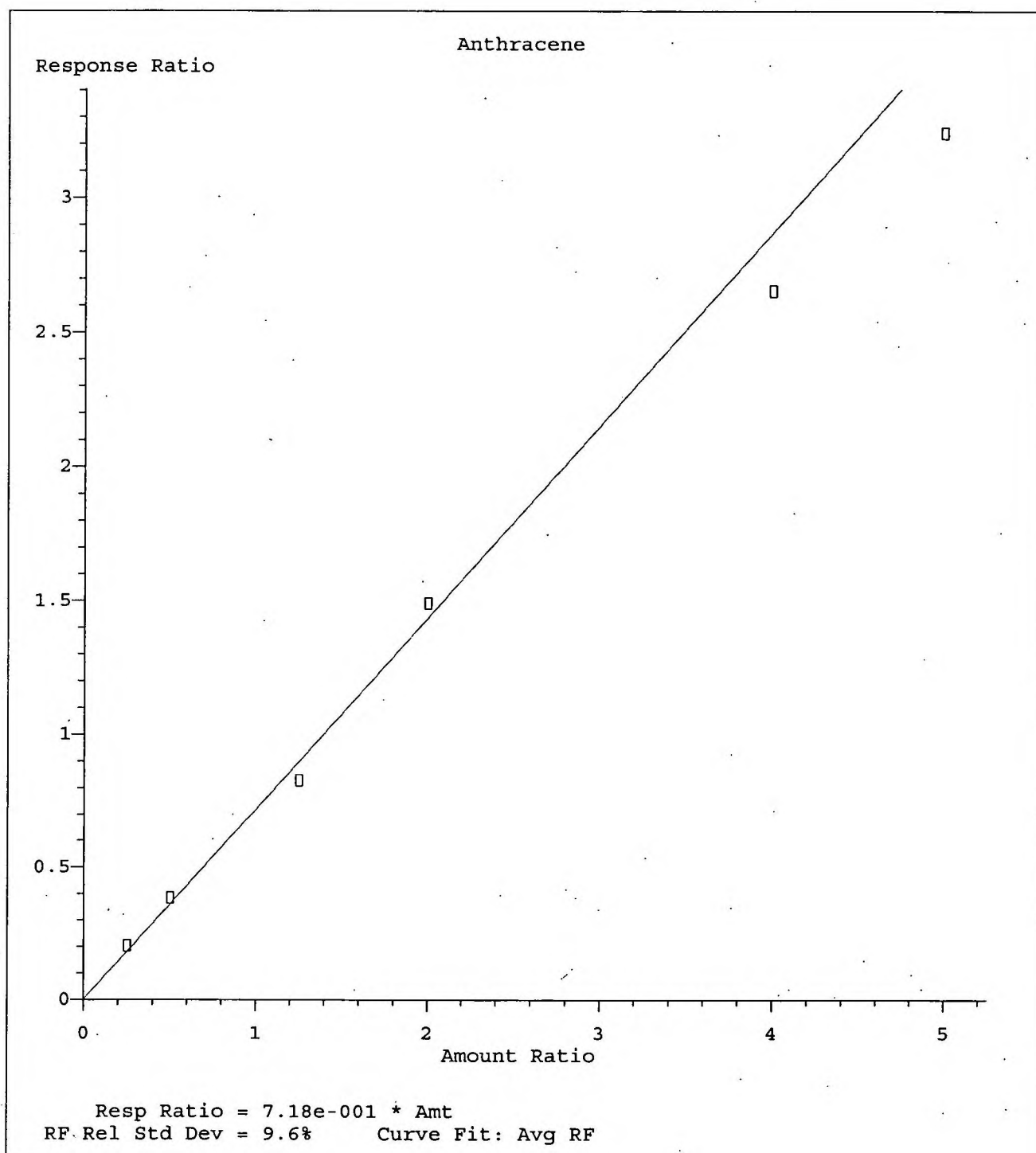
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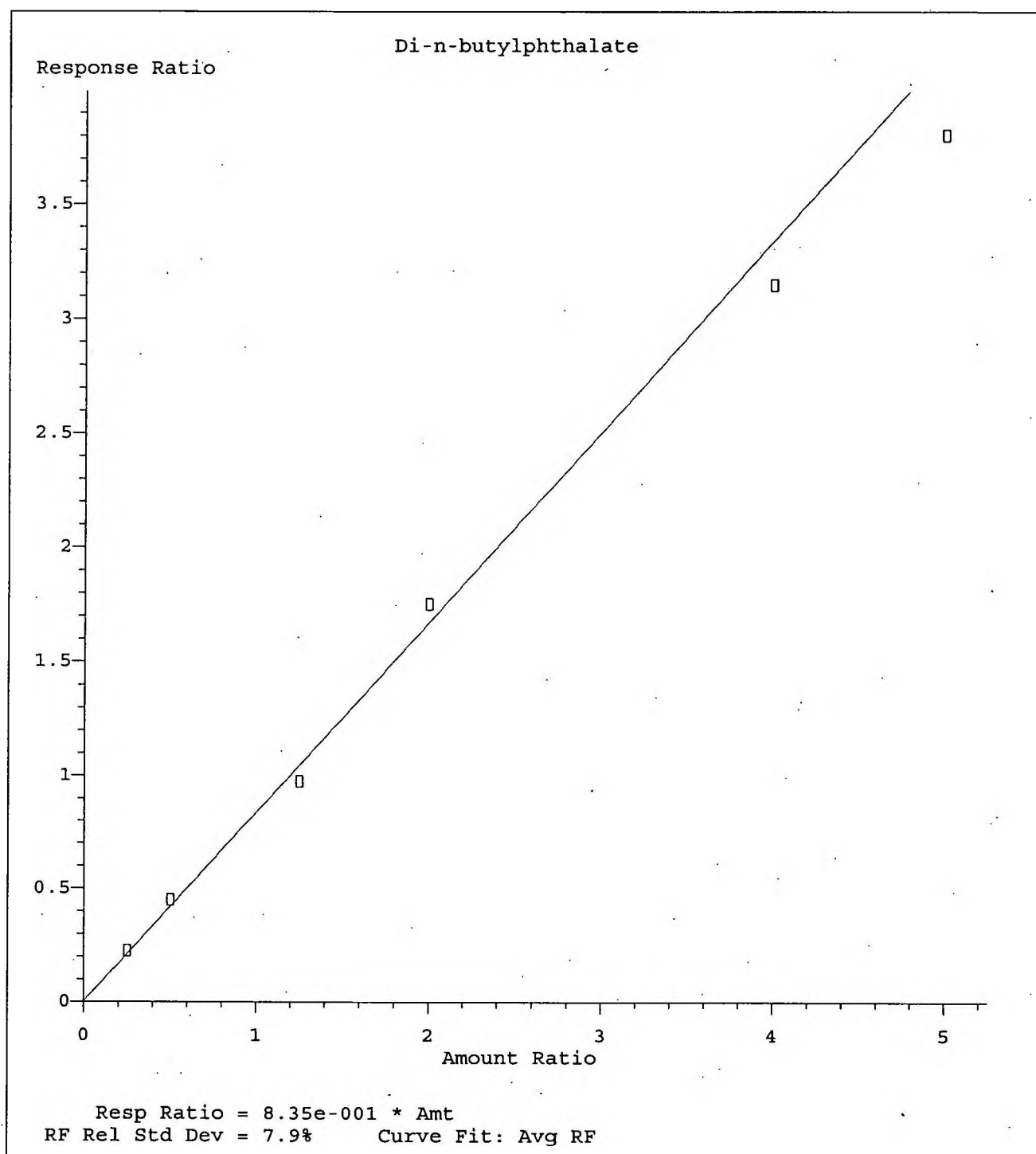
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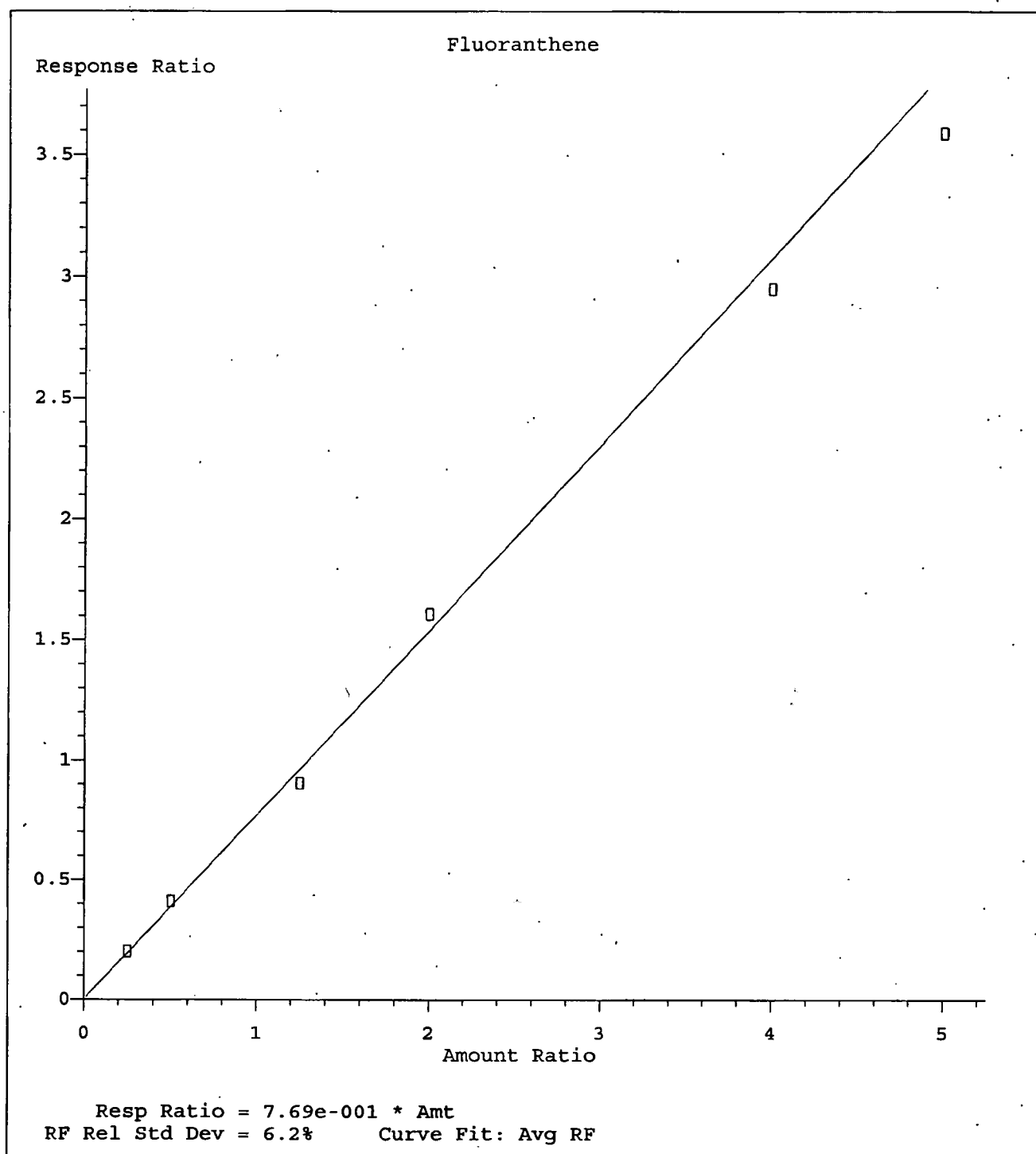
