

DFTPP

Data File : C:\MSDCHEM\1\DATA\041202\DFTA0412.D

Acq On : 12 Apr 2002 2:05 pm

Sample : DFTPP

Misc :

MS Integration Params: BENZO.P

Vial: 3

Operator: Bohlk

Inst : Instrumen

Multiplr: 1.00

Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

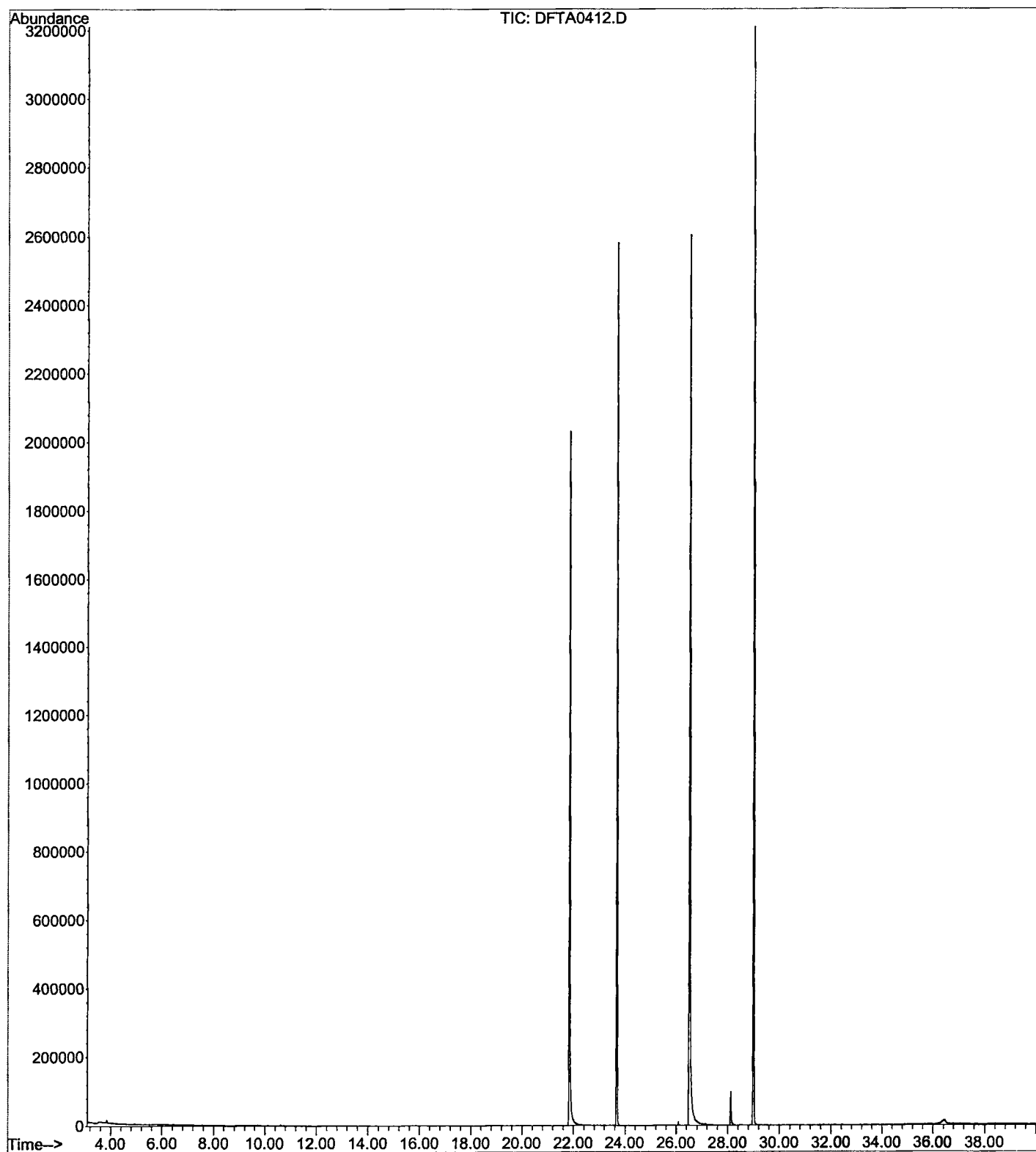
Spectrum Information: Scan 3506

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
51	198	30	60	45.3	145152	PASS
68	69	0.00	2	1.6	2330	PASS
69	198	0.00	100	44.7	143168	PASS
70	69	0.00	2	0.7	946	PASS
127	198	40	60	48.9	156736	PASS
197	198	0.00	1	0.0	0	PASS
198	198	100	100	100.0	320384	PASS
199	198	5	9	6.8	21928	PASS
275	198	10	30	19.5	62544	PASS
365	198	1	100	2.1	6576	PASS
441	443	0.01	100	71.5	26224	PASS
442	198	40	110	58.0	185728	PASS
443	442	17	23	19.7	36680	PASS

DFTA0412.D 5080402BBC.M

Mon Apr 15 10:32:14 2002

File : C:\MSDCHEM\1\DATA\041202\DFTA0412.D
 Operator : Bohlk
 Acquired : 12 Apr 2002 2:05 pm using AcqMethod 508SCAN
 Instrument : Instrumen
 Sample Name: DFTPP
 Misc Info :
 Vial Number: 3



Response Factor Report Instrumen

Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
 Title : Quantitation For Method 508gcscan
 Last Update : Mon Apr 15 14:48:10 2002
 Response via : Initial Calibration

Calibration Files

6 =C100A0412.D 1 =C5A0412.D 2 =C10A0412.D
 3 =C25A0412.D 4 =C40A0412.D 5 =C80A0412.D

	Compound	6	1	2	3	4	5	Avg	%RSD
1) I	1,4-Dichlorobenzene-d	-----ISTD-----							
2)	N-nitroso-dimet	1.190	1.022	1.095	1.128	1.265	1.240	1.157	7.98
3)	bis(2-chloroeth	1.678	1.414	1.568	1.813	1.925	1.779	1.696	10.85
4)	1,3 - Dichlorob	1.250	1.370	1.419	1.458	1.483	1.333	1.386	6.22
5)	1,4 - Dichlorob	1.327	1.526	1.517	1.543	1.565	1.422	1.483	6.12
6)	1,2 - Dichlorob	1.148	1.126	1.171	1.283	1.317	1.229	1.212	6.32
7)	2,2 ' - oxybis	2.145	2.564	2.641	2.662	2.713	2.456	2.530	8.26
8)	N-Nitroso-di-n-	1.186	1.051	1.170	1.280	1.324	1.243	1.209	7.97
9) I	Hexachloroethan	0.568	0.714	0.701	0.681	0.682	0.613	0.660	8.61
10)	Naphthalene-d8	-----ISTD-----							
11)	Nitrobenzene	0.411	0.281	0.326	0.432	0.449	0.430	0.388	17.62
12)	Isophorone	0.711	0.738	0.758	0.833	0.834	0.759	0.772	6.54
13)	bis(2-Chloroeth	0.451	0.427	0.428	0.512	0.527	0.482	0.471	9.00
14)	1,2,4-Trichloro	0.230	0.278	0.259	0.271	0.270	0.245	0.259	7.04
15)	Naphthalene	0.821	1.078	1.039	1.030	1.013	0.889	0.978	10.25
16)	Hexachlorobutad	0.113	0.129	0.122	0.131	0.132	0.121	0.125	5.67
17)	Acenaphthene-d10	-----ISTD-----							
18)	Hexachlorocyclo	0.219	0.124	0.135	0.200	0.221	0.224	0.187	24.33
19)	2-chloronaphtha	0.801	0.686	0.706	0.899	0.921	0.842	0.809	12.04
20) I	Dimethylphthala	1.107	1.104	1.109	1.202	1.223	1.157	1.150	4.56
21)	2,6-Dinitrotolu	0.243	0.162	0.193	0.239	0.257	0.251	0.224	16.98
22)	Acenaphthylene	1.372	1.431	1.457	1.569	1.577	1.469	1.479	5.41
23)	Acenaphthene	0.798	1.016	0.992	0.986	0.972	0.850	0.936	9.53
24)	2,4-Dinitrotolu	0.317	0.191	0.235	0.266	0.297	0.314	0.270	18.34
25)	Diethylphthalat	0.793	1.022	1.018	1.023	1.008	0.866	0.955	10.46
26)	4-Chlorophenyl-	0.472	0.550	0.537	0.558	0.555	0.497	0.528	6.68
27)	Fluorene	0.896	1.157	1.164	1.156	1.130	0.974	1.079	10.69
28) s	TCMX (SURR)	0.158	0.163	0.153	0.166	0.156	0.161	0.160	3.06
29)	Azobenzen/ 1,2-	1.541	1.704	1.668	1.737	1.747	1.569	1.661	5.25
30)	Phenanthrene-d10	-----ISTD-----							
31)	N-Nitrosodiphen	0.346	0.467	0.458	0.464	0.445	0.379	0.426	12.02
32)	4-Bromophenyl-p	0.172	0.180	0.184	0.191	0.201	0.182	0.185	5.45
33)	Hexachlorobenze	0.157	0.180	0.177	0.177	0.182	0.165	0.173	5.51
34)	Phenanthrene	0.947	1.210	1.183	1.150	1.150	1.011	1.108	9.44
35)	Anthracene	0.997	1.166	1.187	1.182	1.185	1.067	1.131	7.08
36)	Di-n-butylphtha	1.124	1.234	1.296	1.364	1.374	1.213	1.268	7.59
37)	Fluoranthene	1.096	1.102	1.151	1.244	1.296	1.165	1.176	6.76
38) s	2,4-DDD (SURR)	0.256	0.285	0.347	0.299	0.311	0.289	0.298	10.11
39)	Chrysene-d12	-----ISTD-----							
40)	Pyrene	1.470	1.502	1.491	1.520	1.576	1.493	1.508	2.44
41)	Butylbenzylphth	0.713	0.580	0.643	0.738	0.776	0.731	0.697	10.35
42)	Benzo (a) antra	1.238	1.109	1.144	1.248	1.295	1.277	1.218	6.15
43)	Bis (2-ethylhex	0.942	0.924	0.984	1.023	1.074	0.992	0.990	5.51
44)	Chrysene	1.140	1.203	1.229	1.233	1.226	1.204	1.206	2.88
45)	Perylene-d12	-----ISTD-----							
46)	Di-n-Octylphtha	2.522	2.042	2.283	2.668	2.836	2.578	2.488	11.42
47)	Benzo (b) fluor	1.541	1.346	1.482	1.602	1.680	1.594	1.541	7.53

(#) = Out of Range

5080402BBC.M

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Page 1

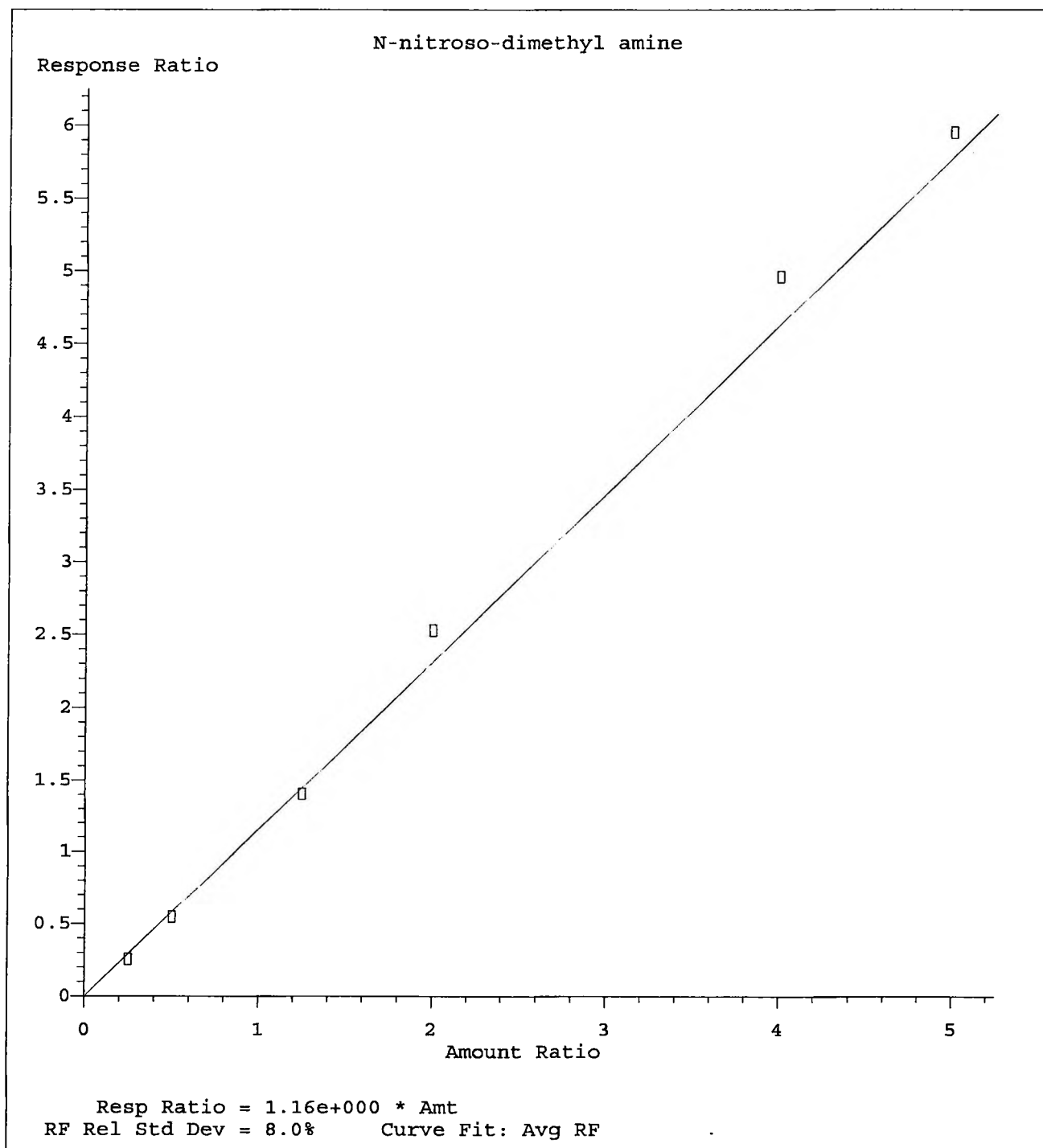
Response Factor Report Instrumen

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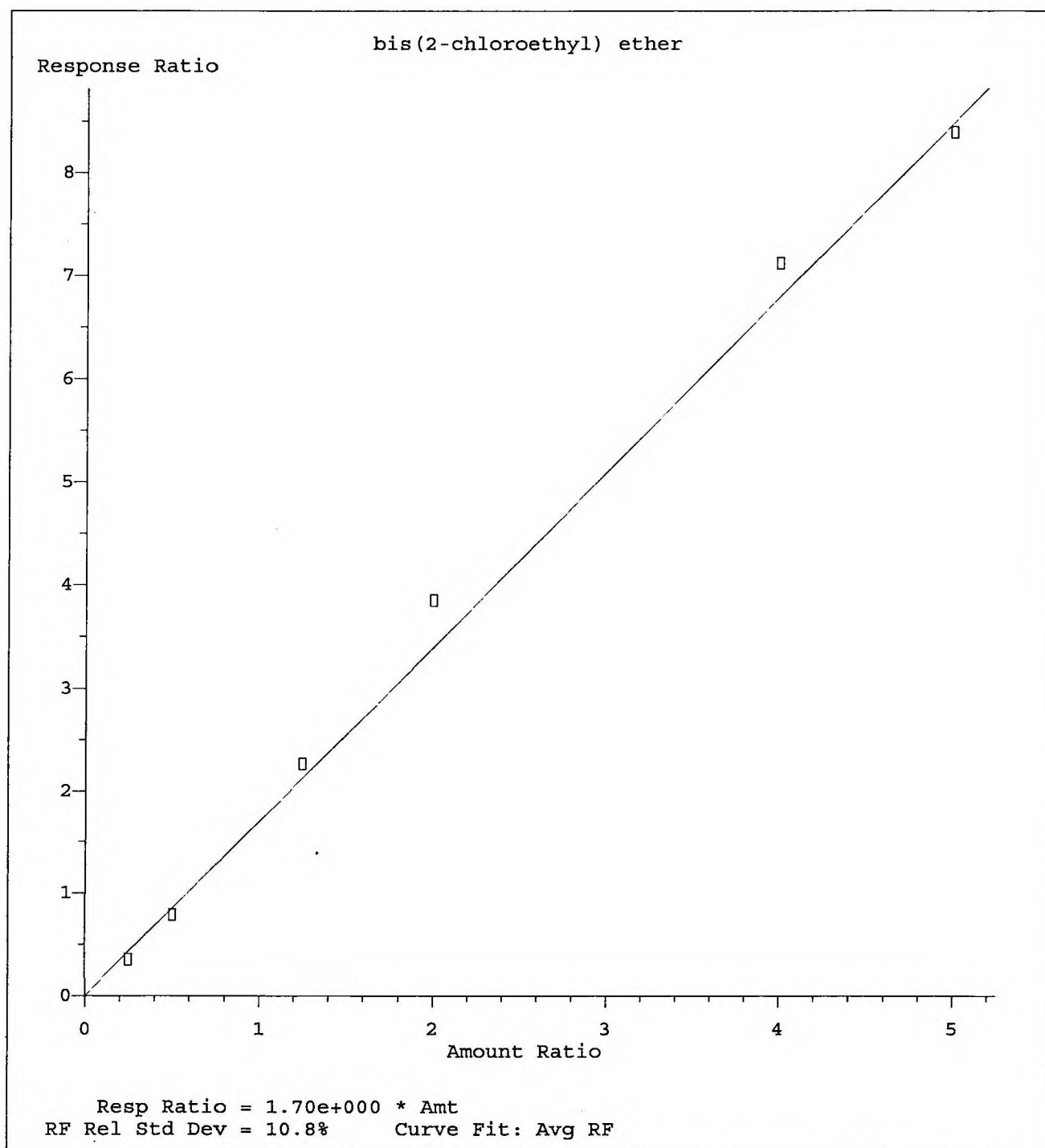
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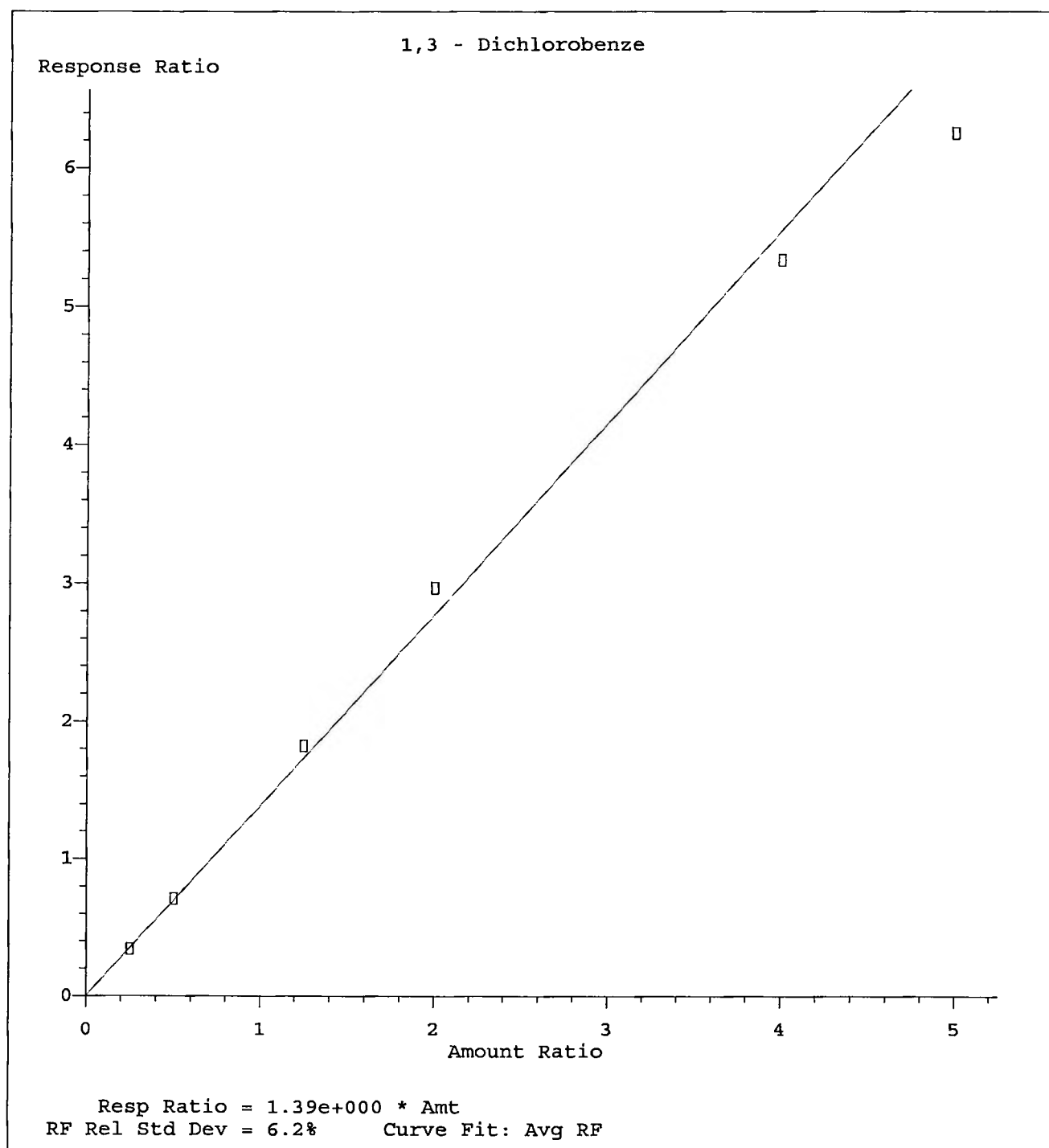
	Compound	6	1	2	3	4	5	Avg	%RSD
48)	Benzo (k) fluor	1.583	1.752	1.825	1.825	1.820	1.676	1.747	5.70
49)	Benzo (a) pyren	1.500	1.352	1.489	1.580	1.614	1.536	1.512	6.06
50)	Indeno (1,2,3-c	0.856	0.569	0.701	0.843	0.886	0.906	0.794	16.56
51)	Dibenz (a,h) an	0.835	0.645	0.752	0.829	0.890	0.865	0.802	11.25
52)	Benzo (g,h,i) p	0.814	0.770	0.849	0.882	0.906	0.858	0.846	5.76



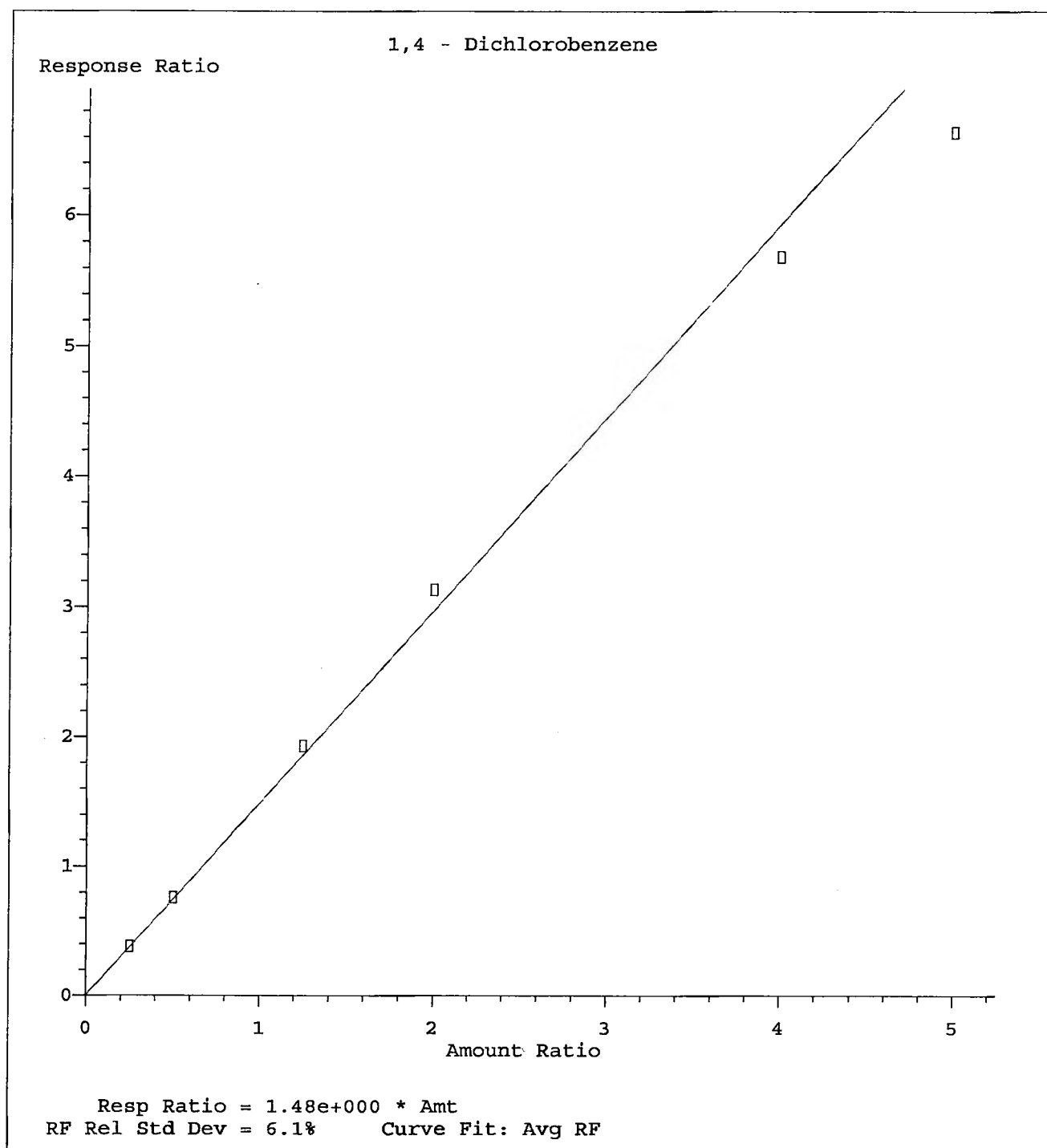
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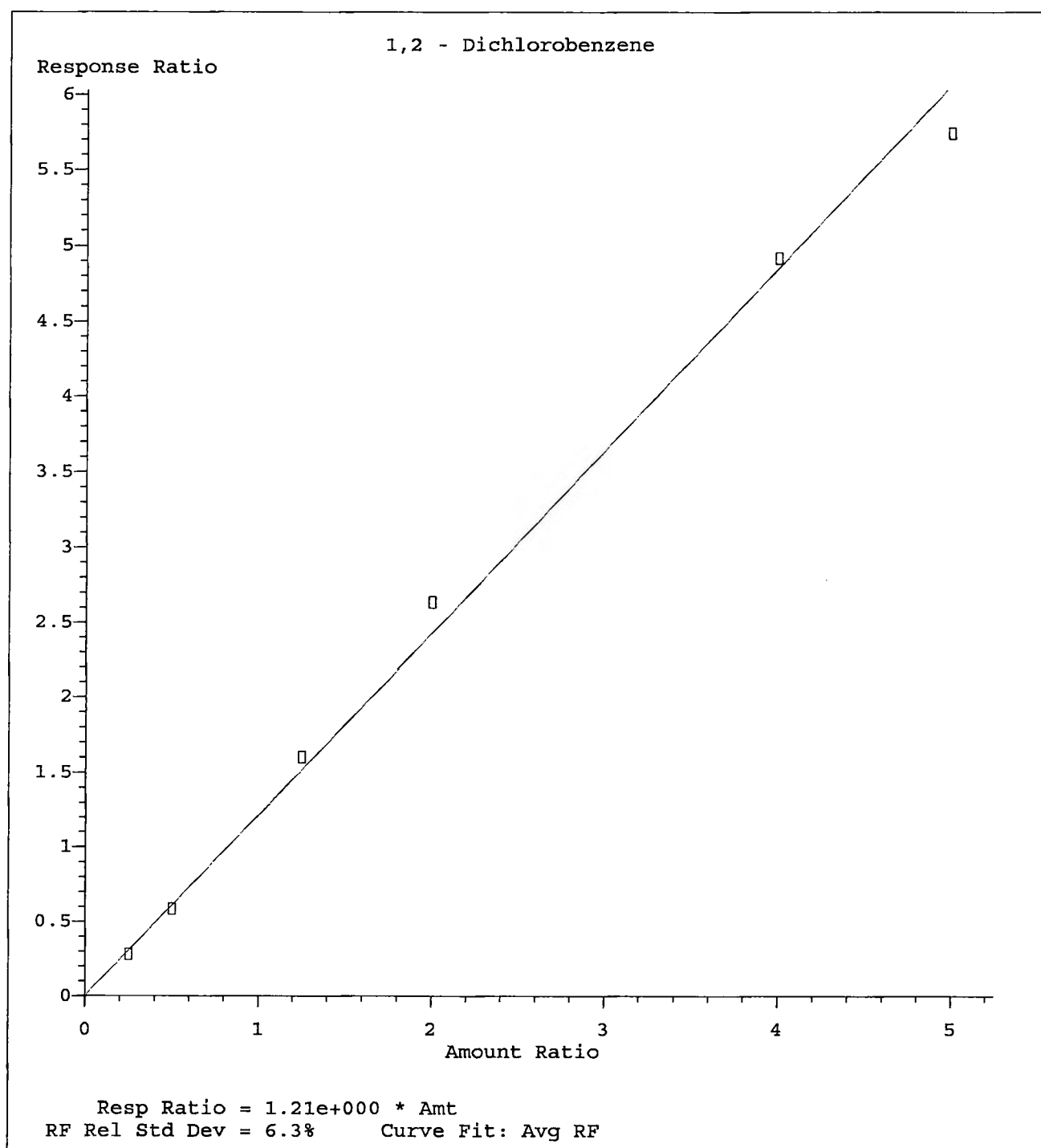
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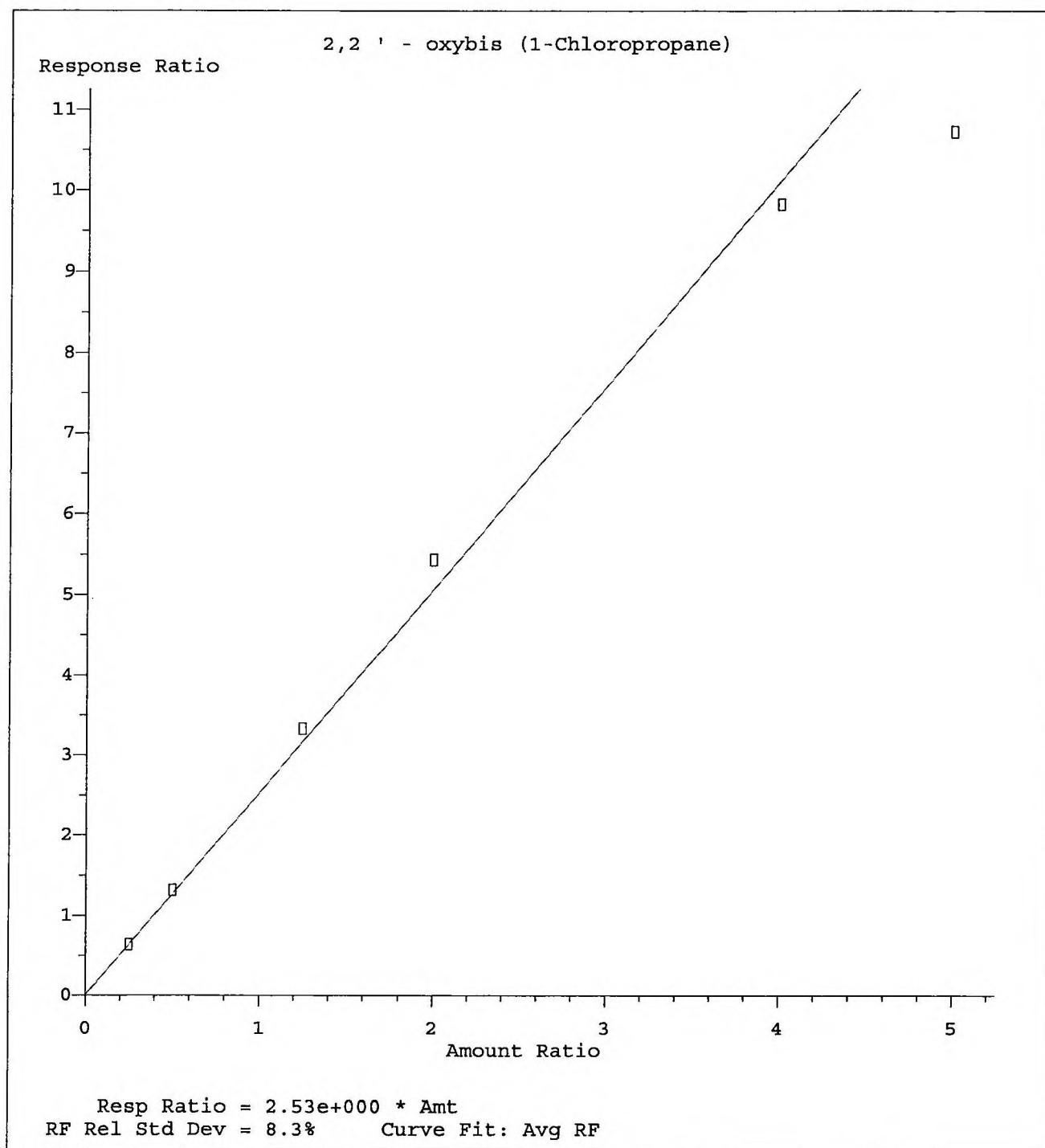
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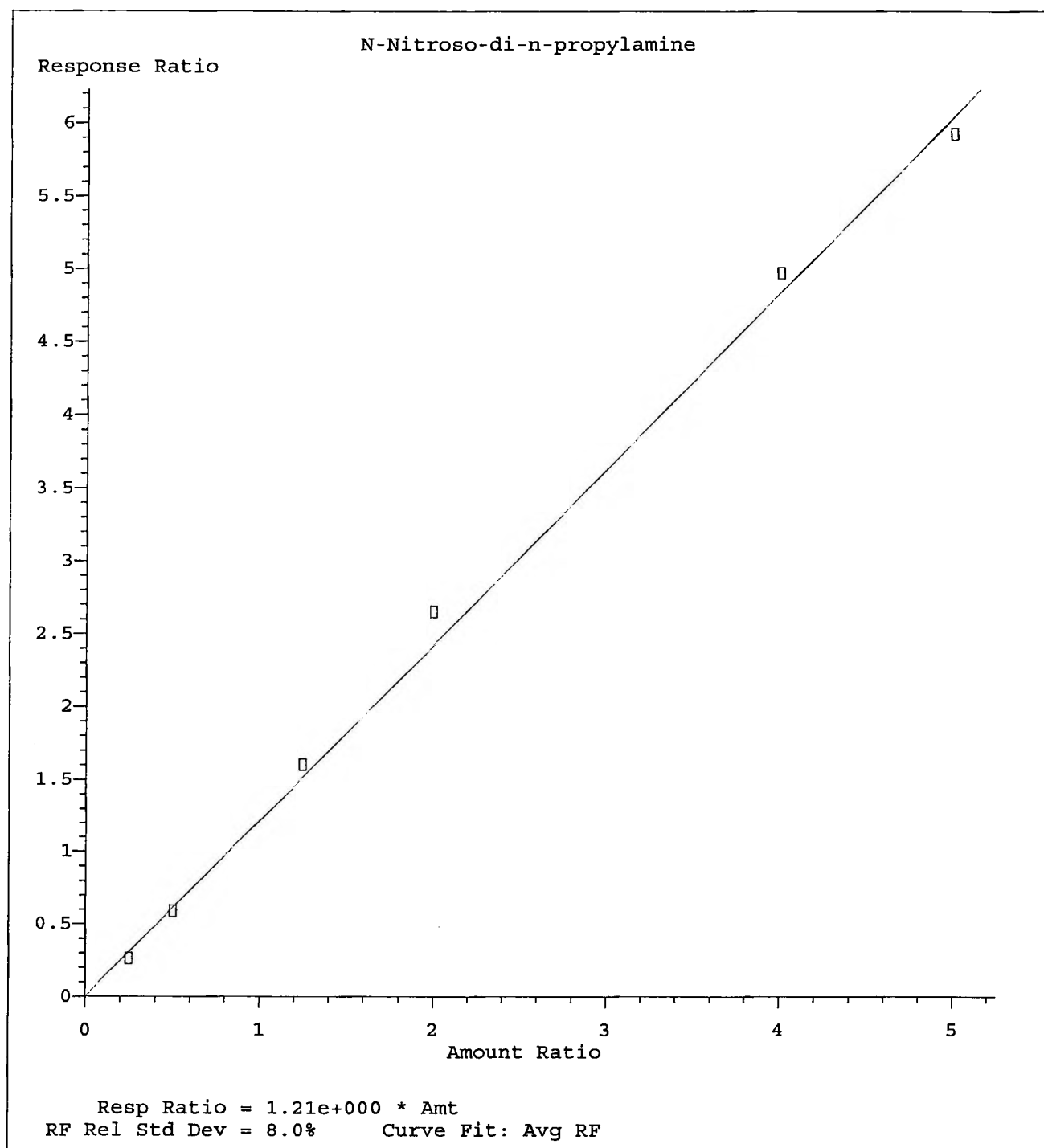
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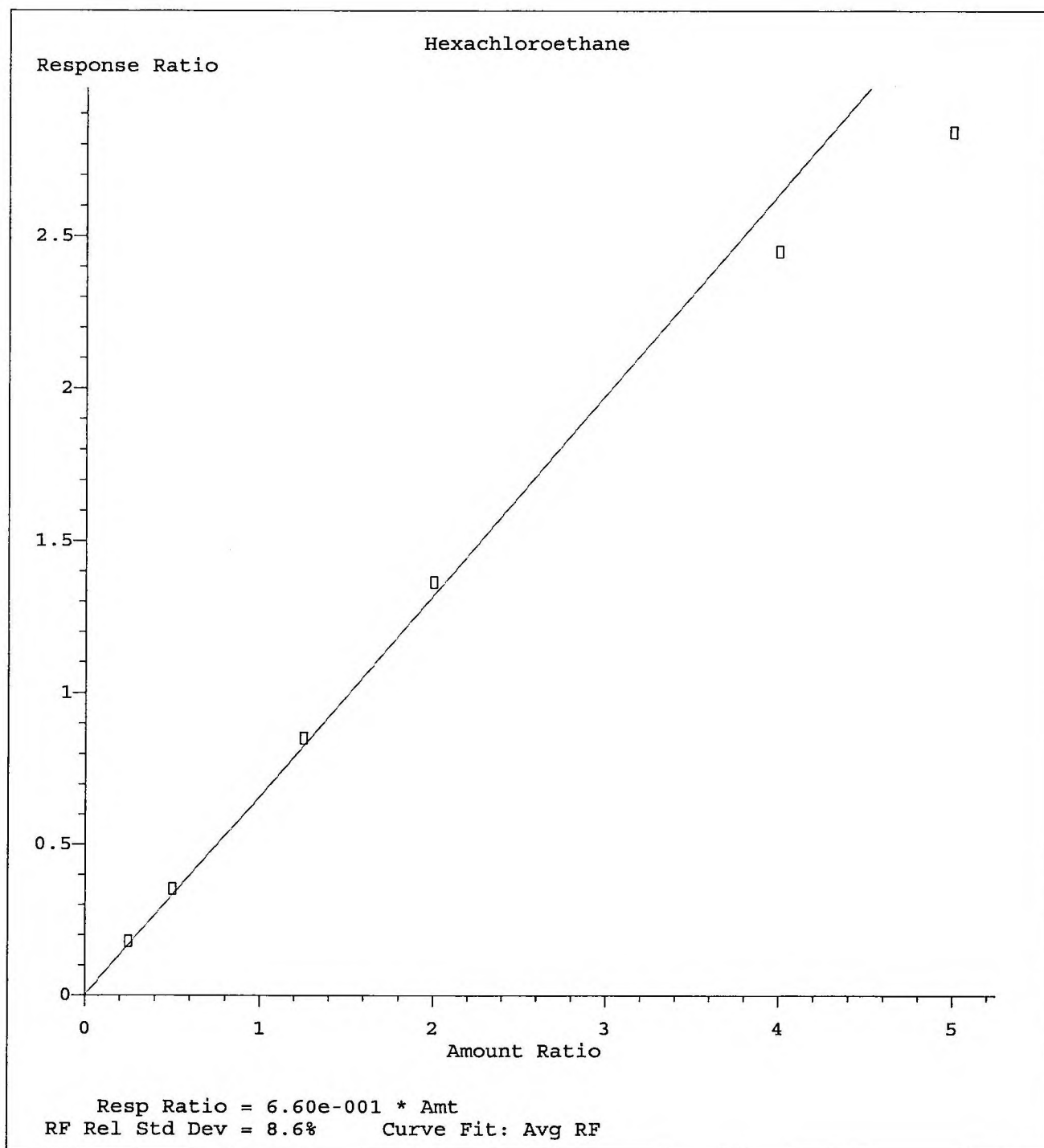
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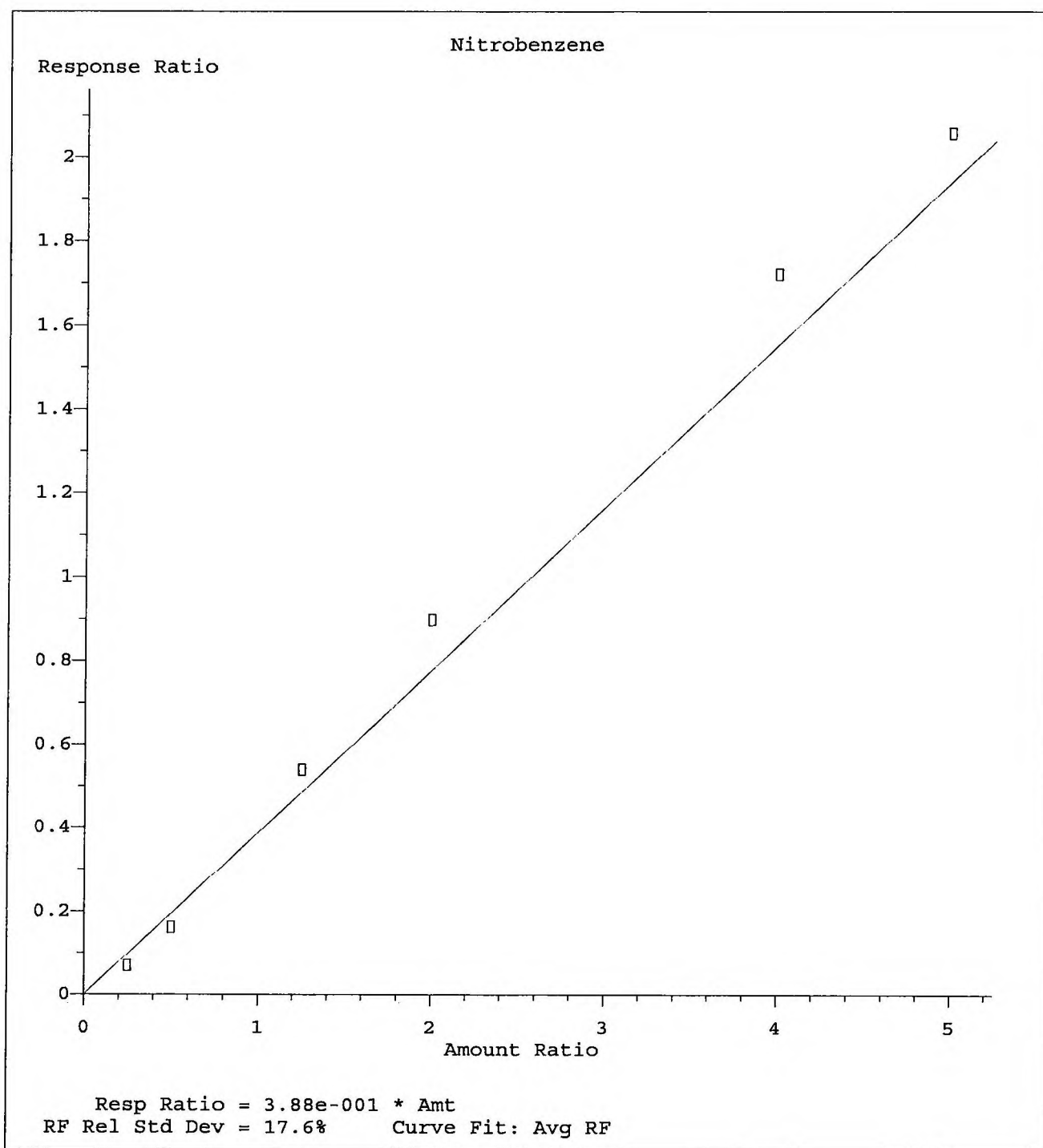