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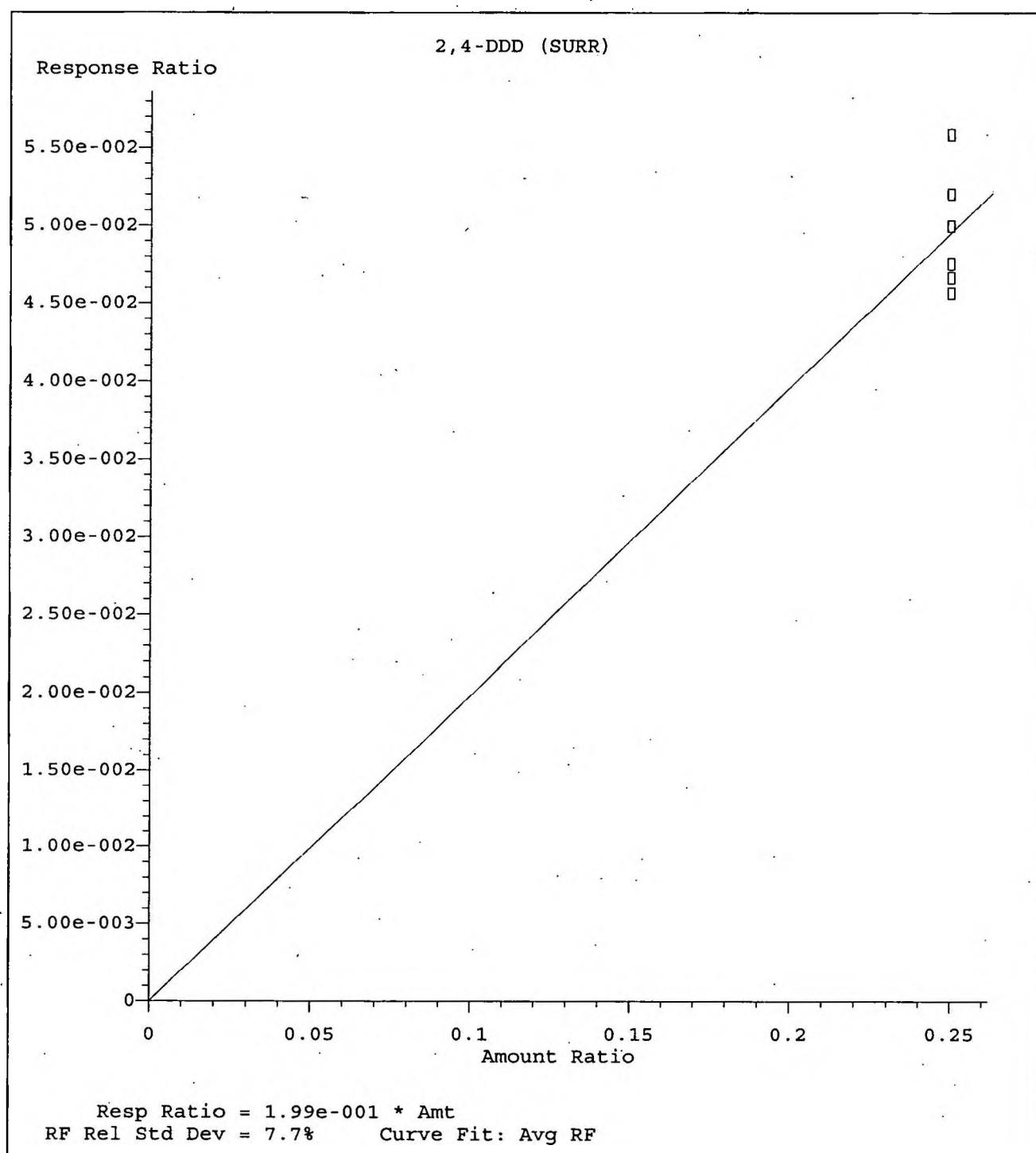
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 Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