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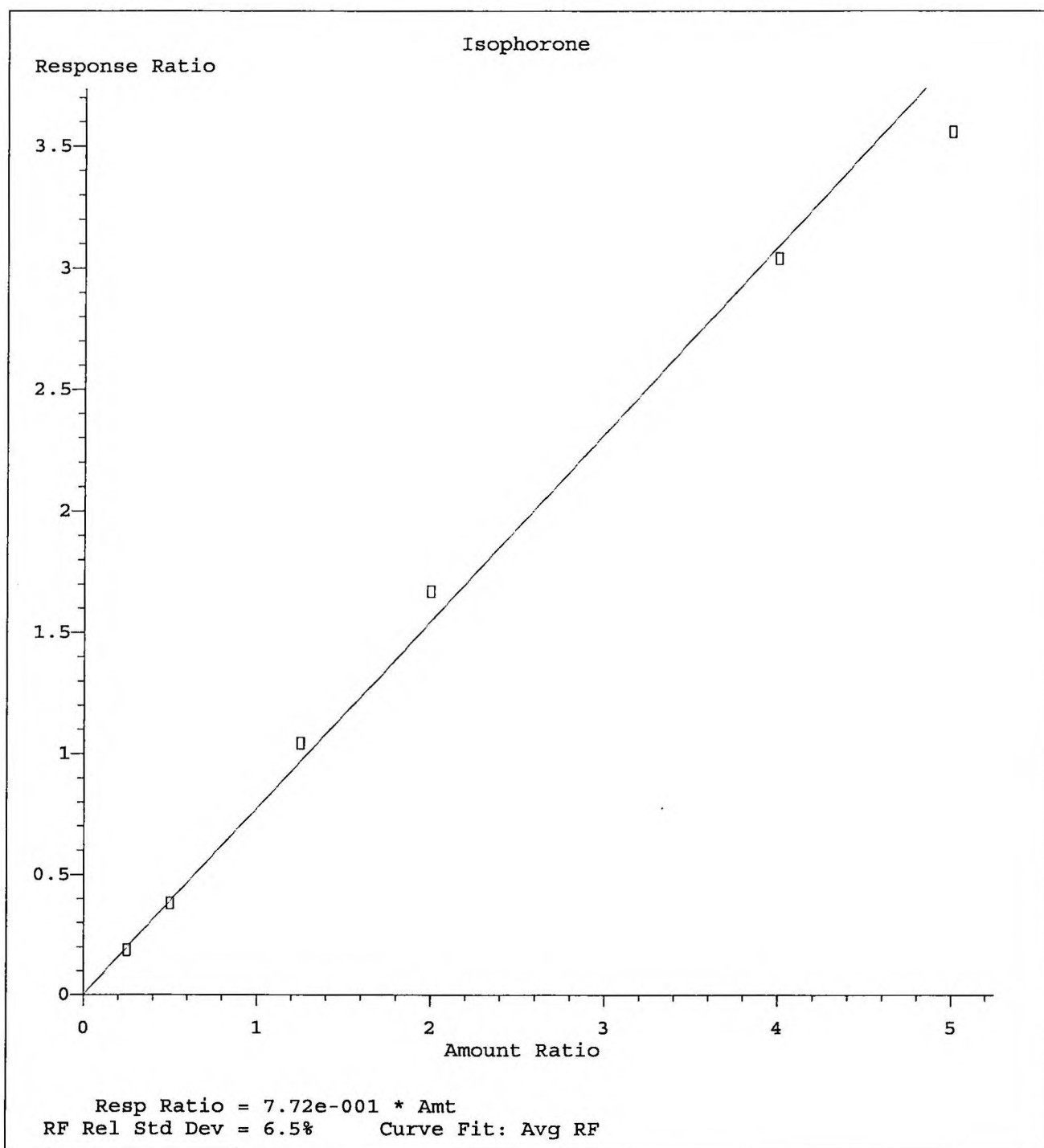
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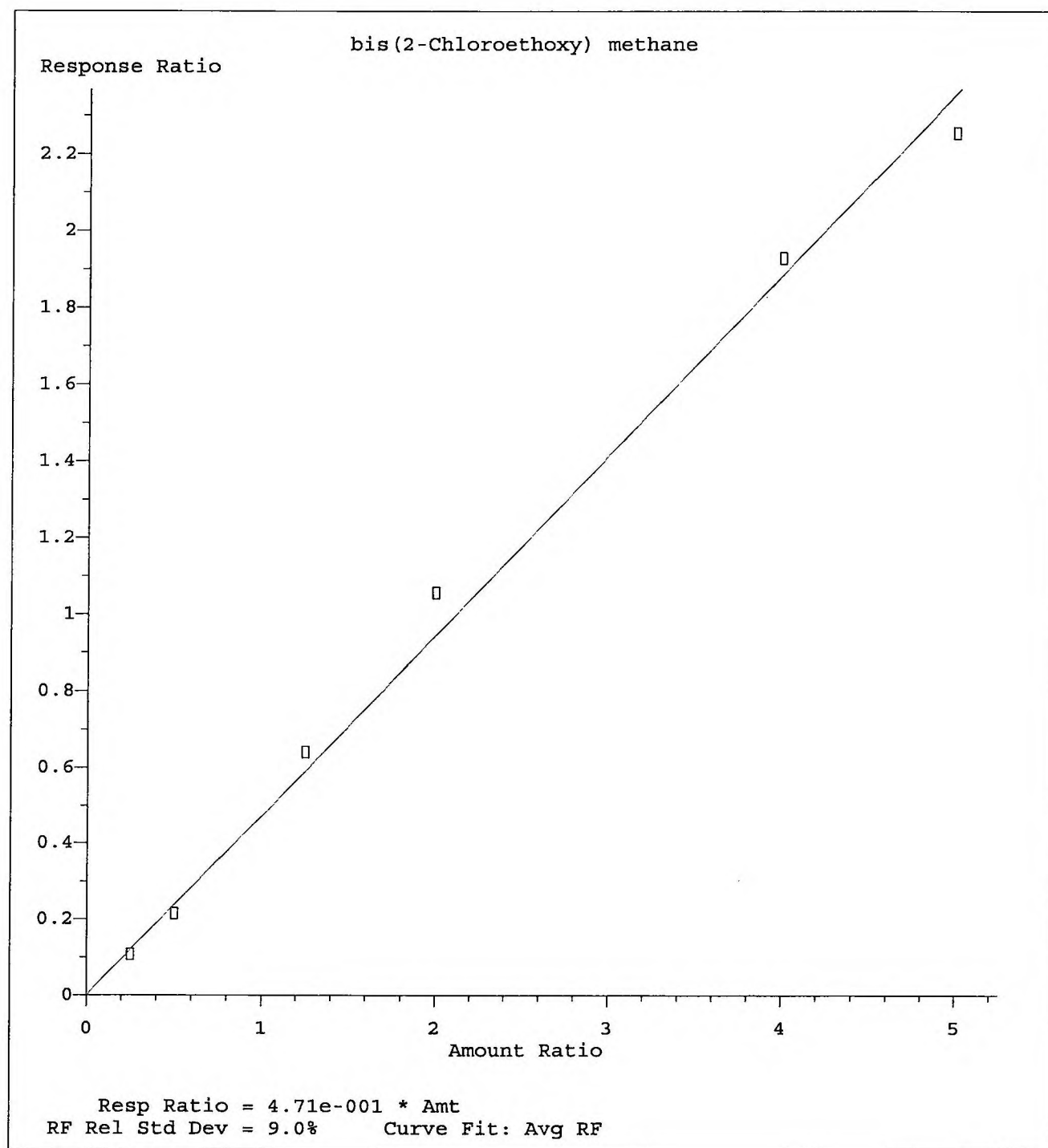
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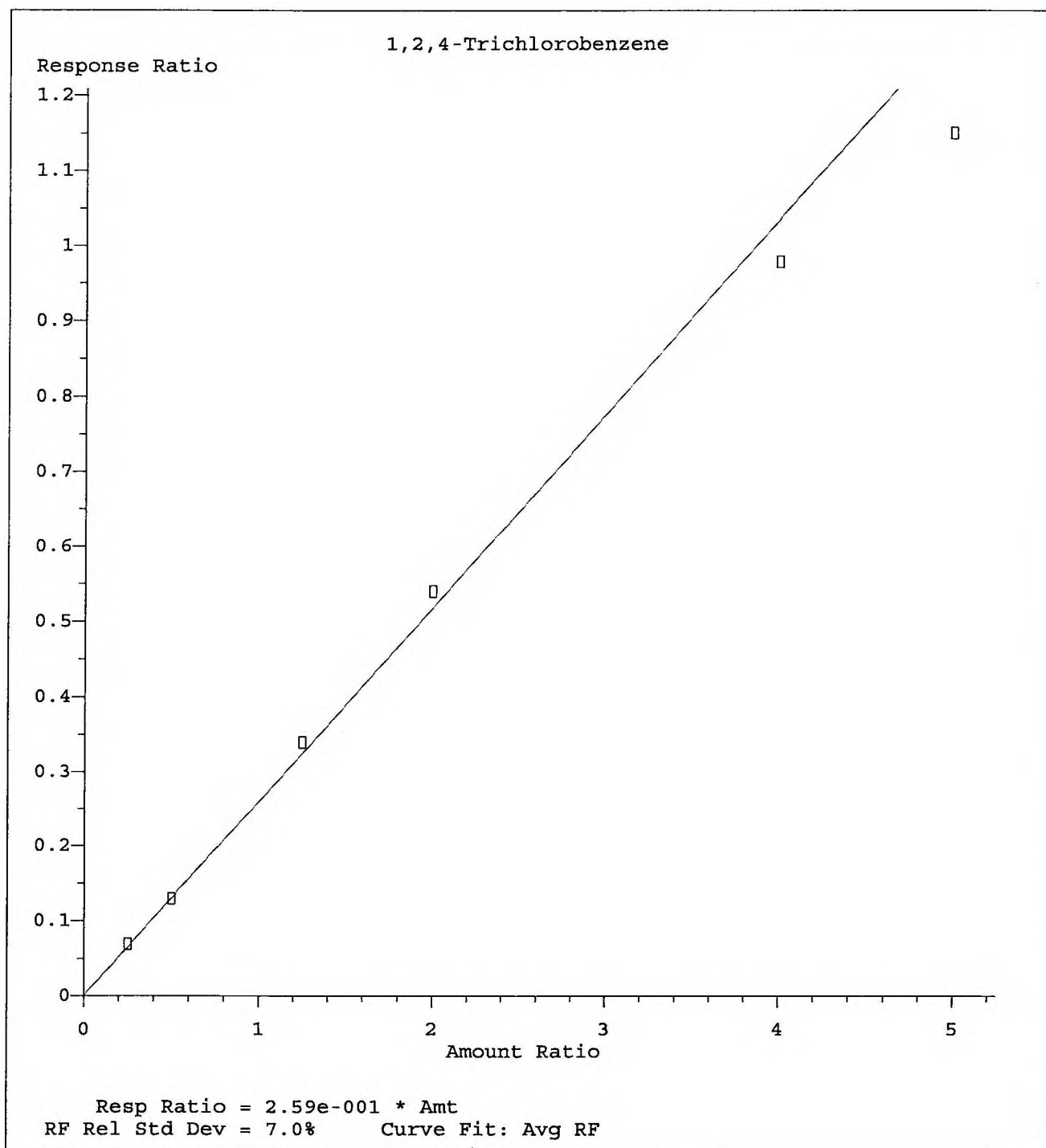
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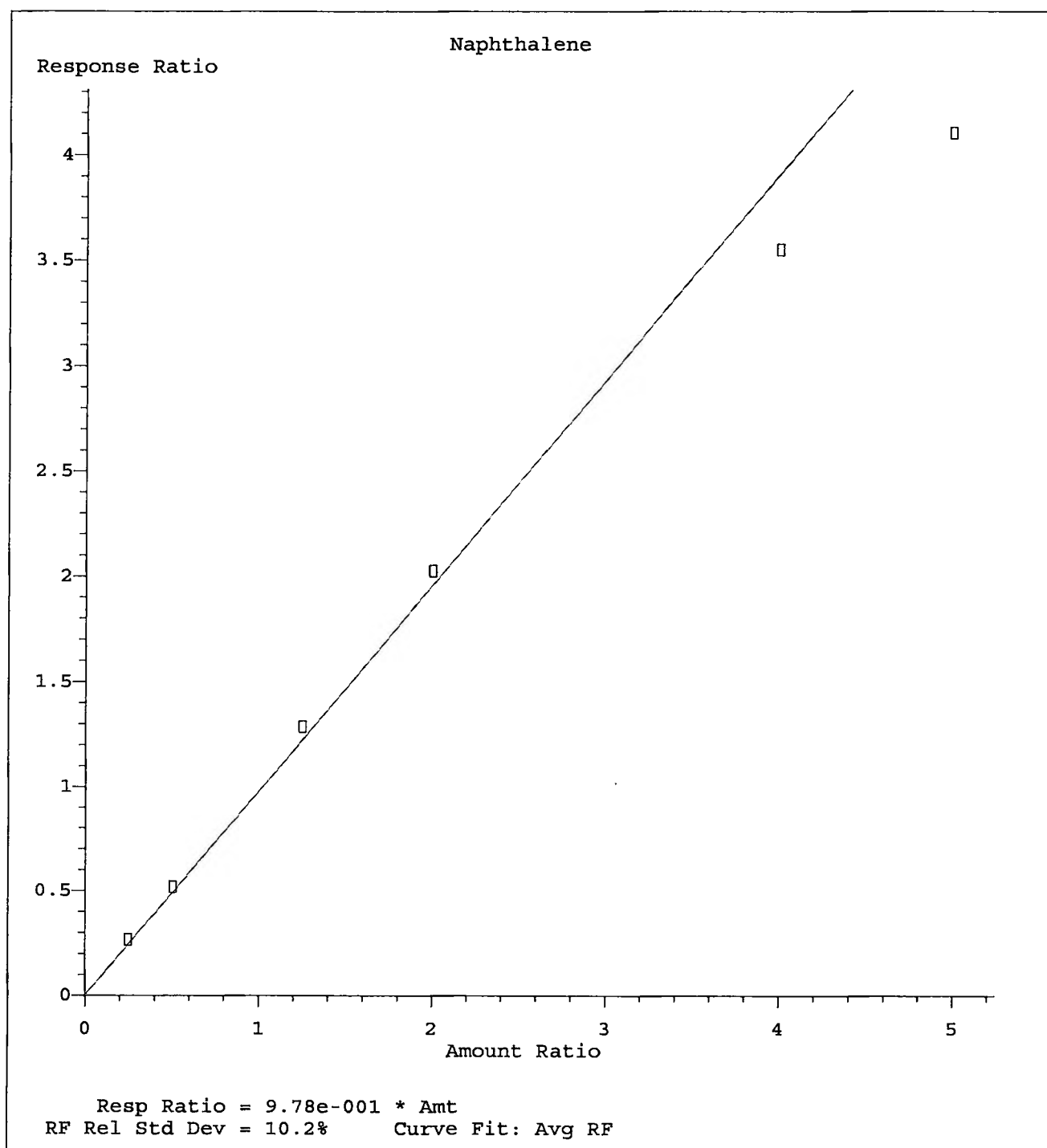
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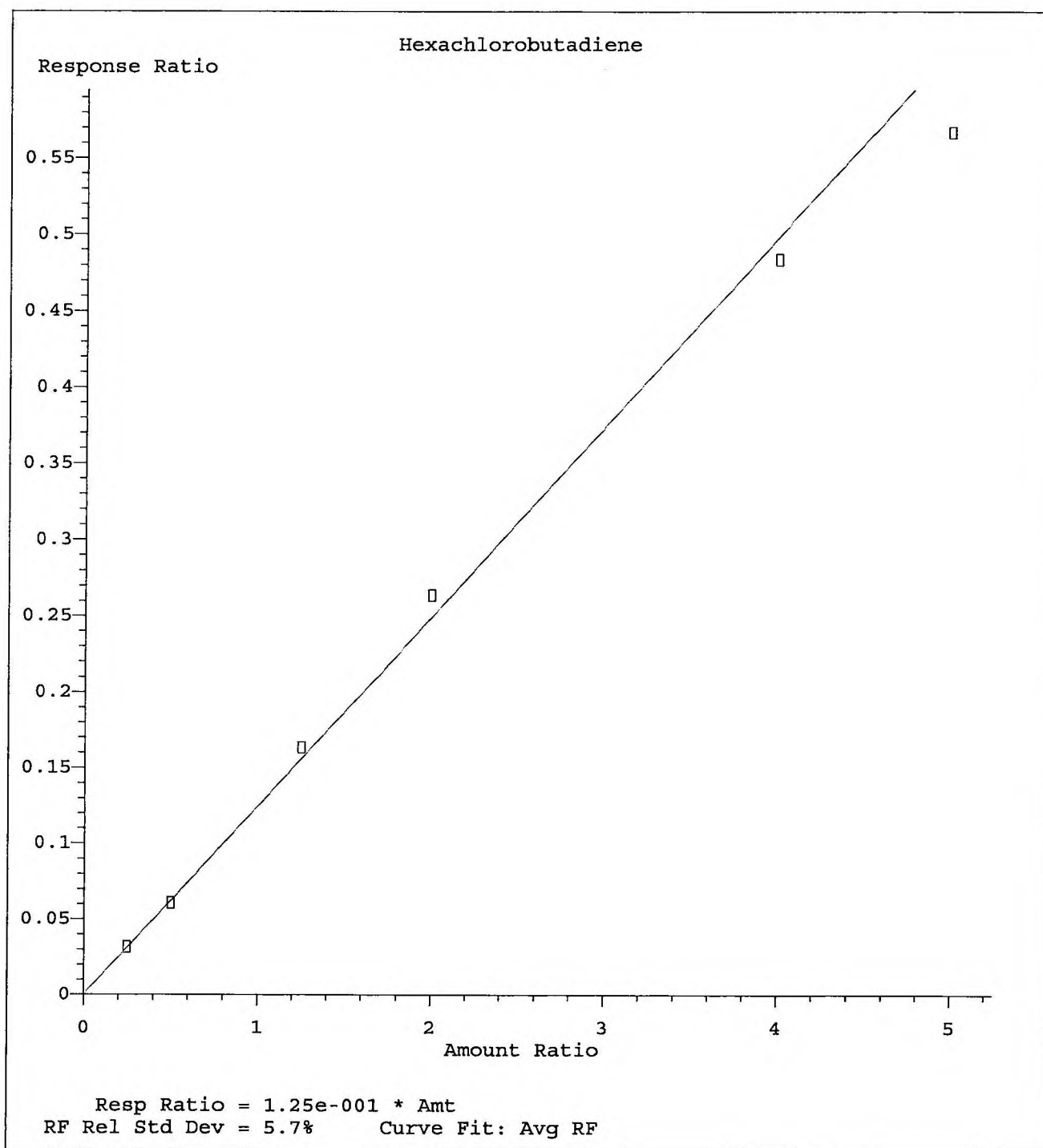