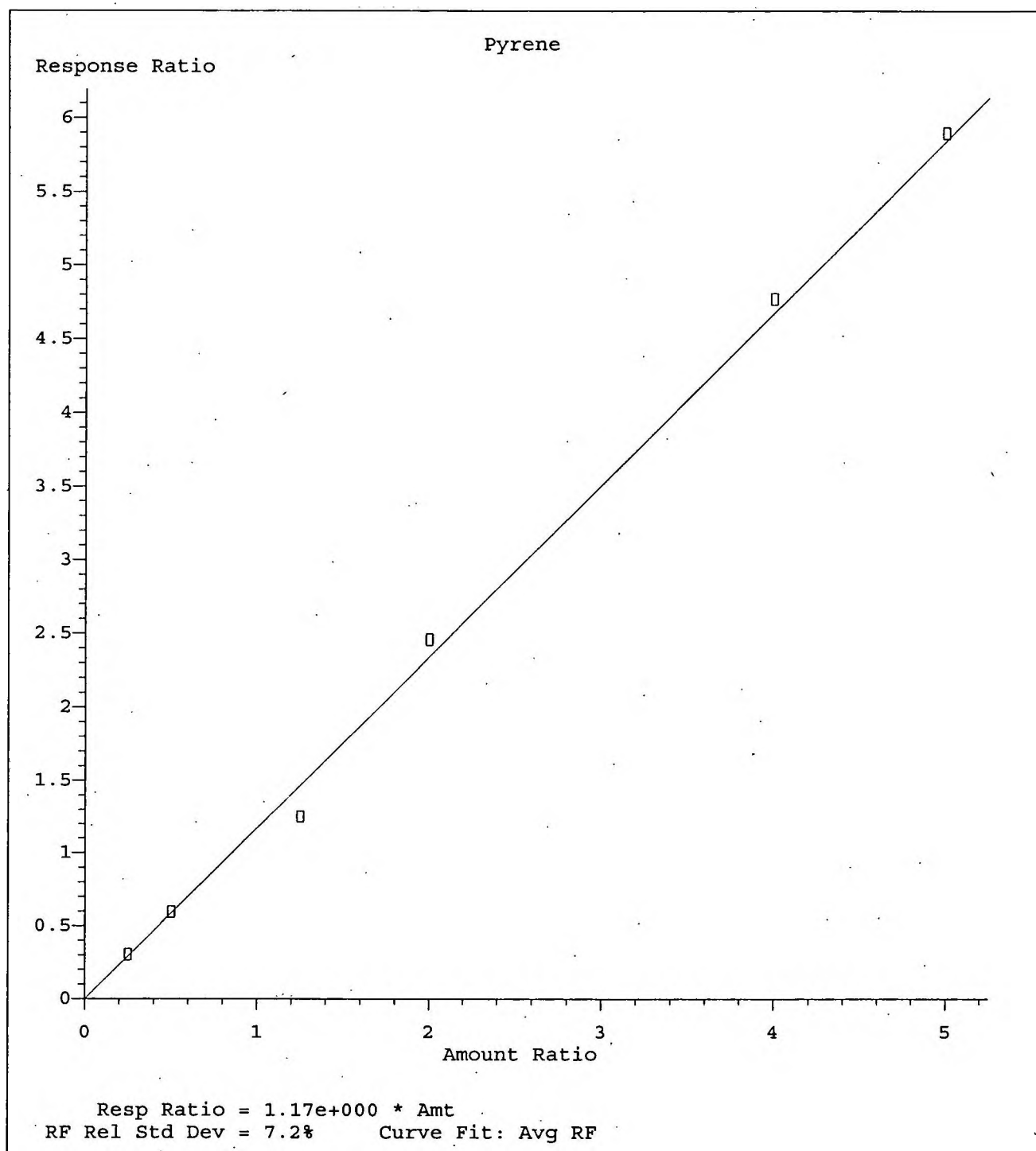
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



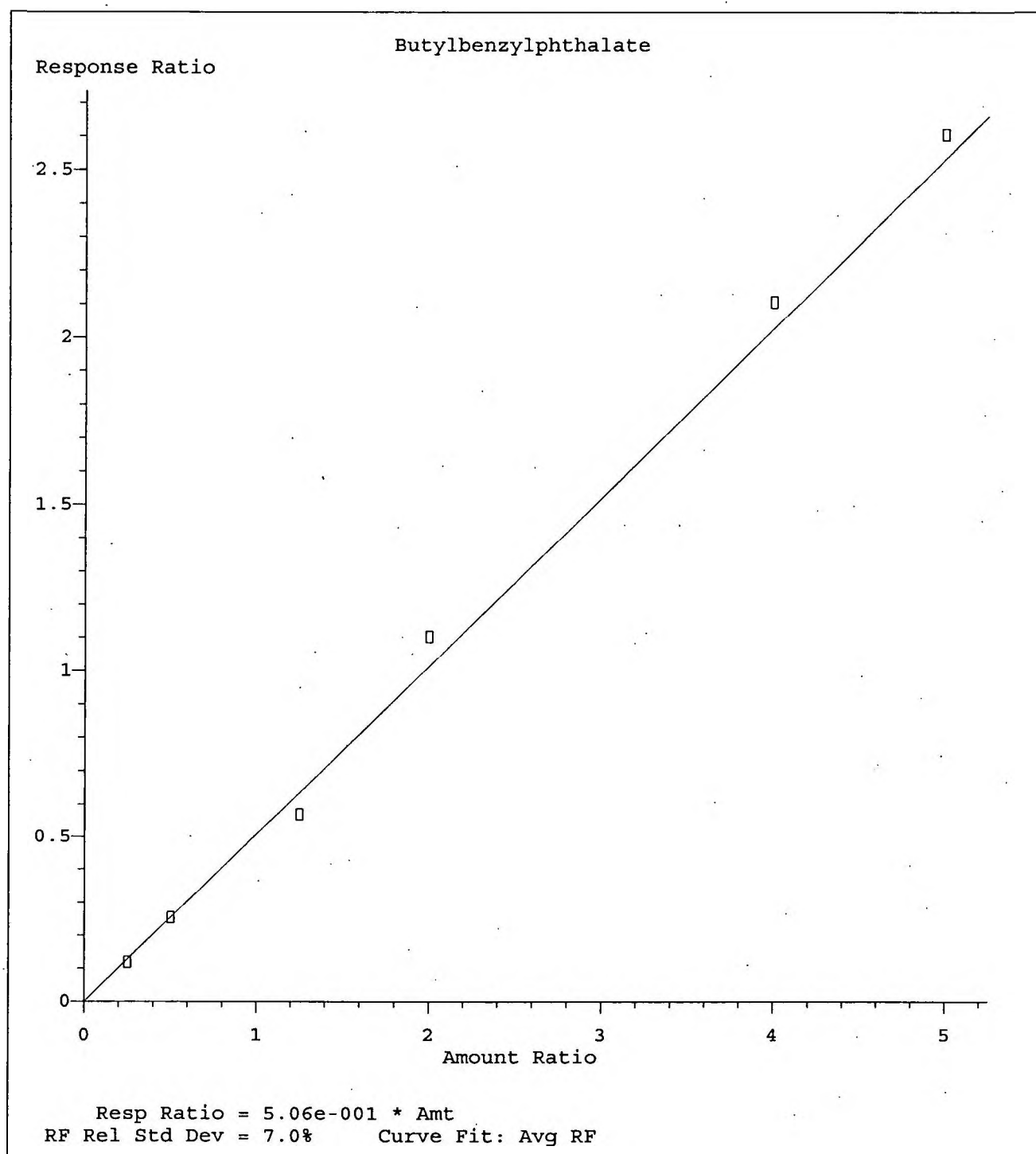
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



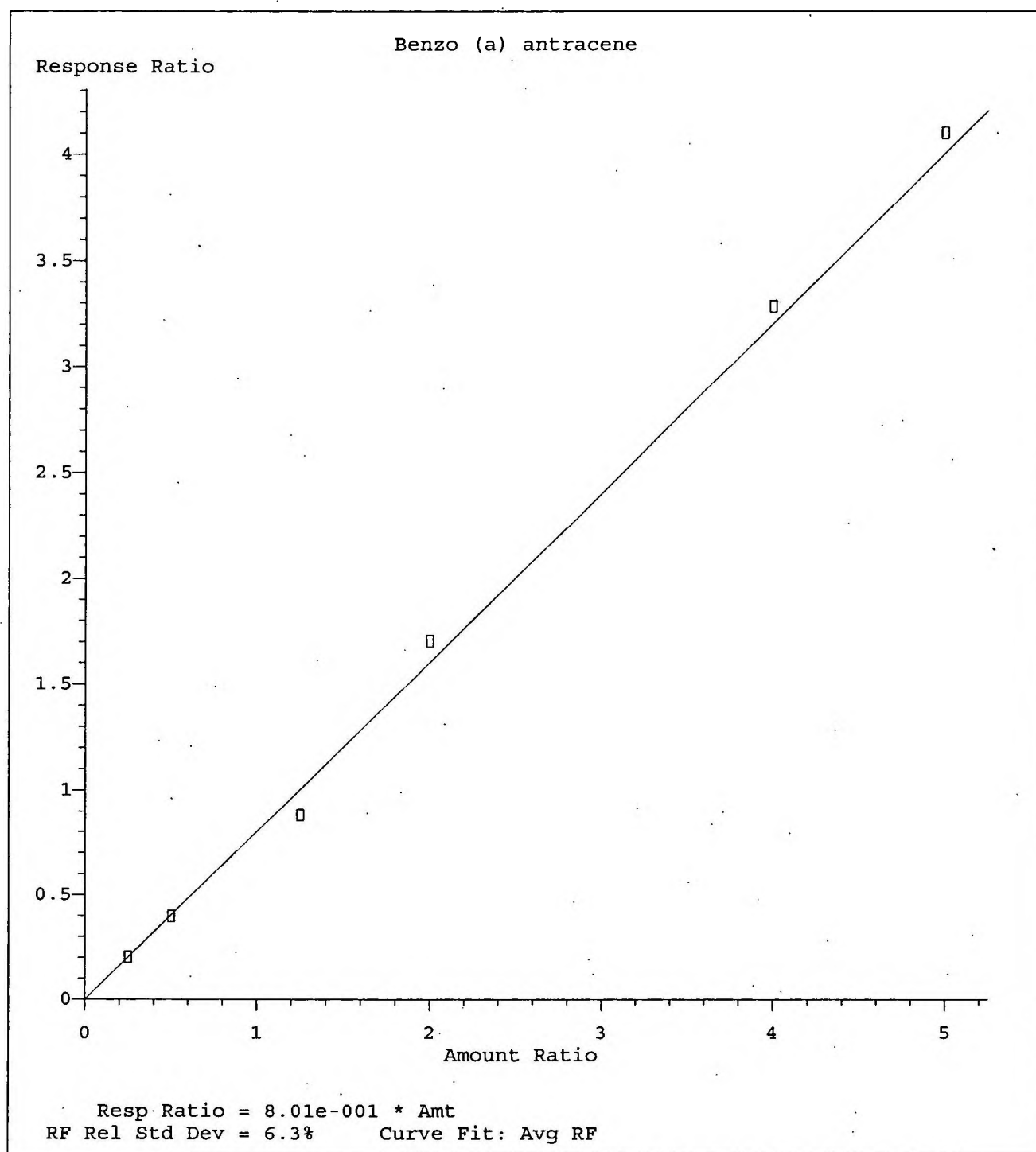