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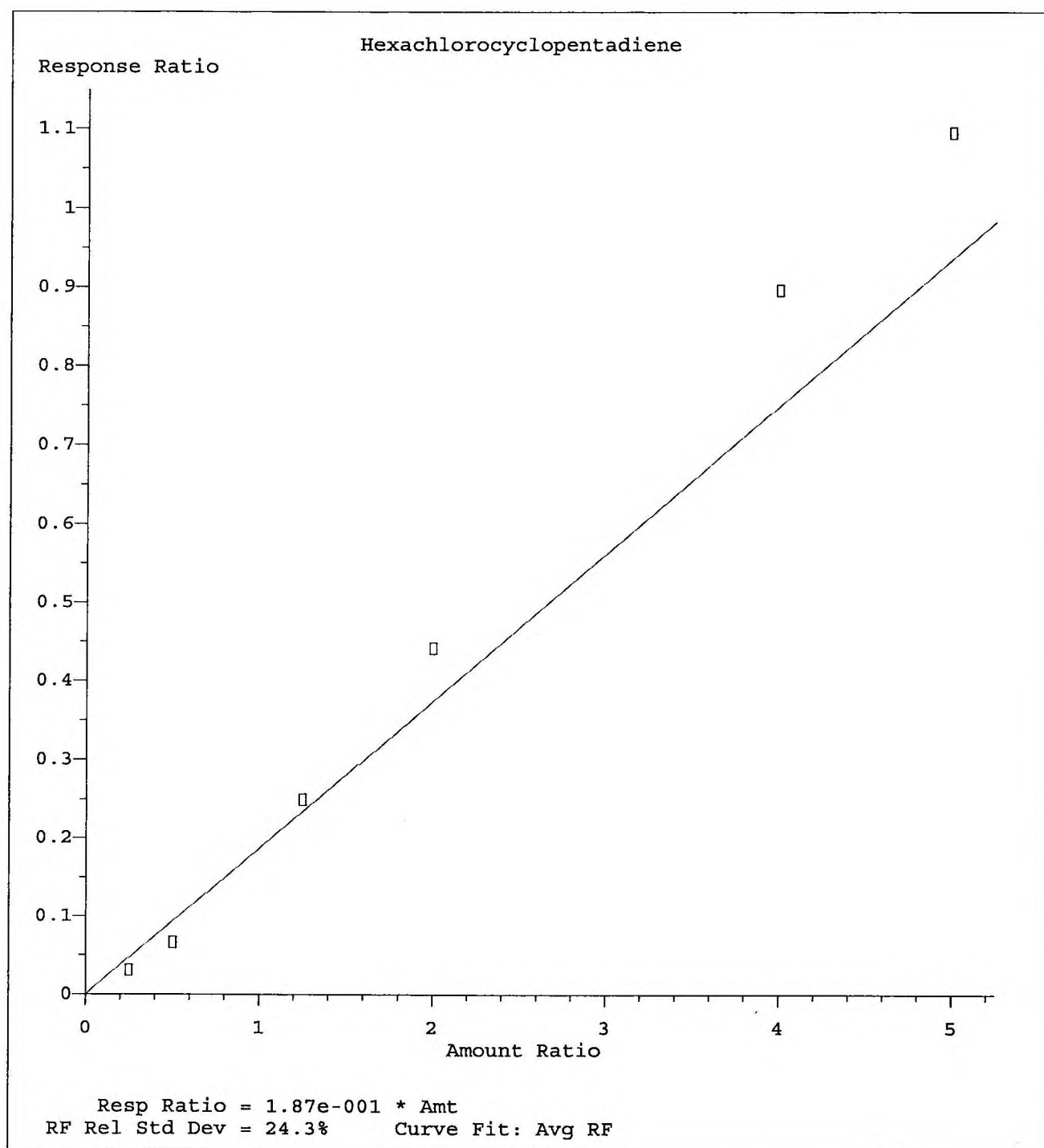
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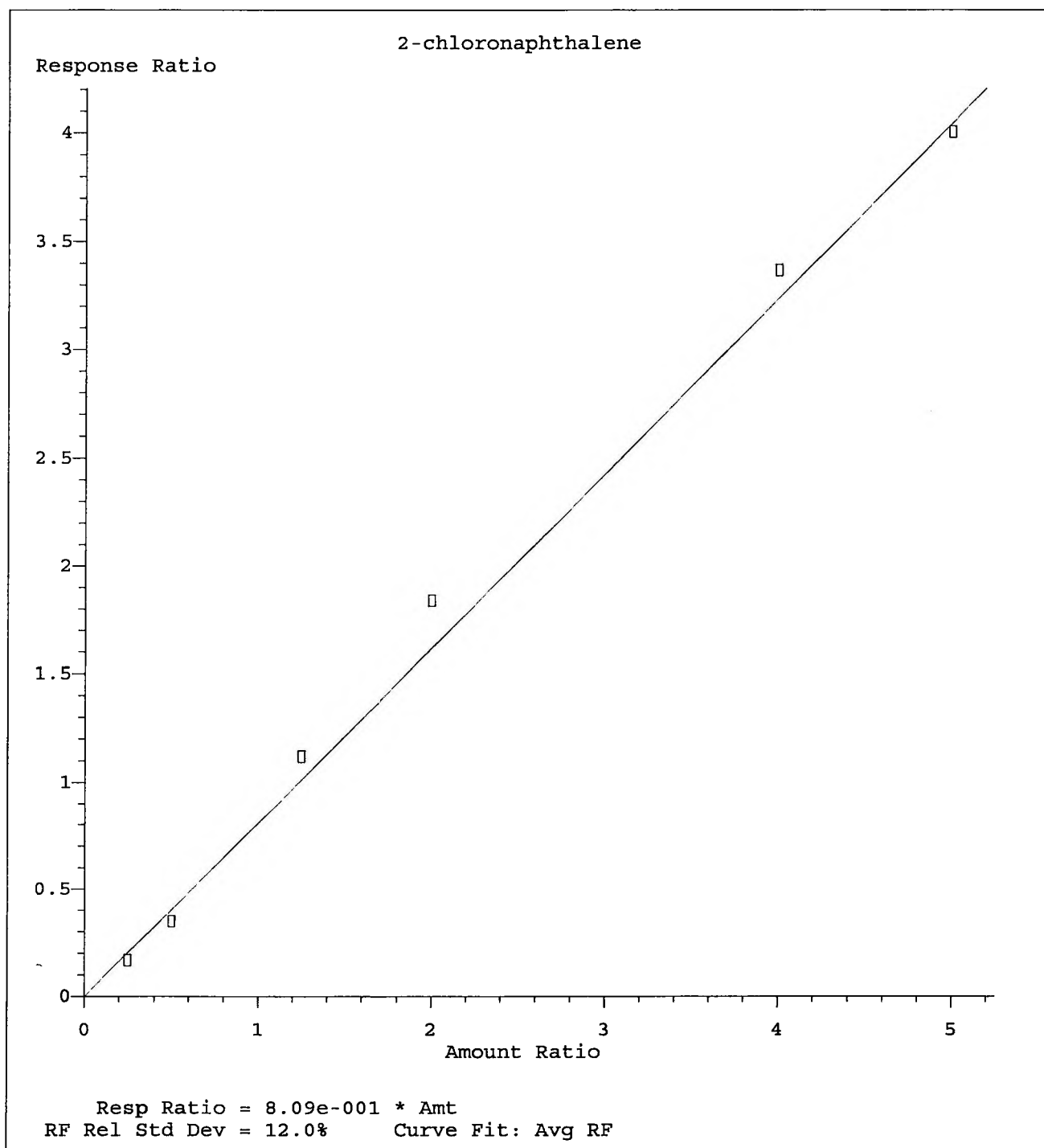
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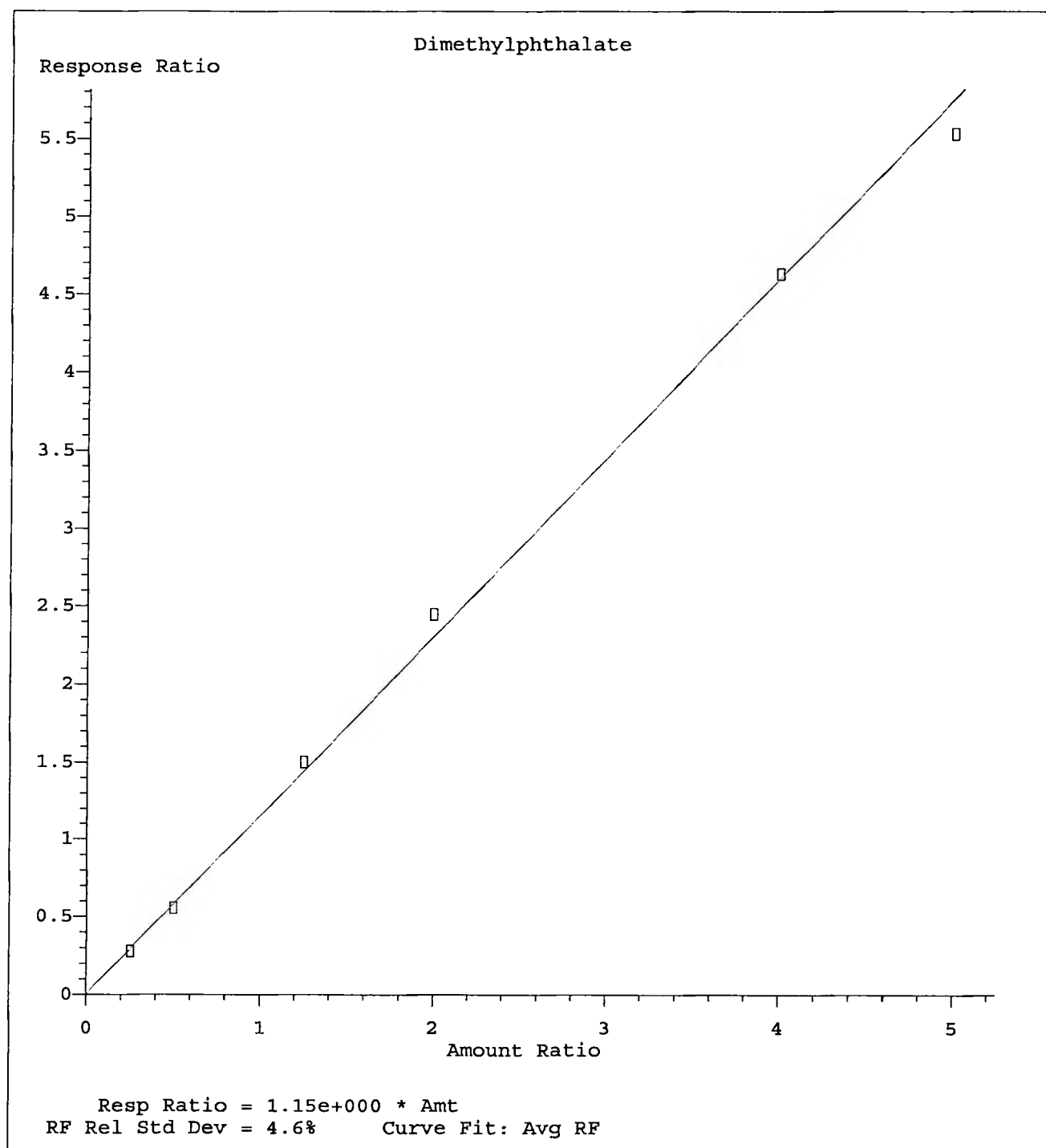
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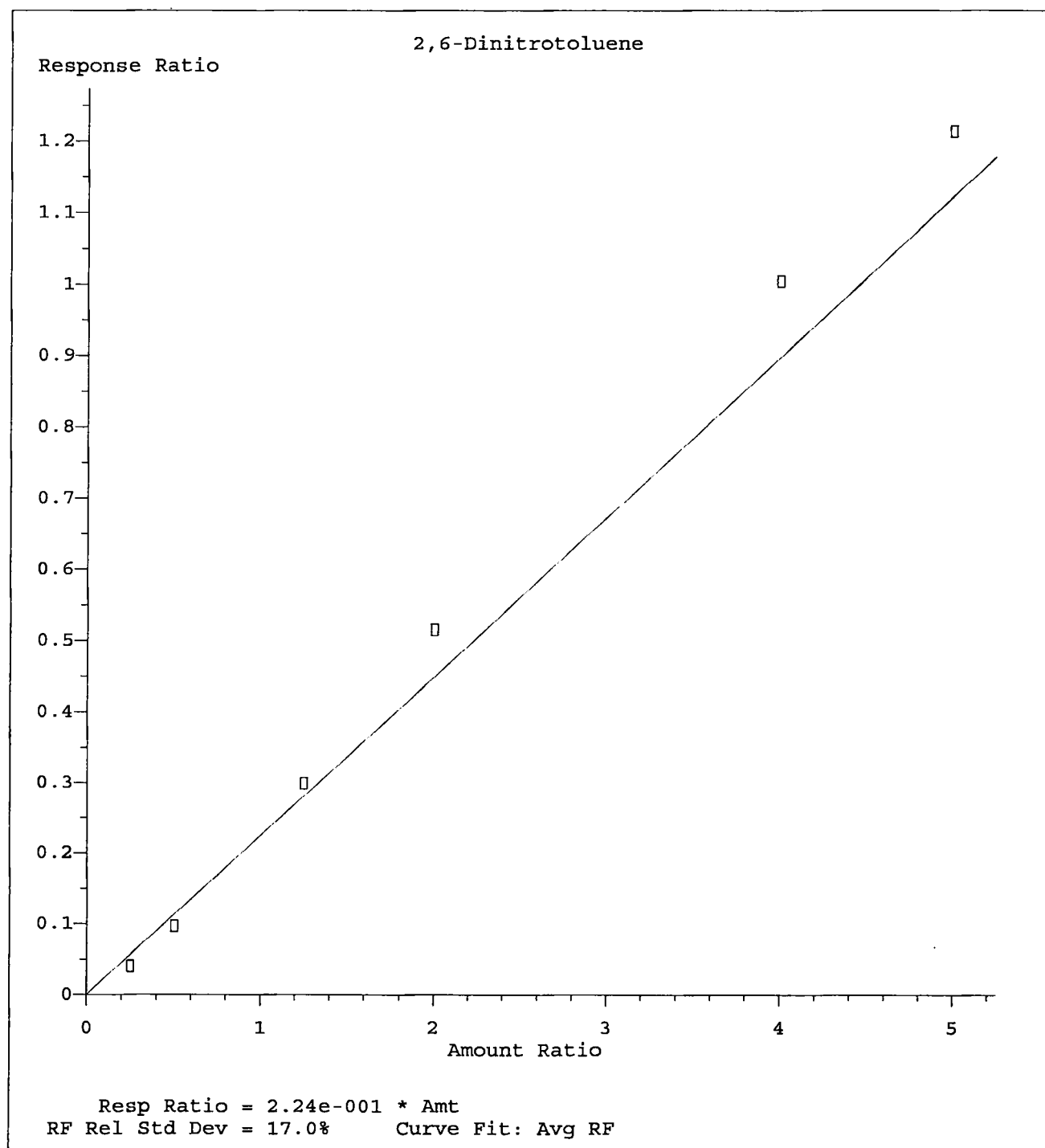
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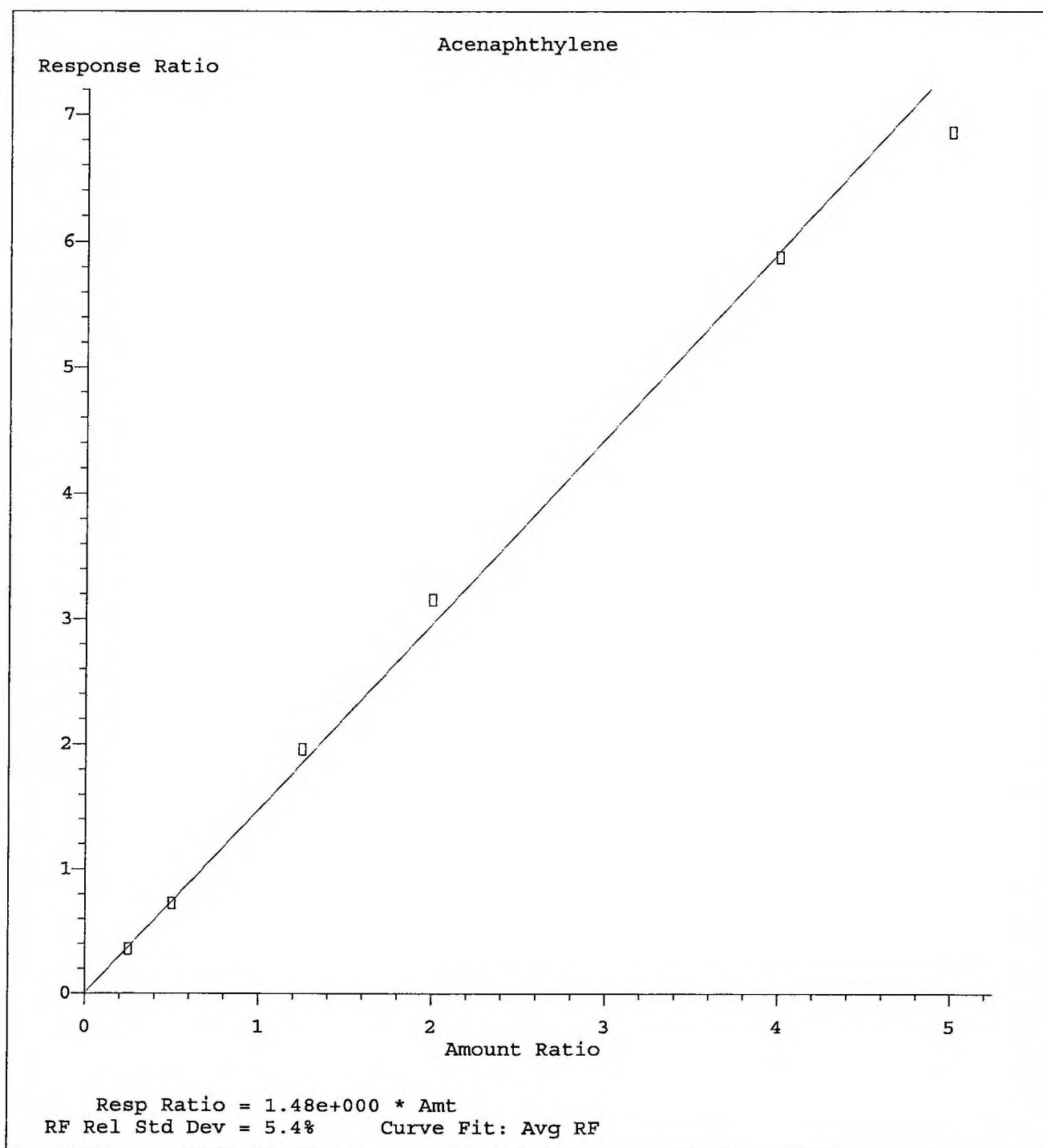