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



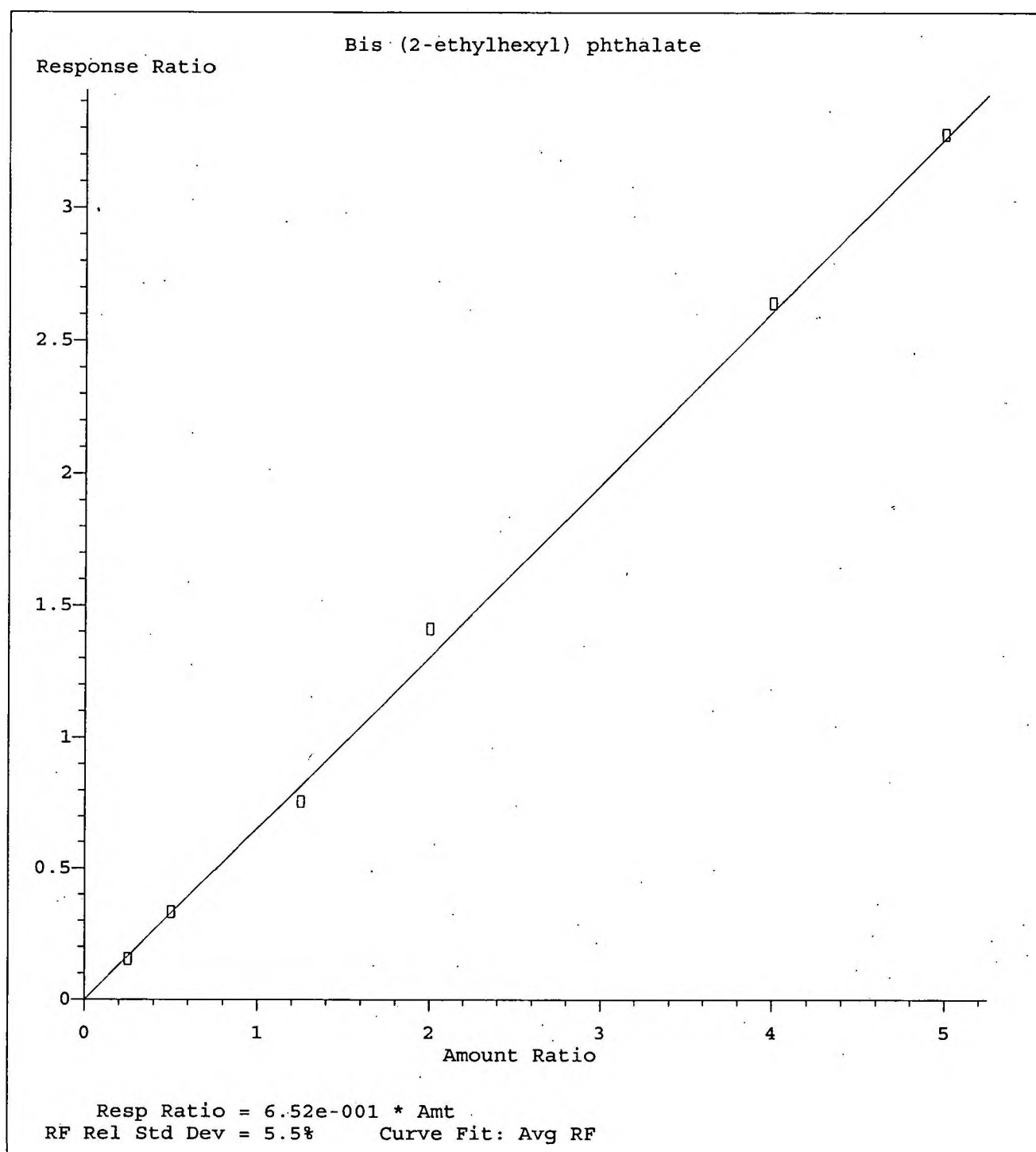
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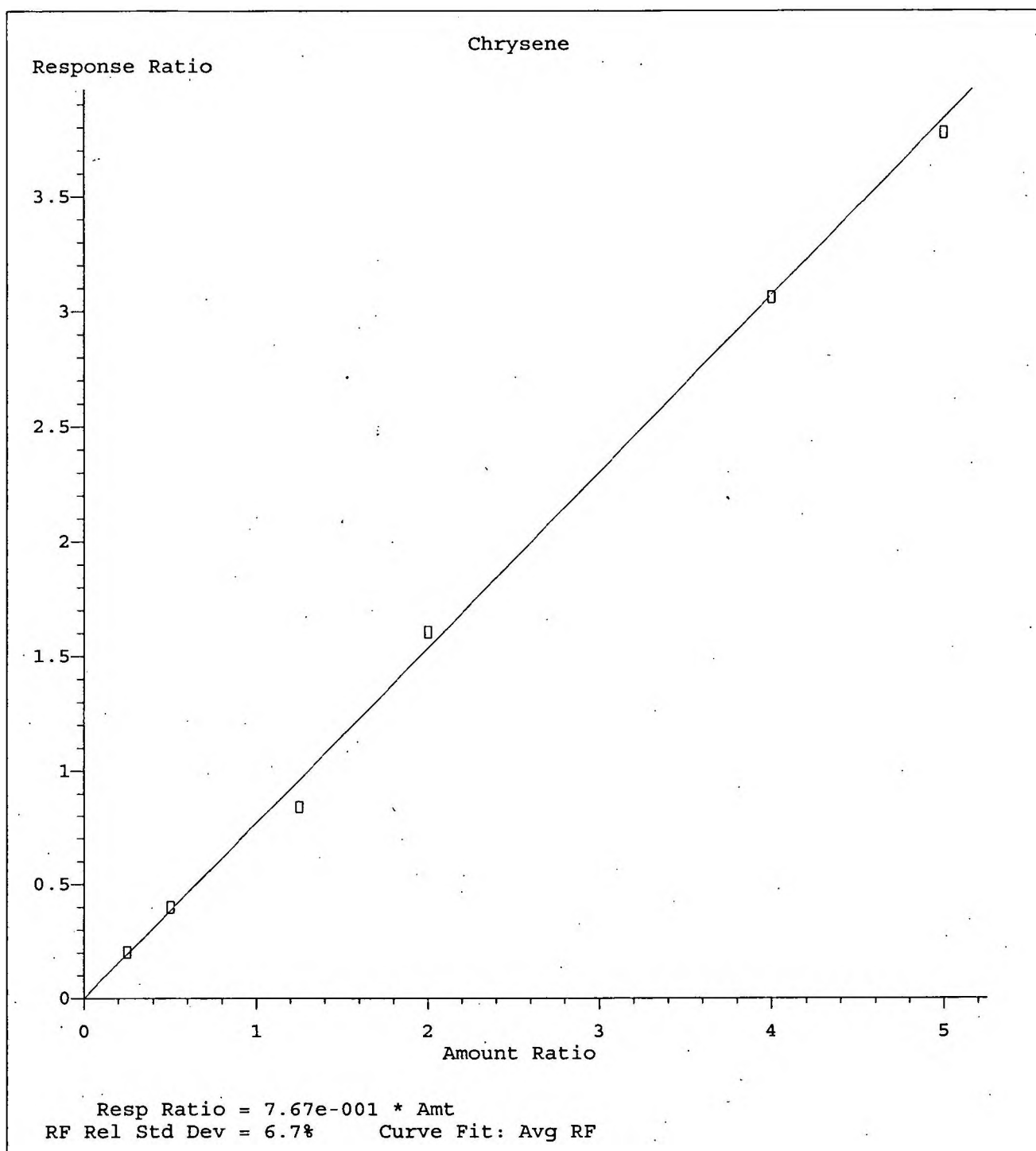
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