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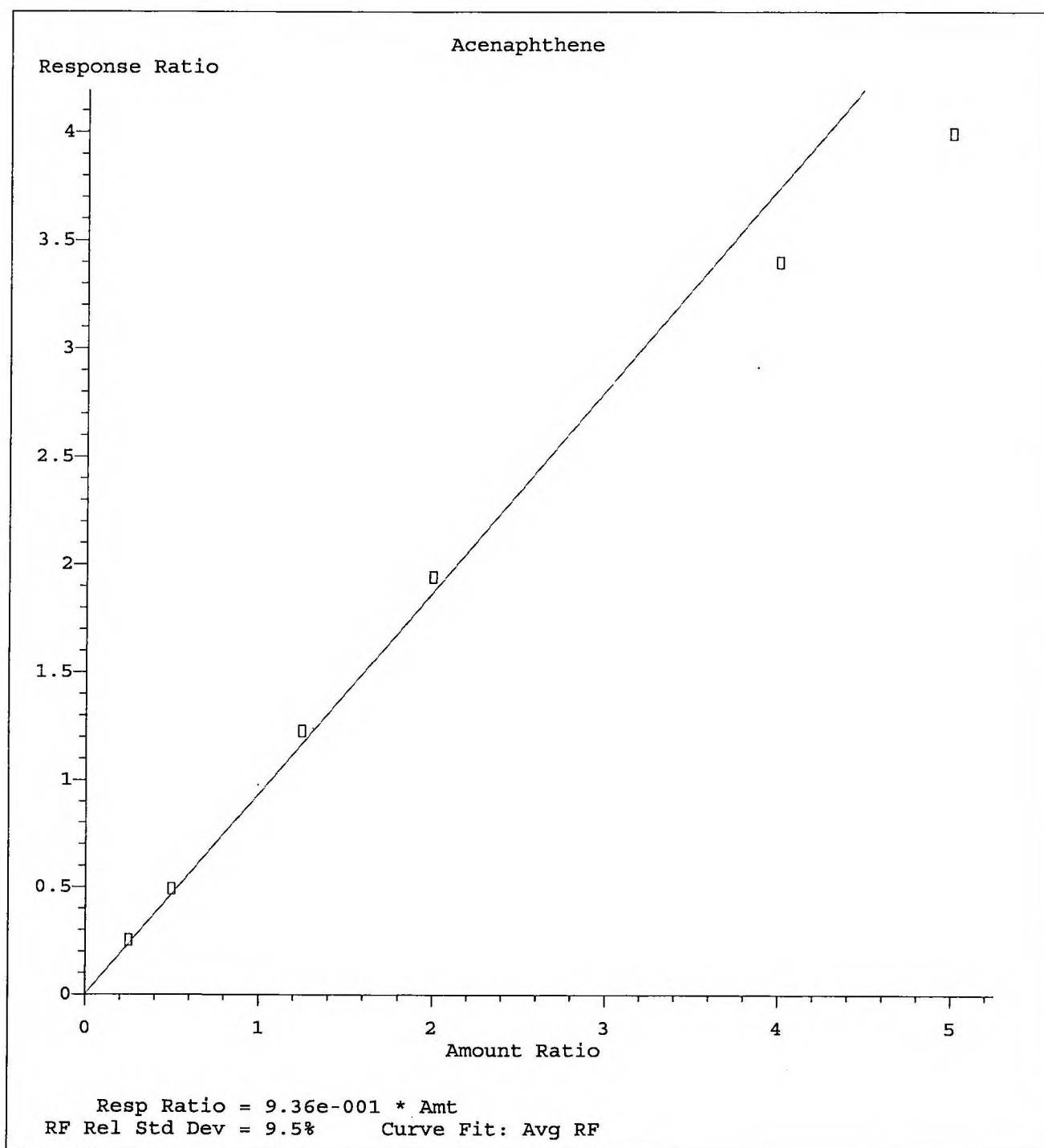
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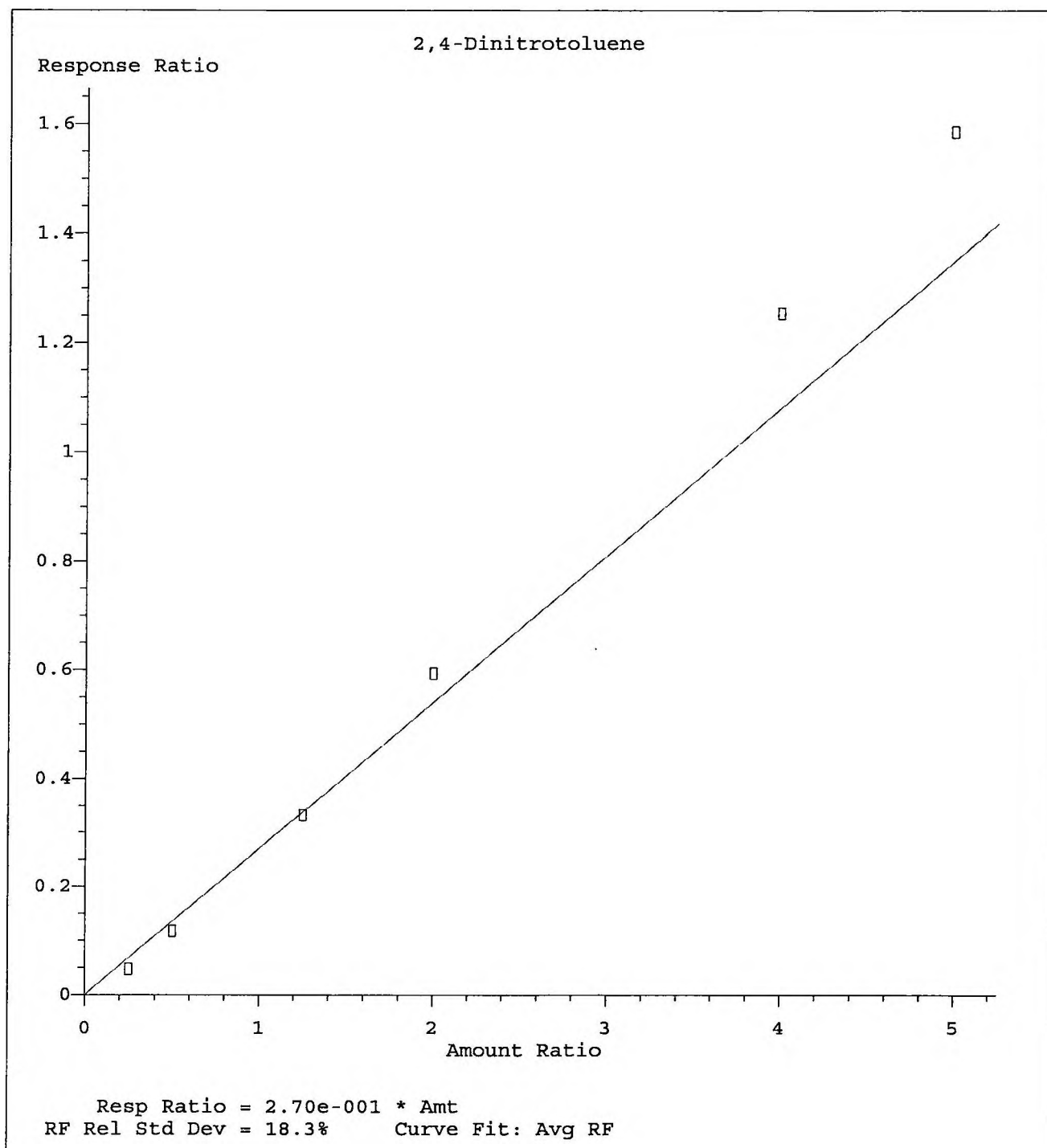
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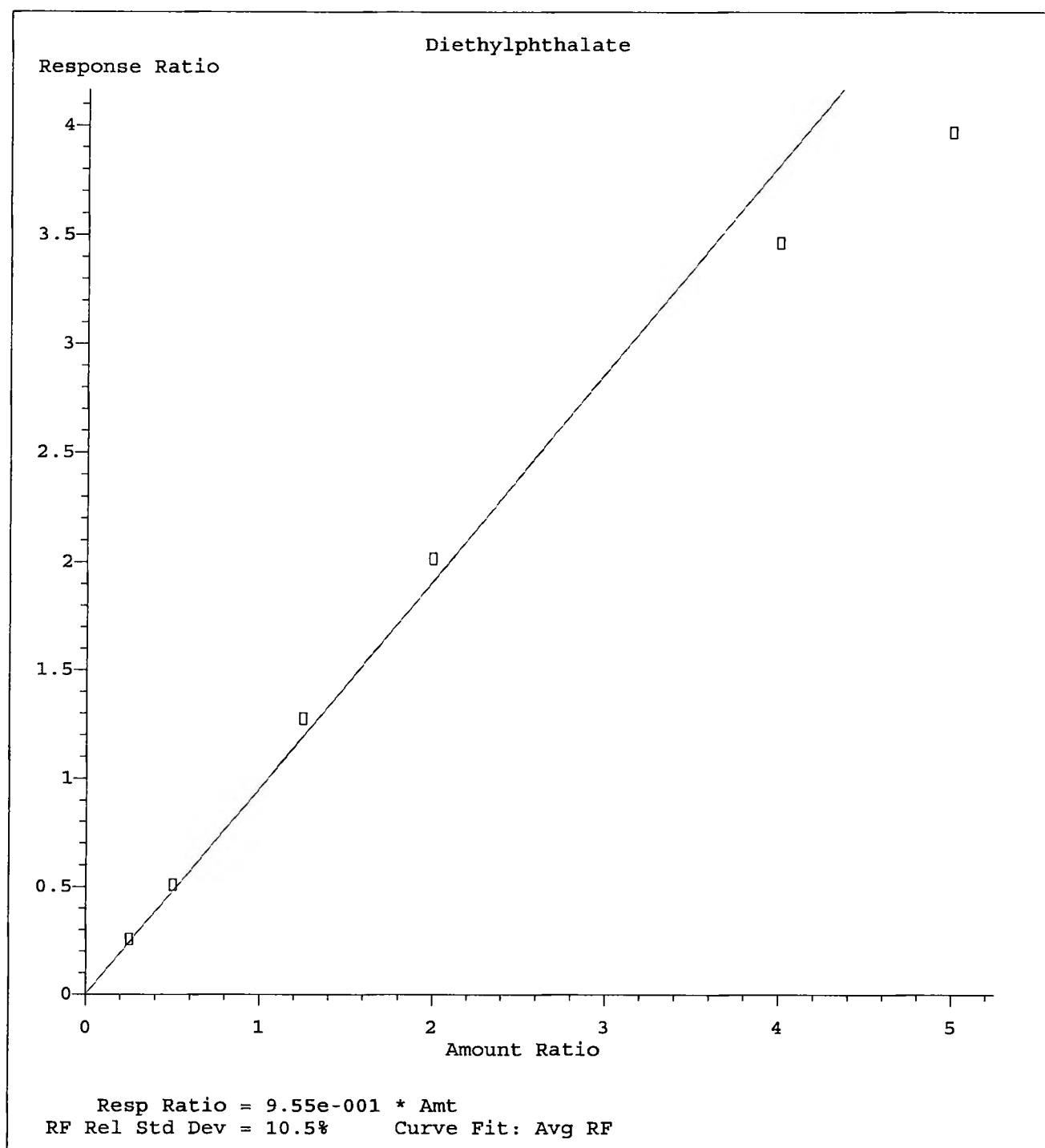
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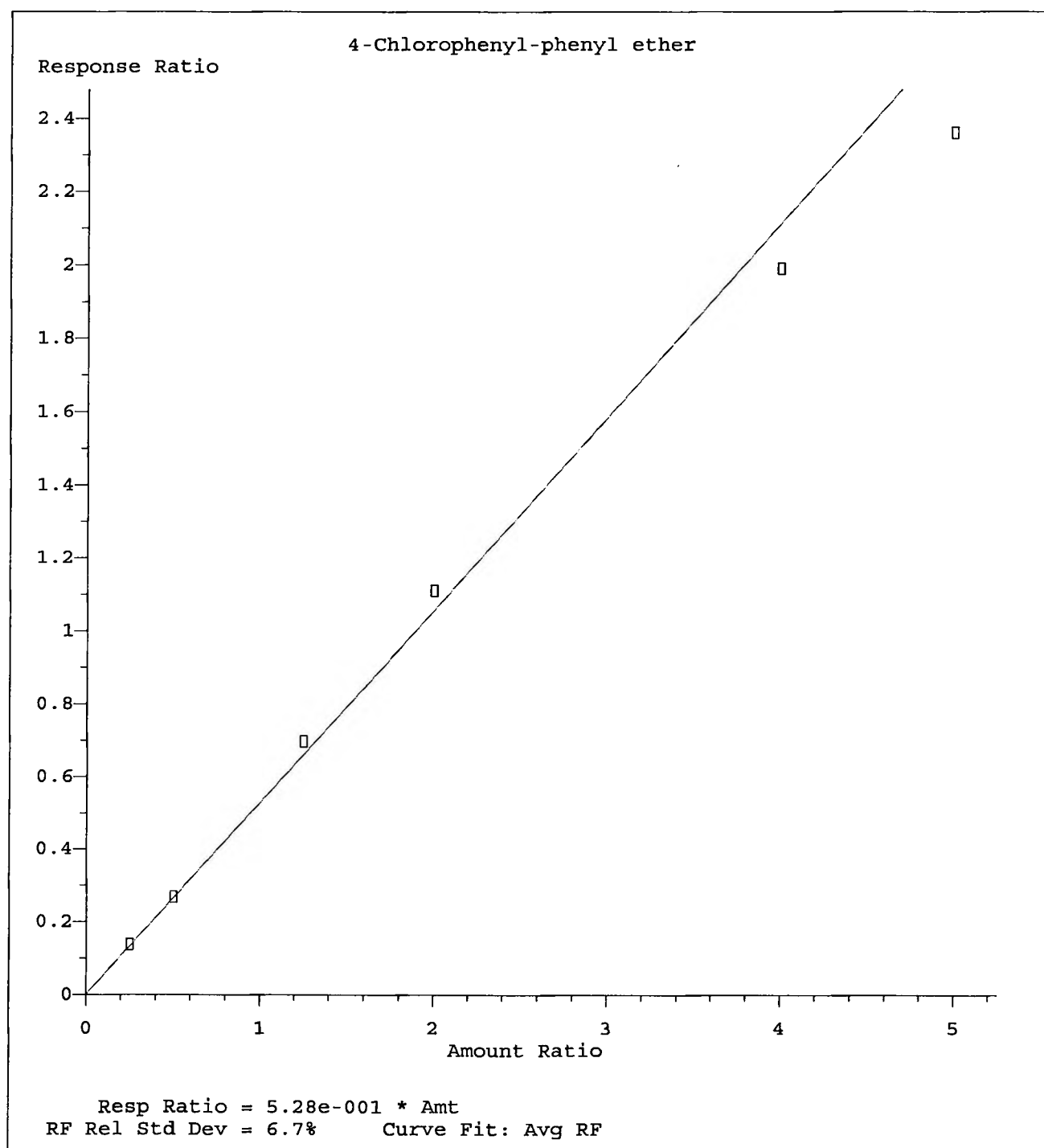
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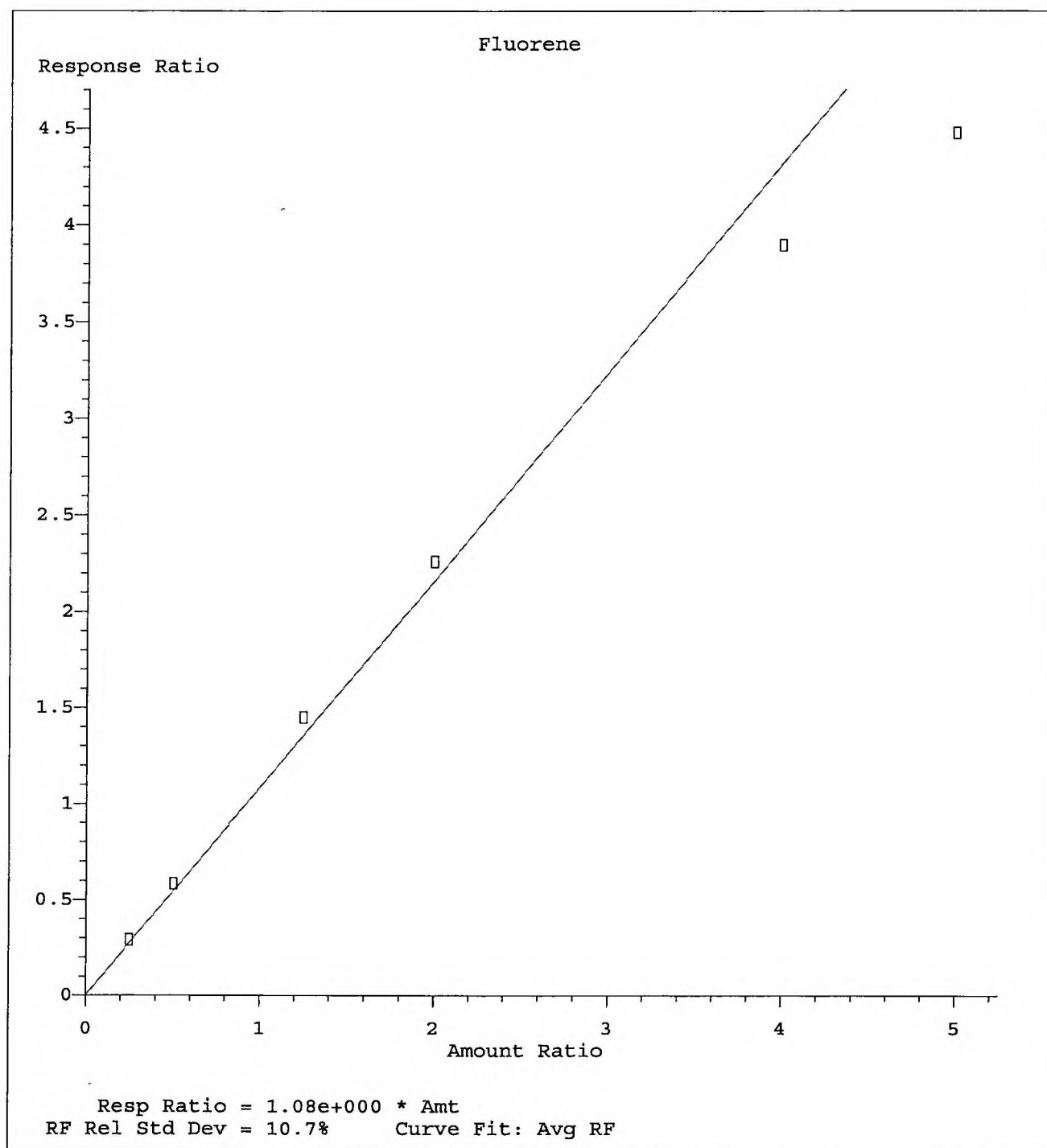