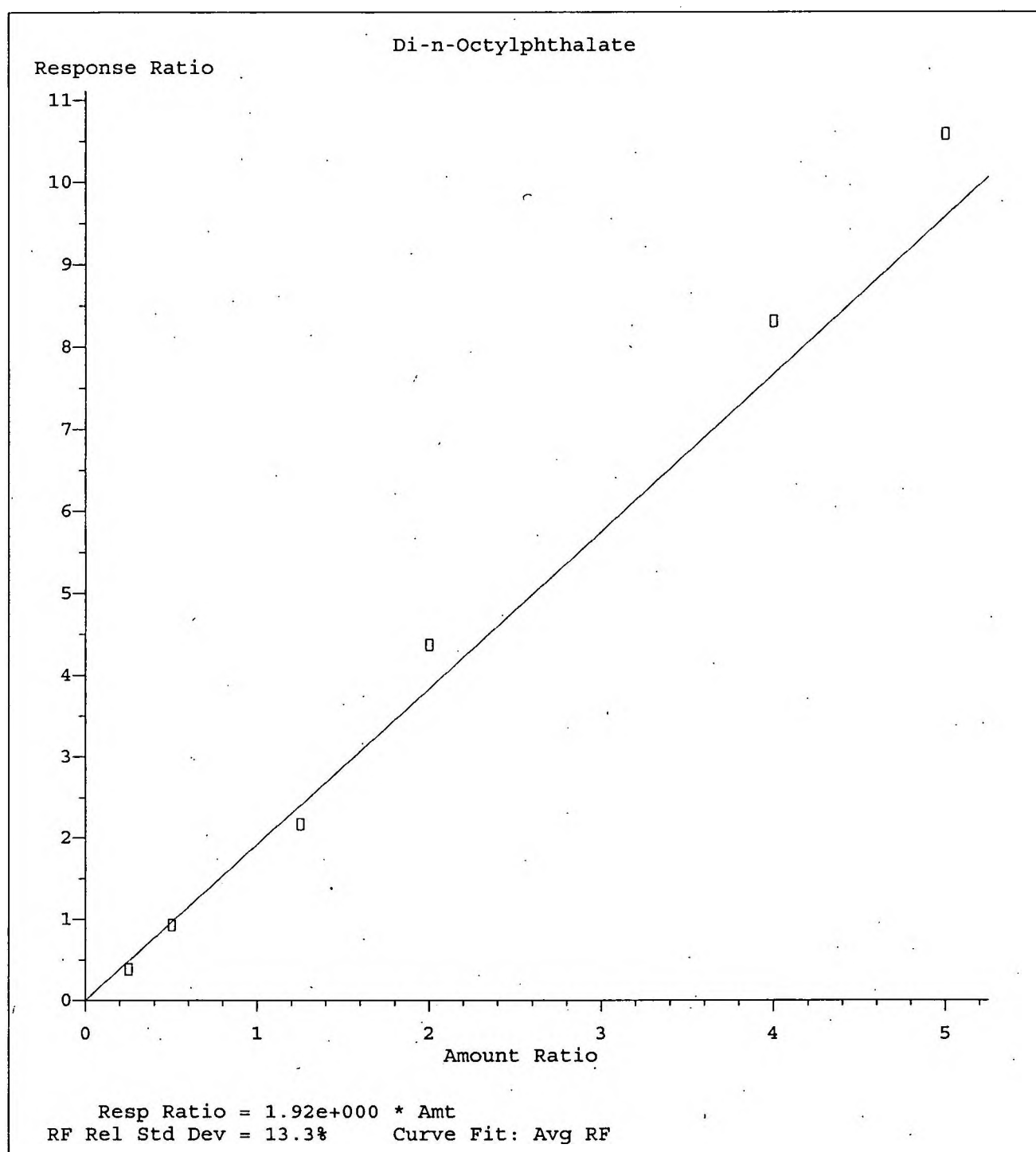
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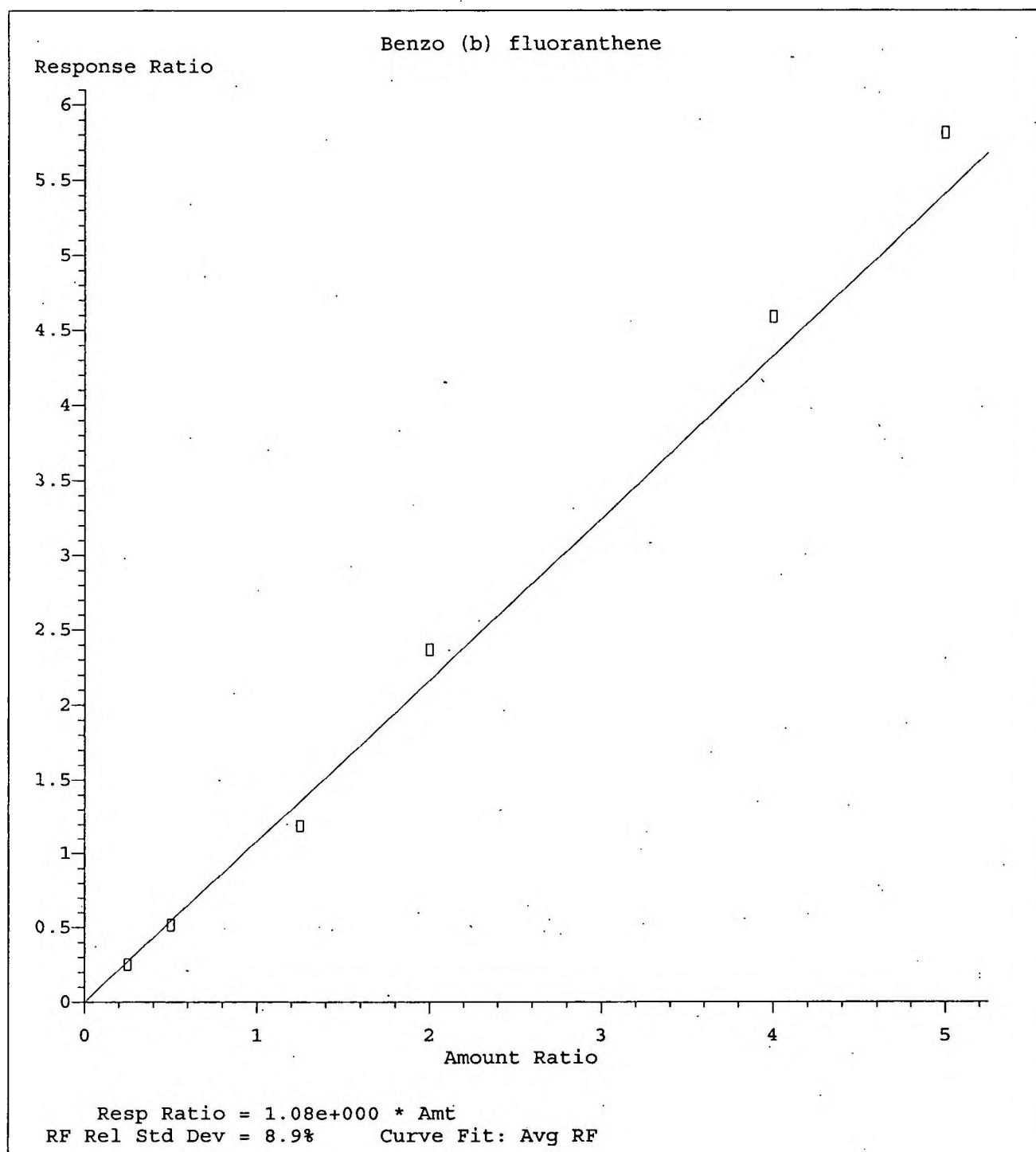
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Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



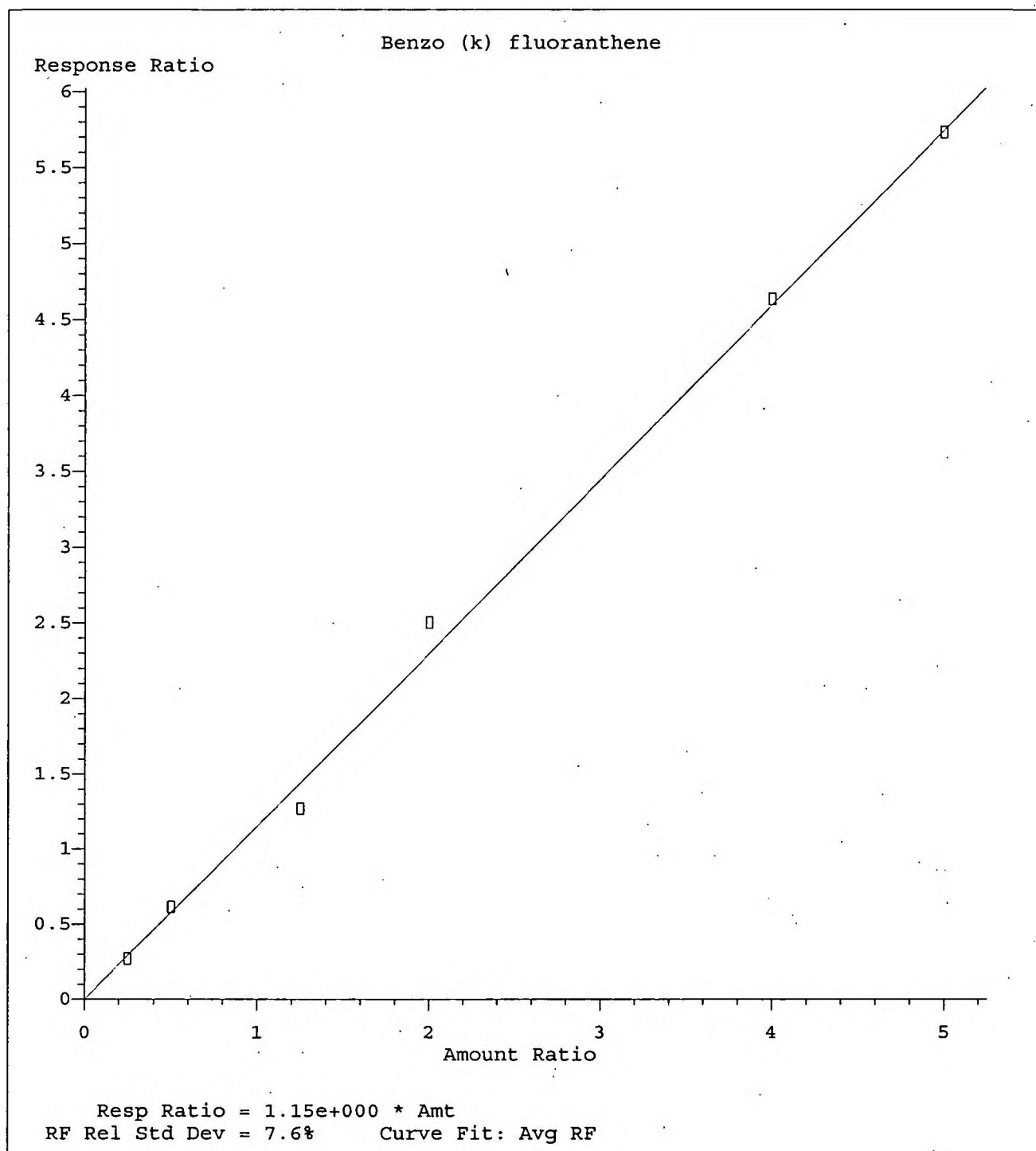
Method Name: C:\MSDCHEM\1\5973N\5080402.M
 Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



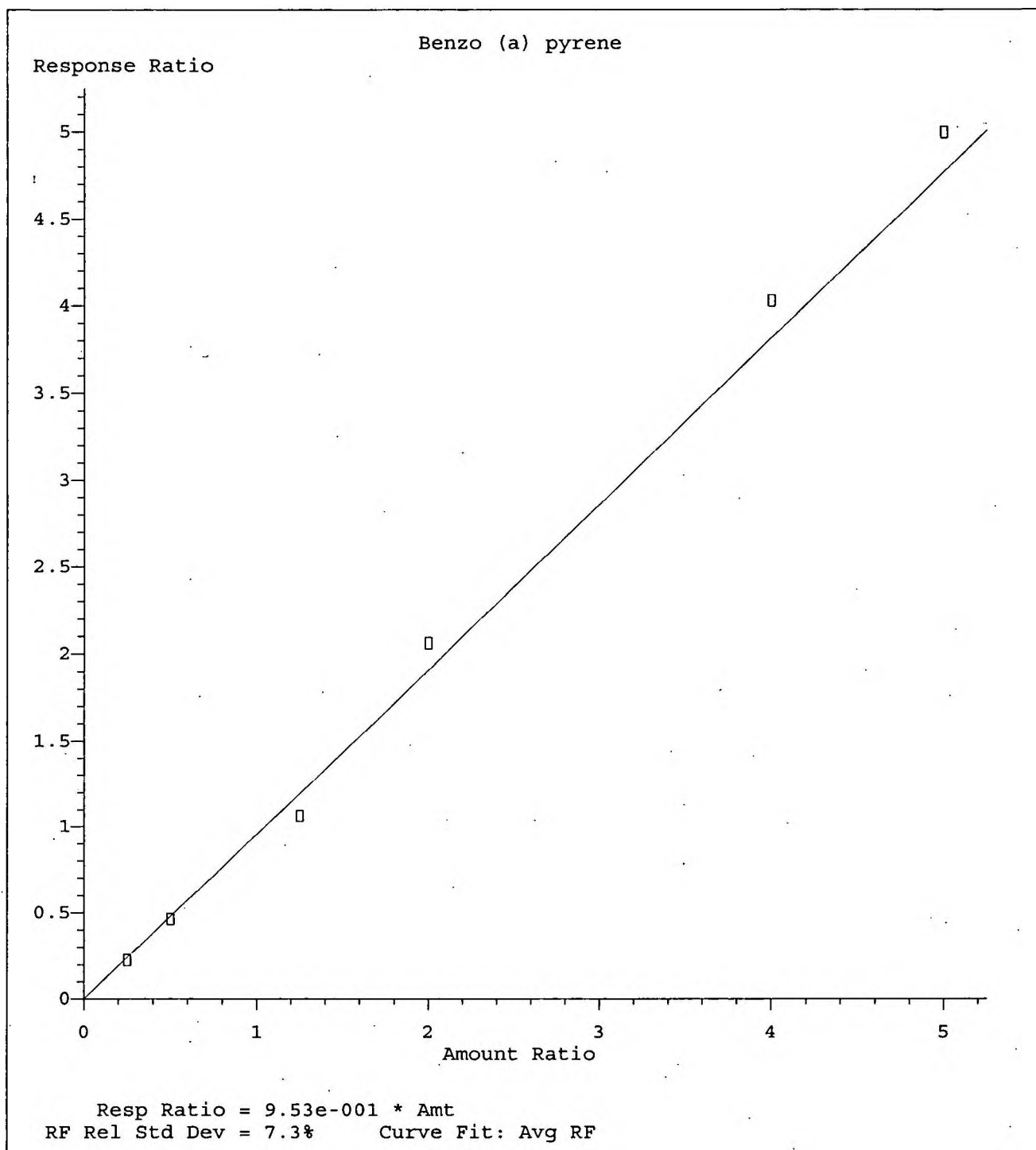