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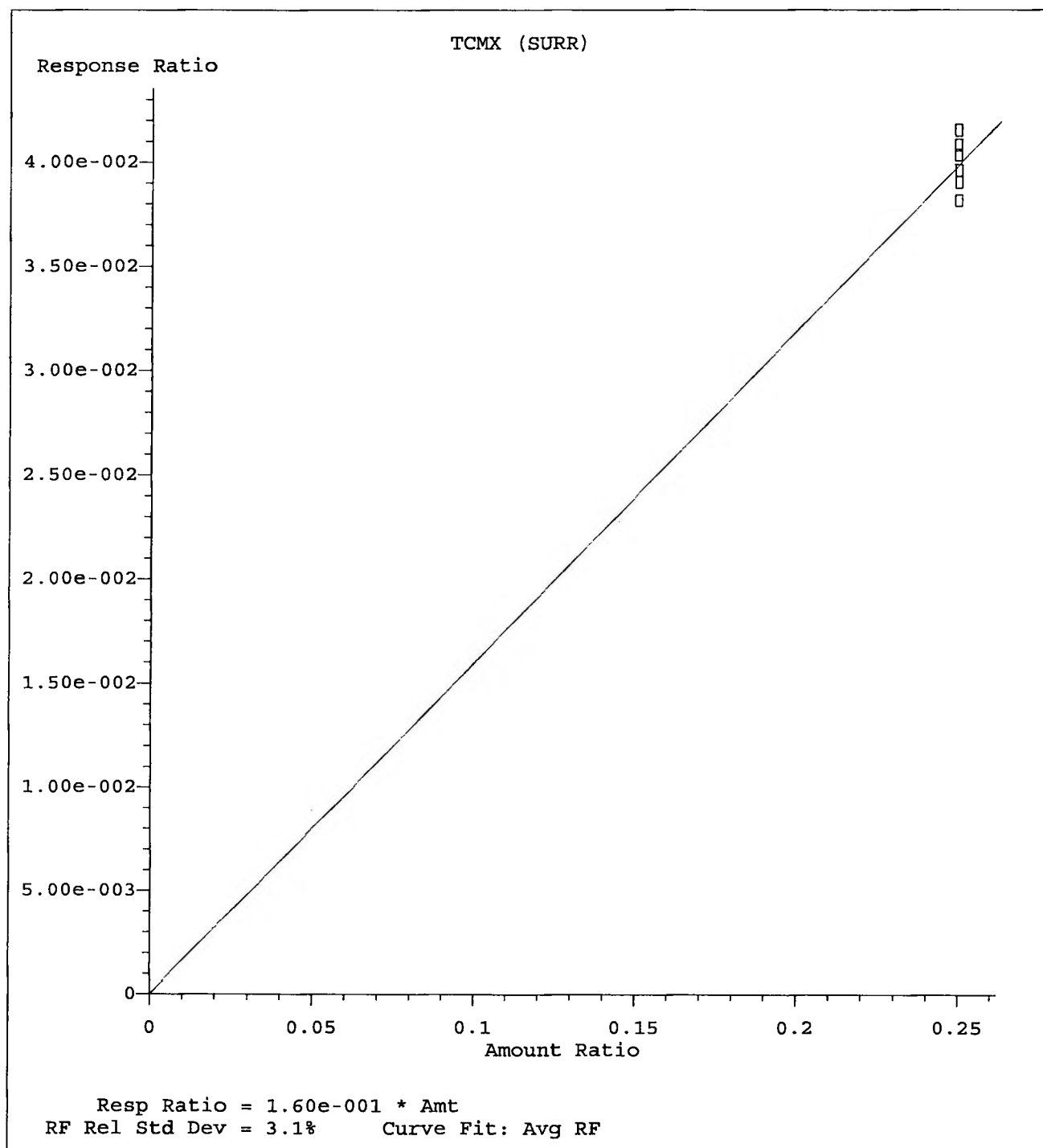
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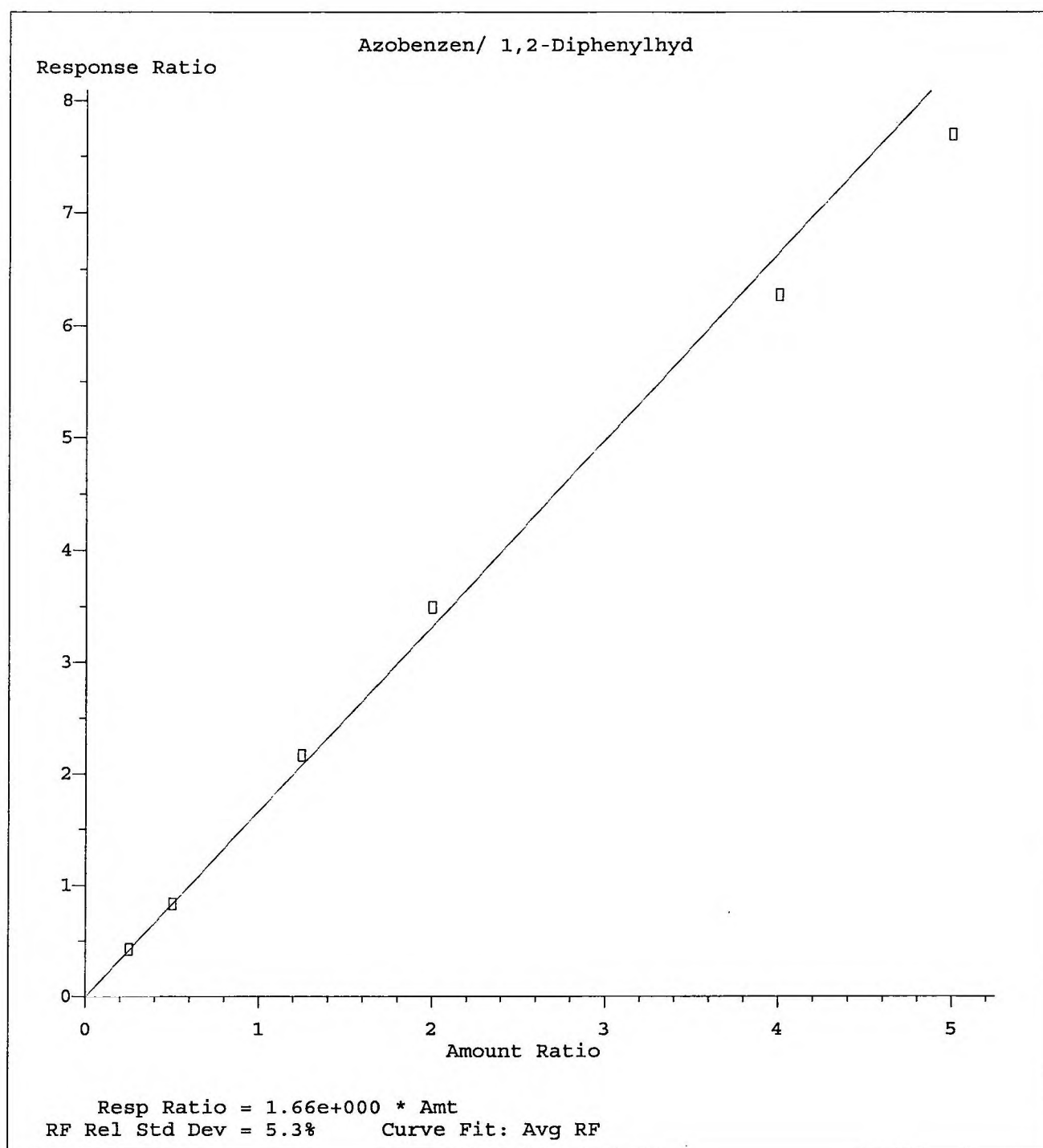
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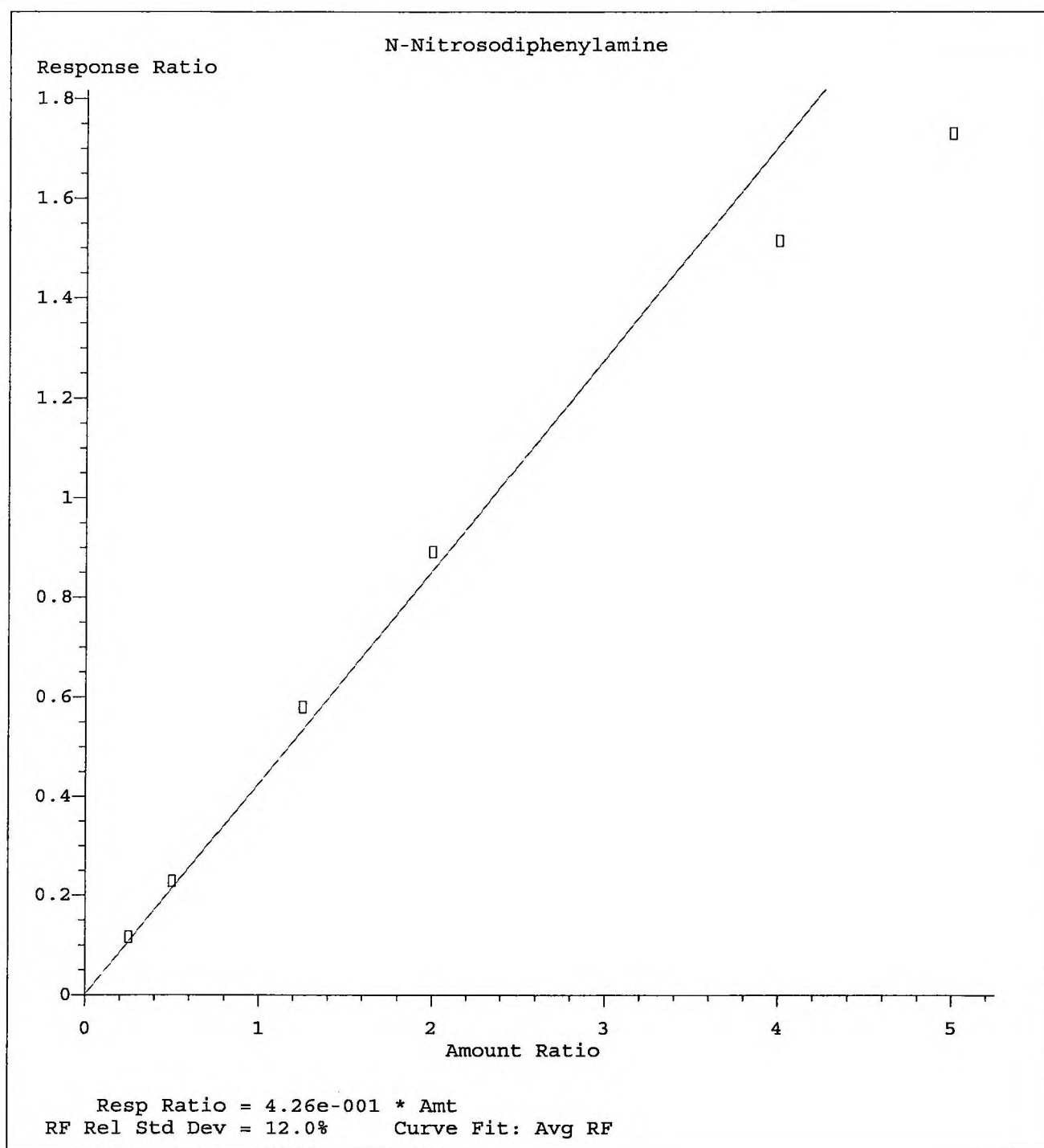
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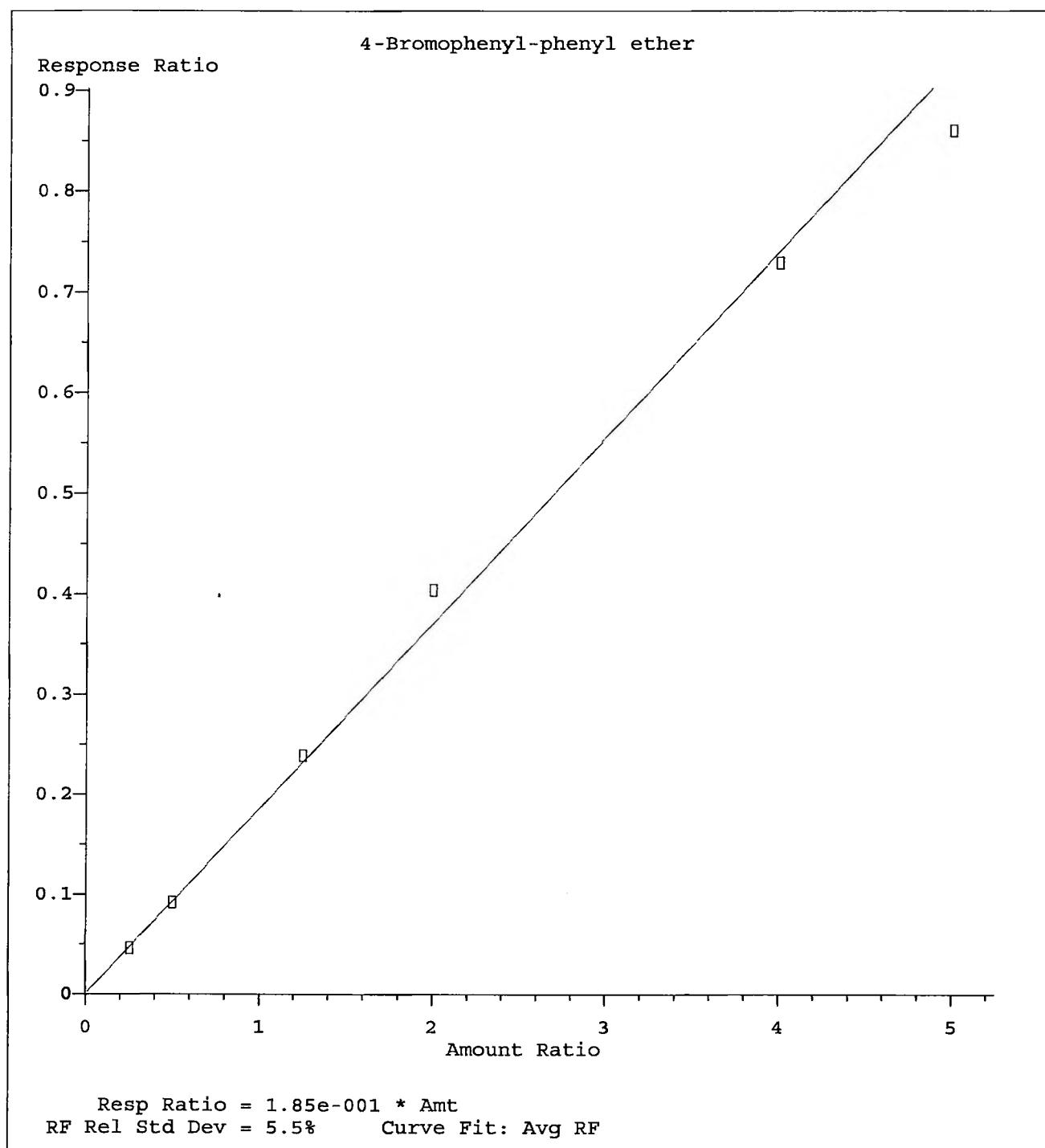
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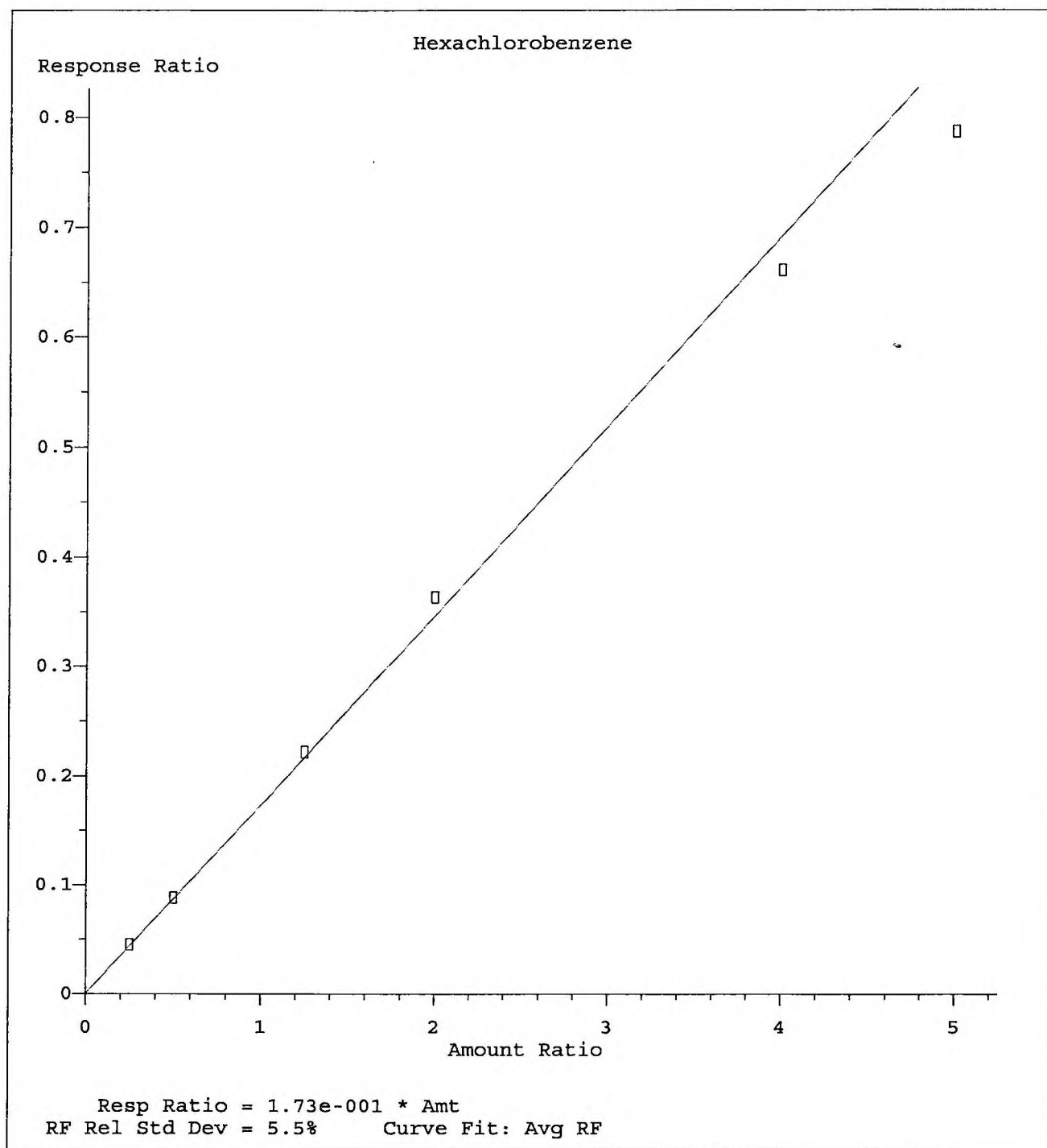
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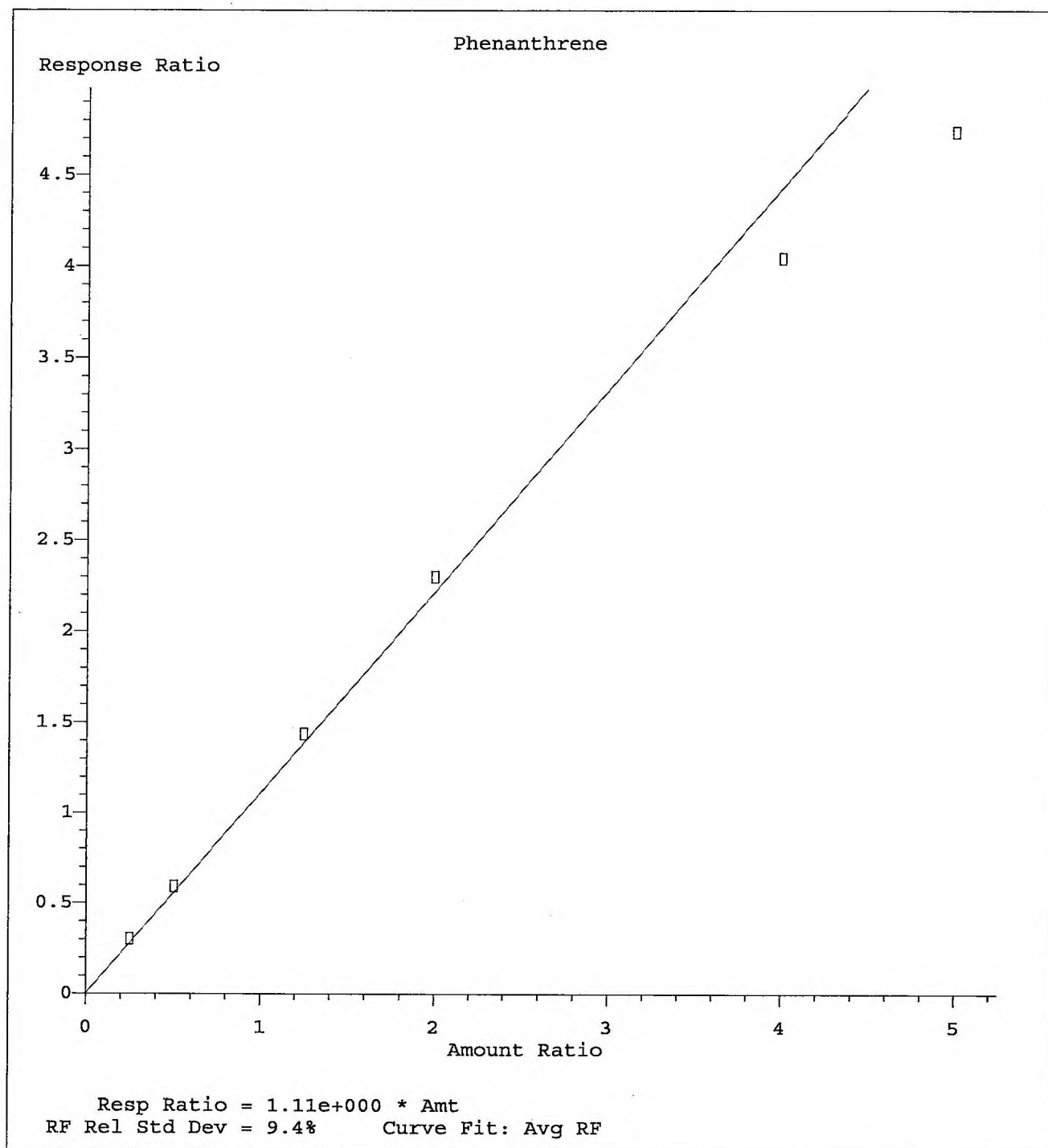
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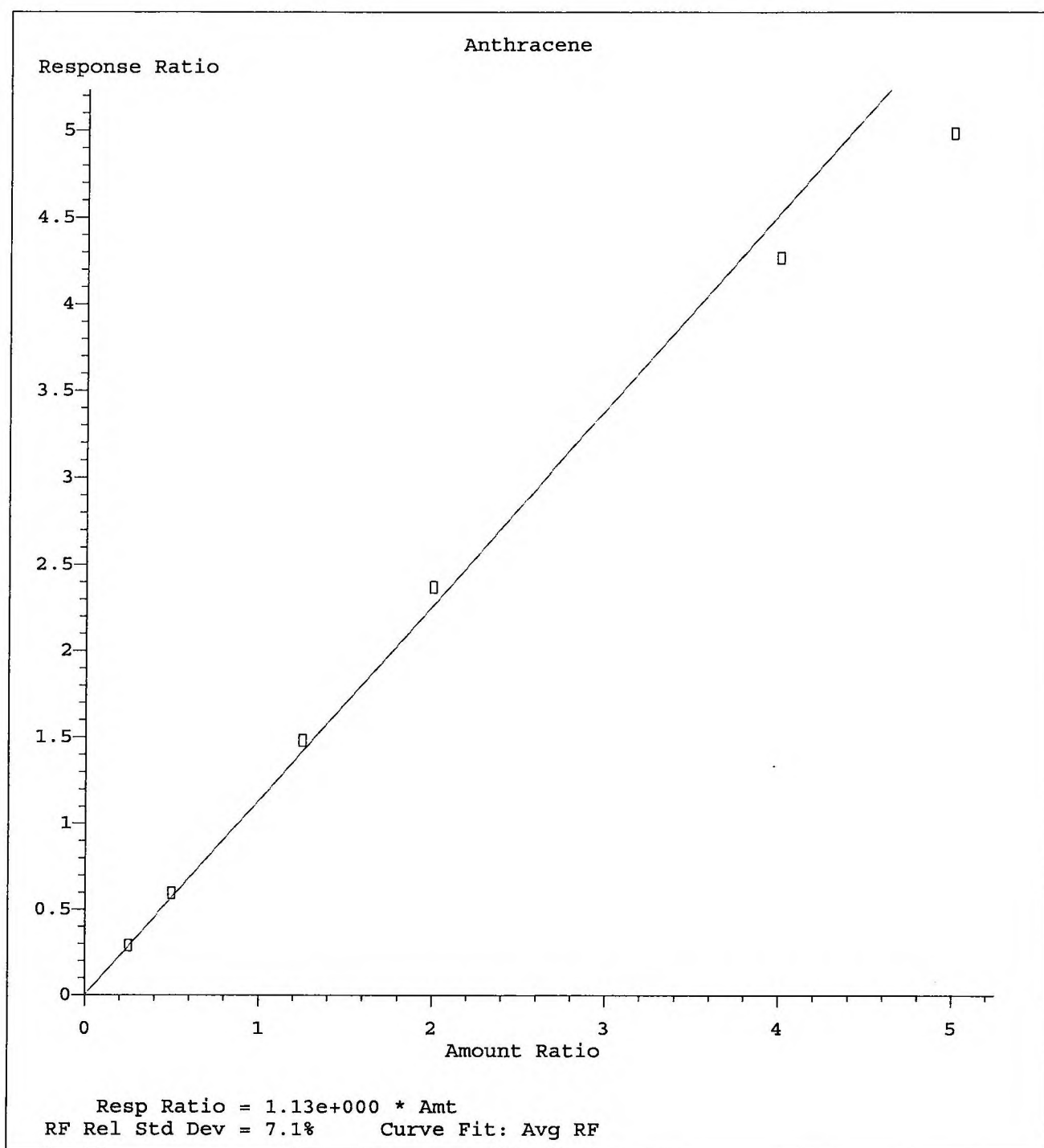
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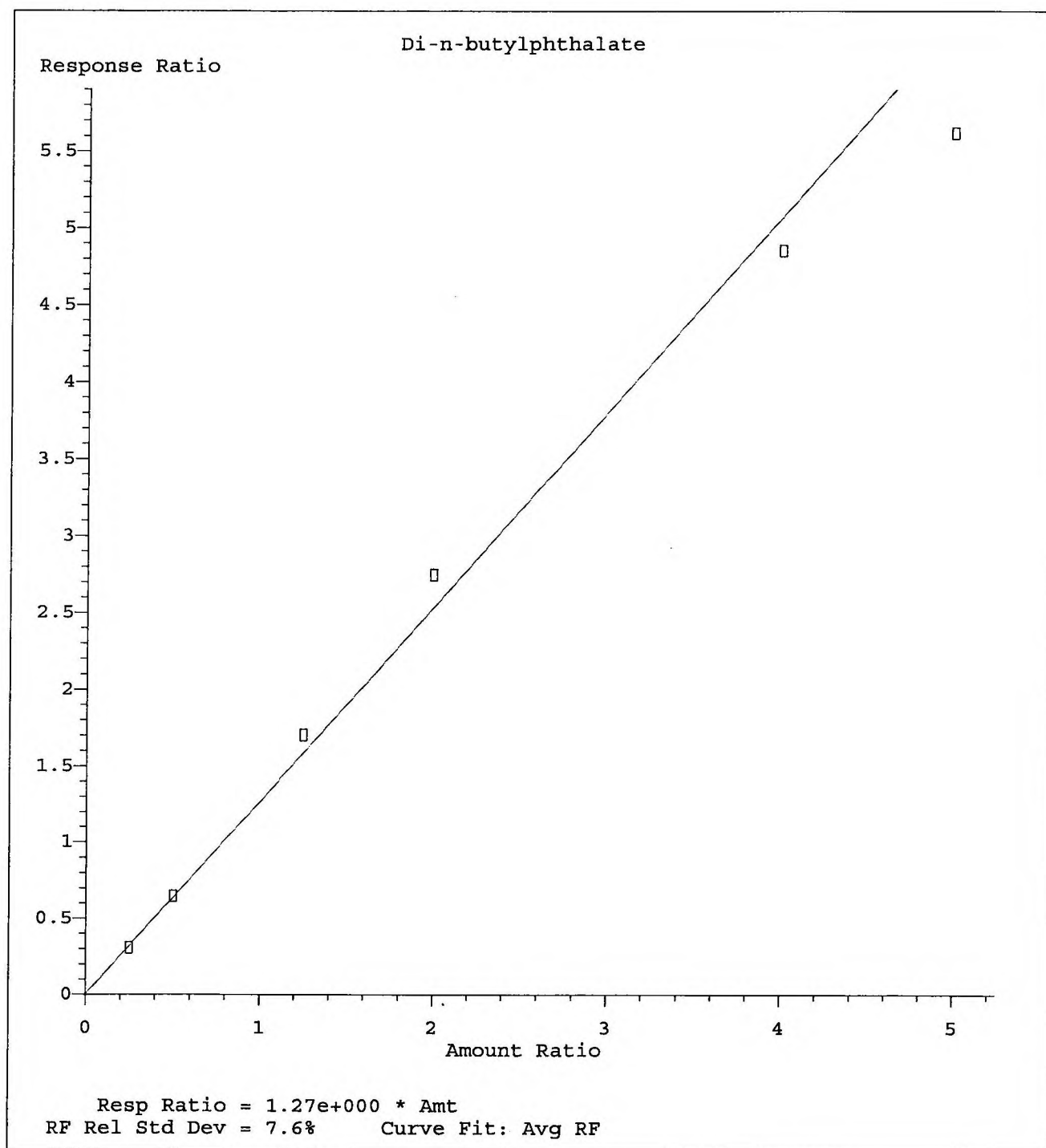
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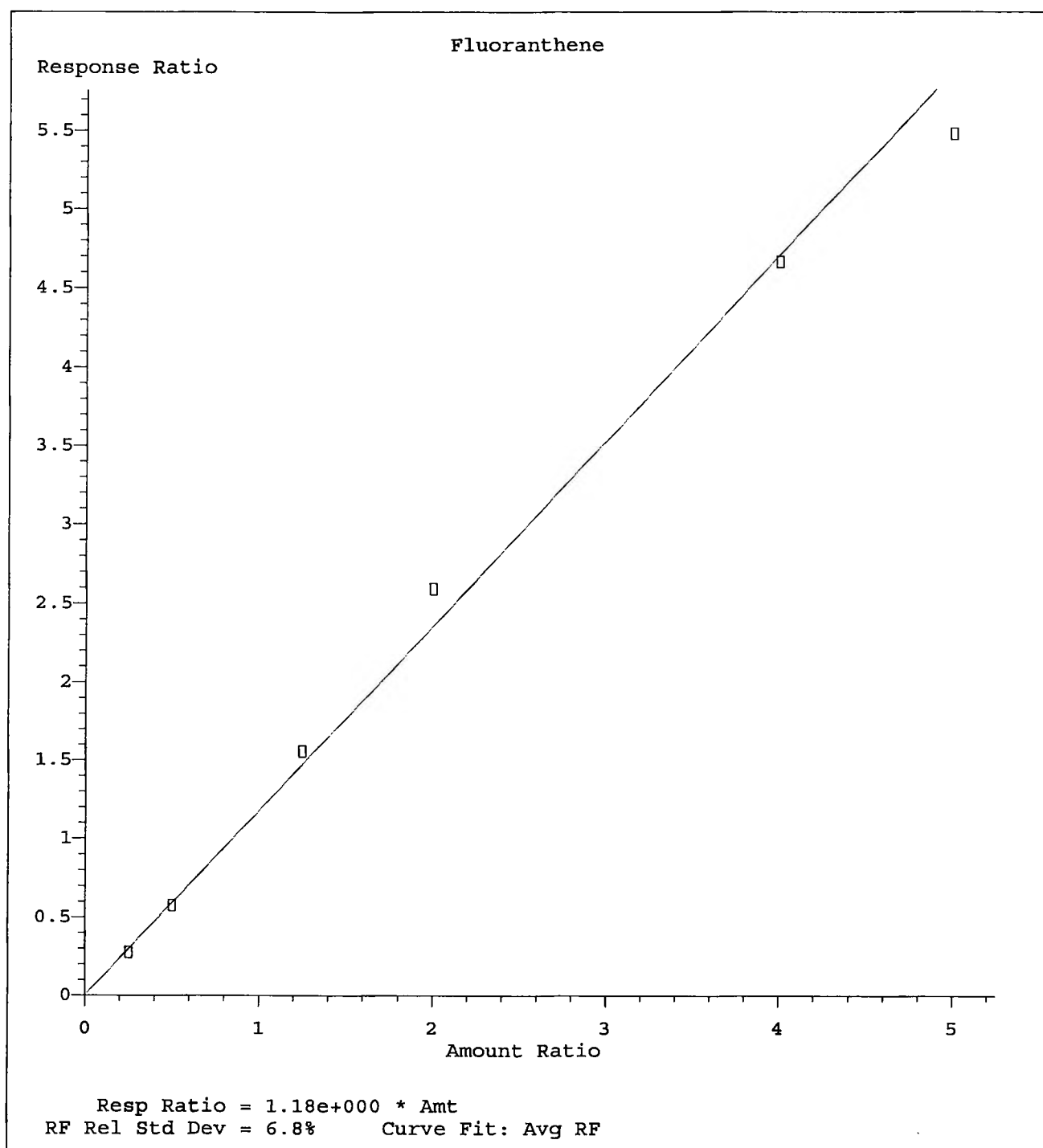
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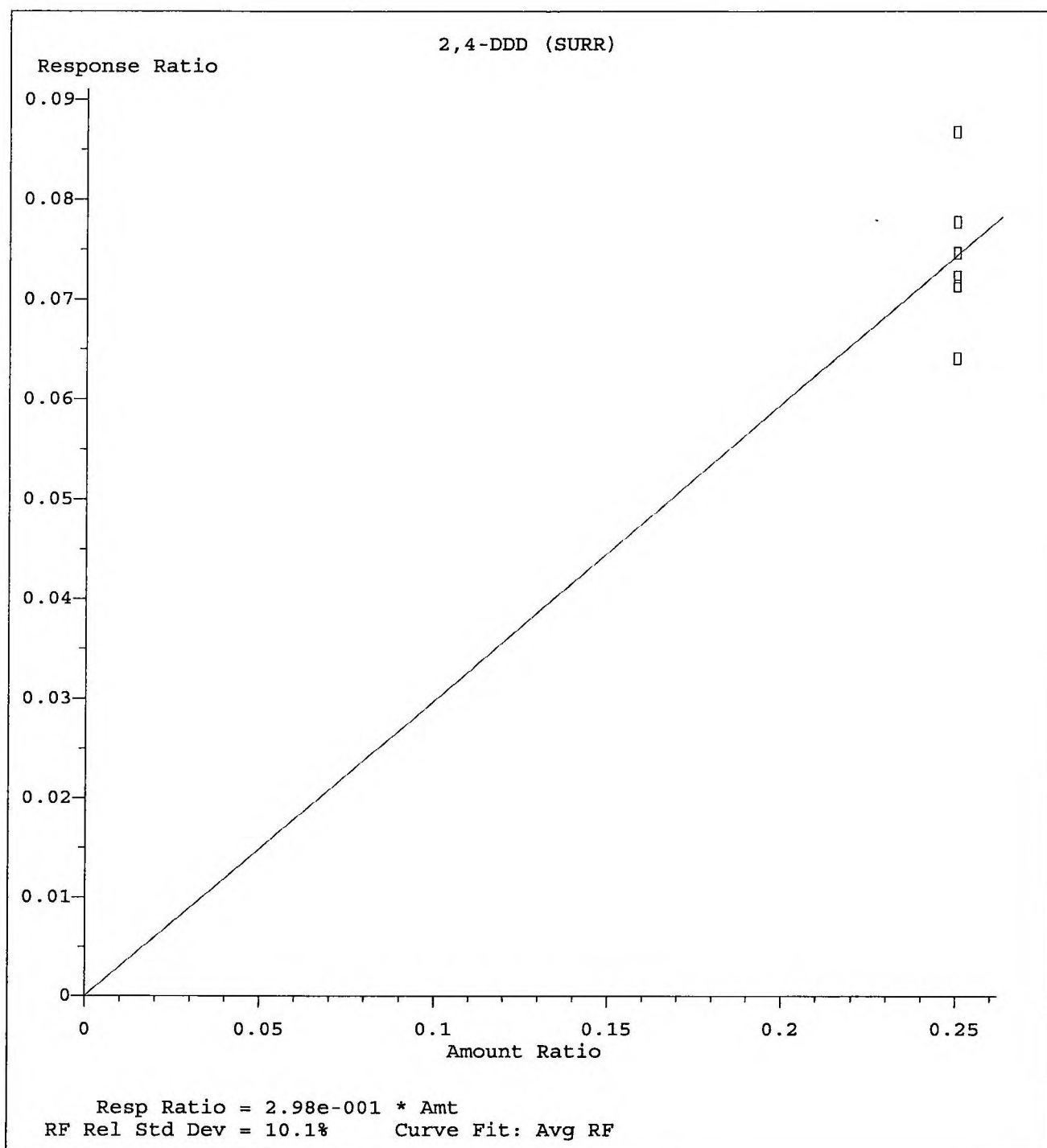