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 Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



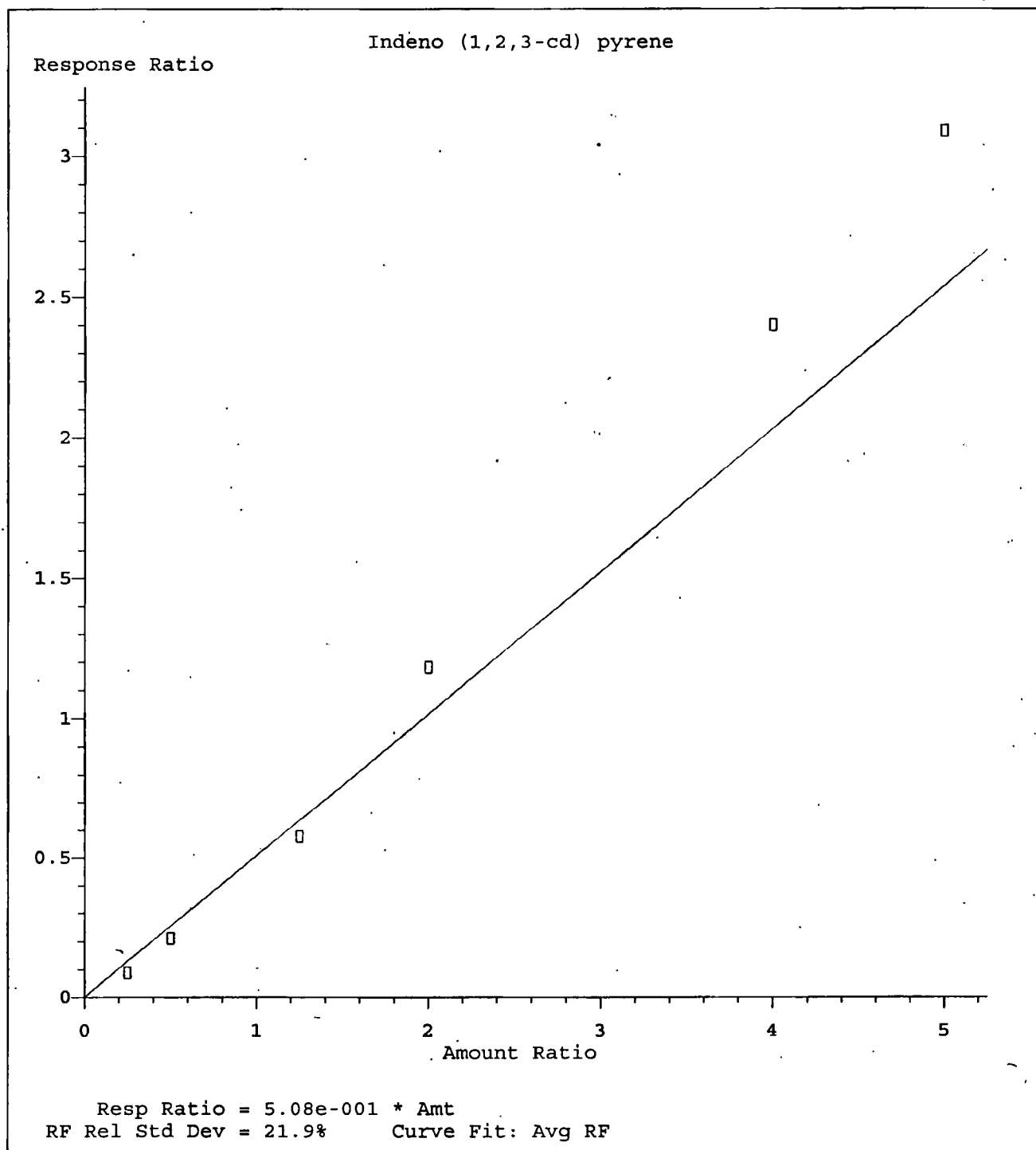
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 Calibration Table Last Updated: Mon Apr 08 06:47:22 2002



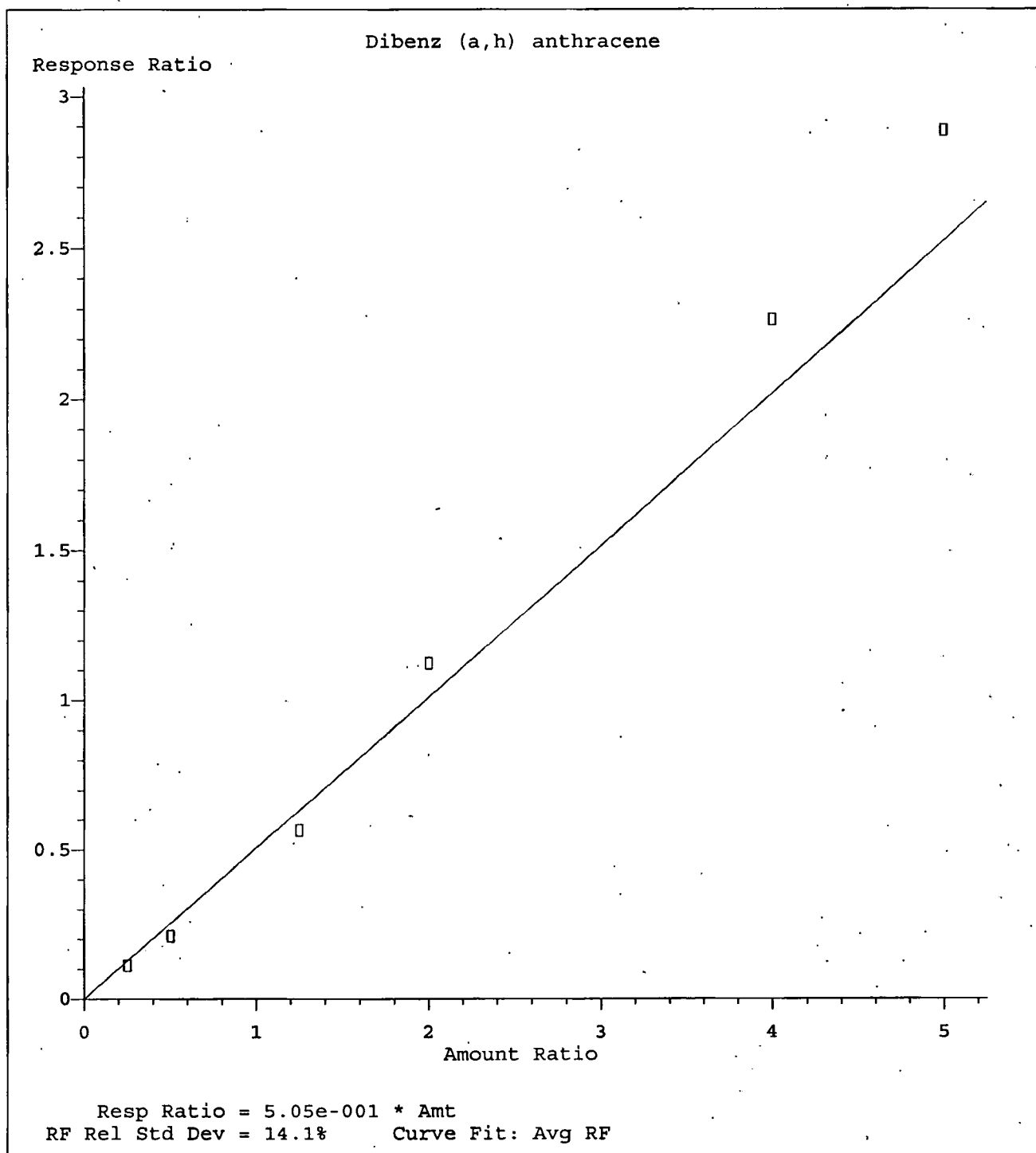
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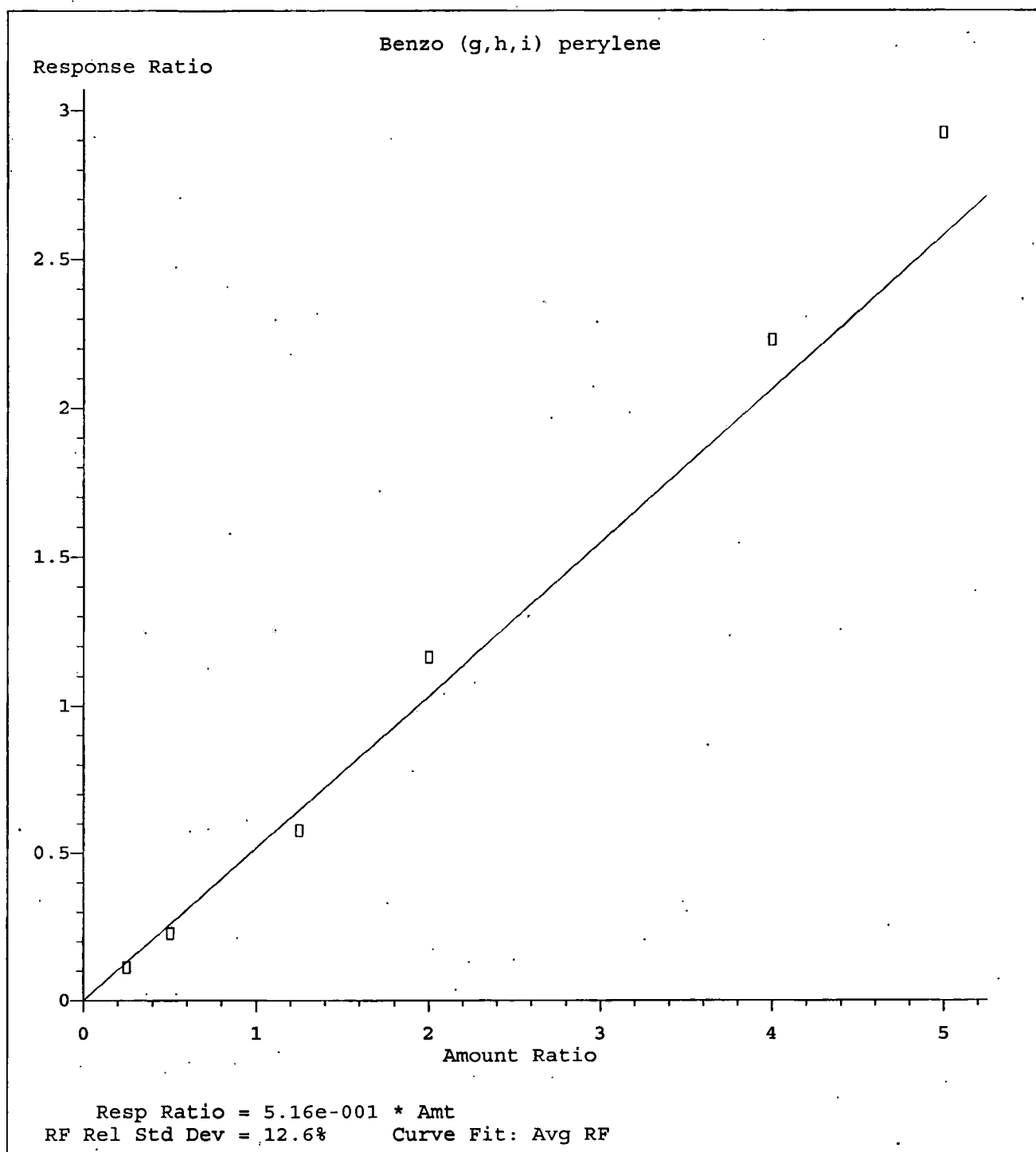
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