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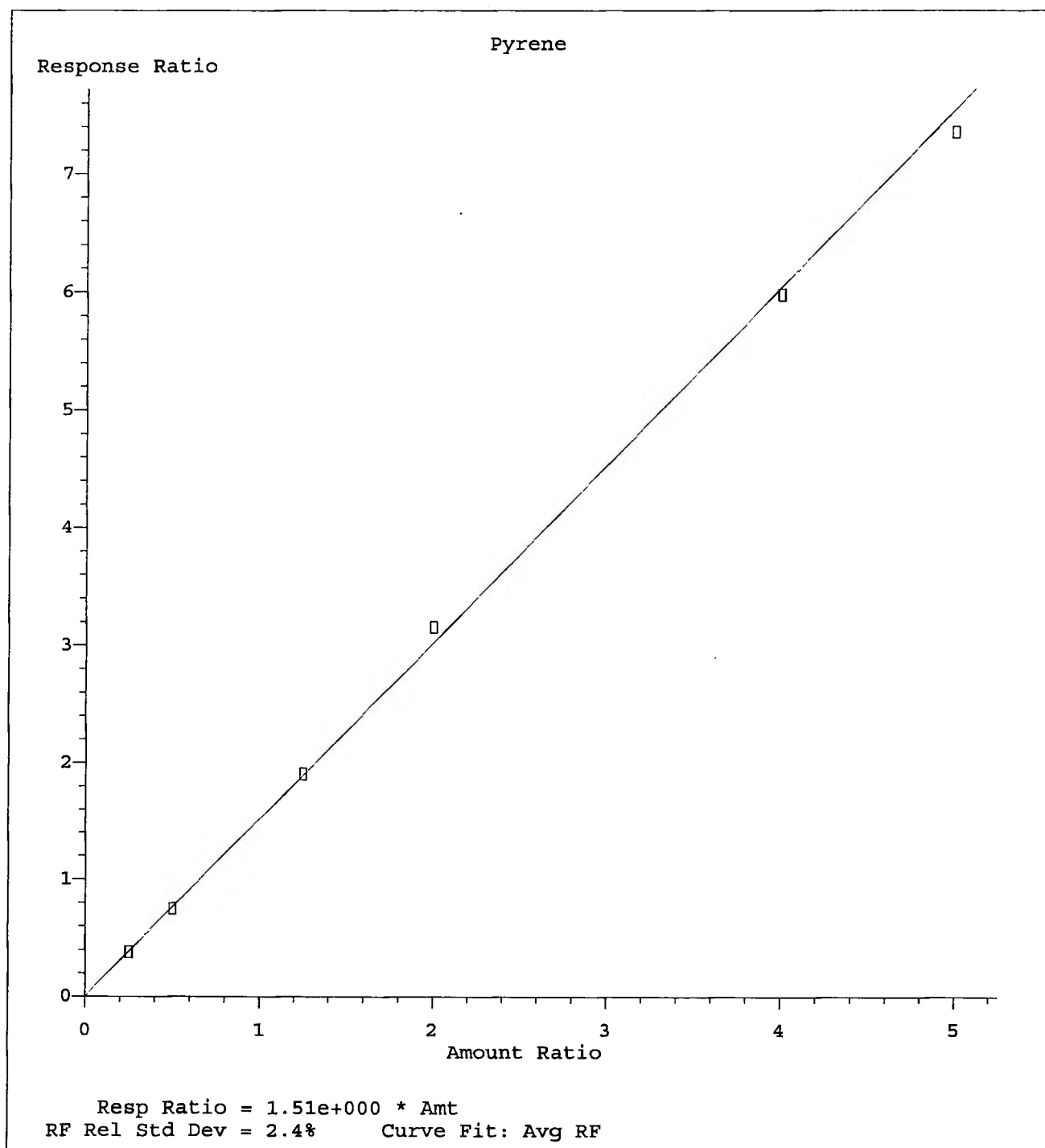
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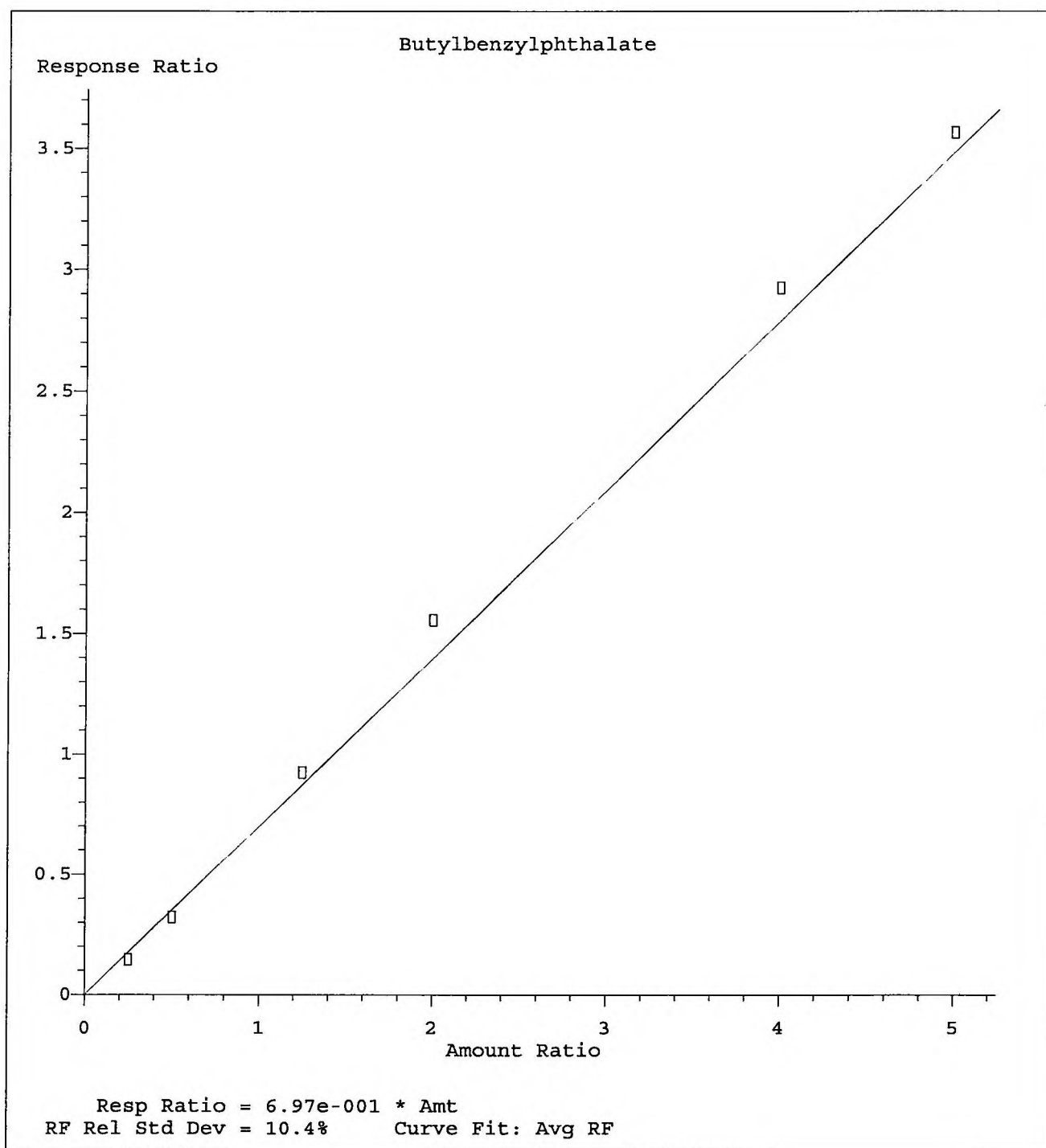
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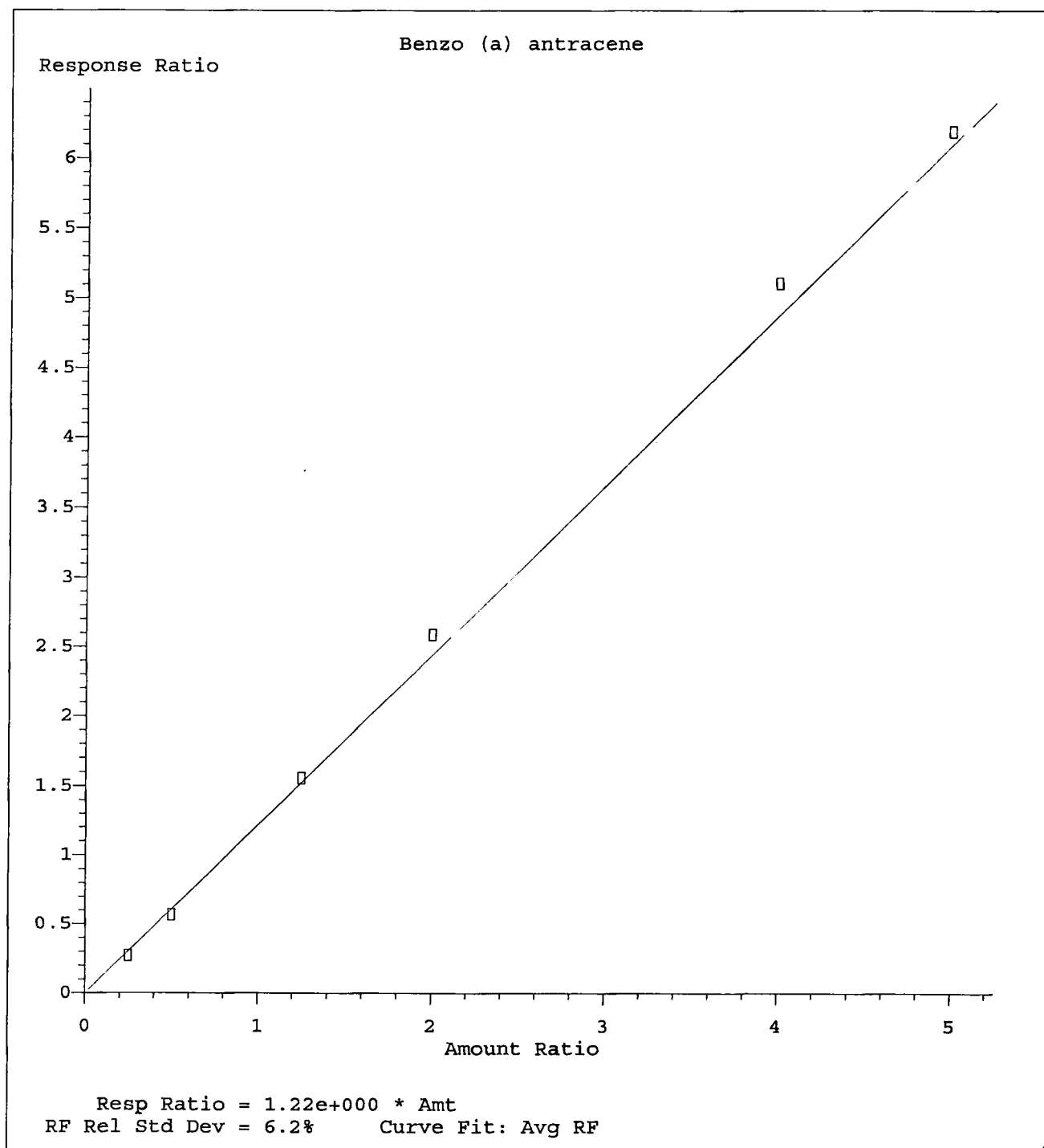
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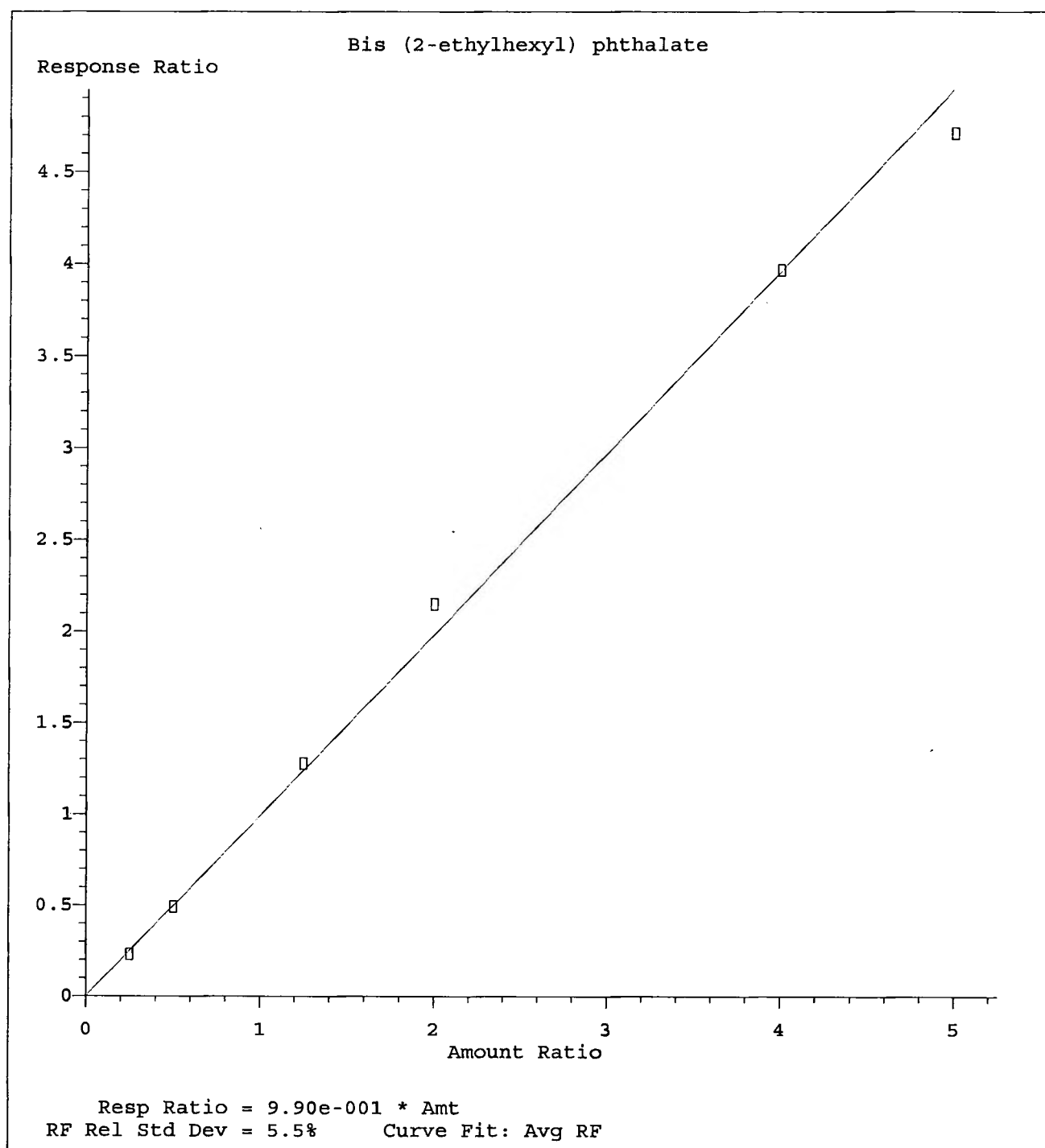
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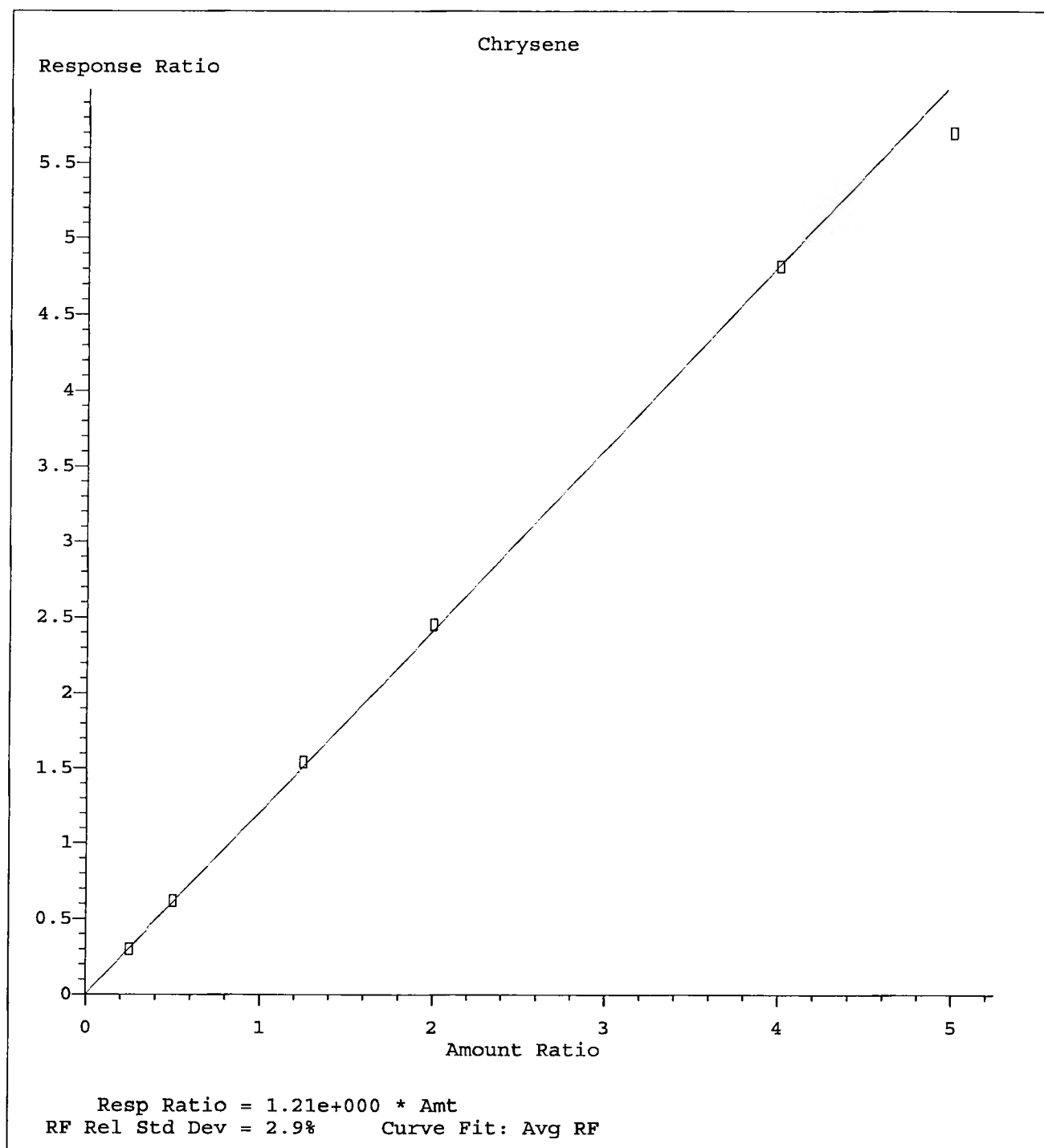
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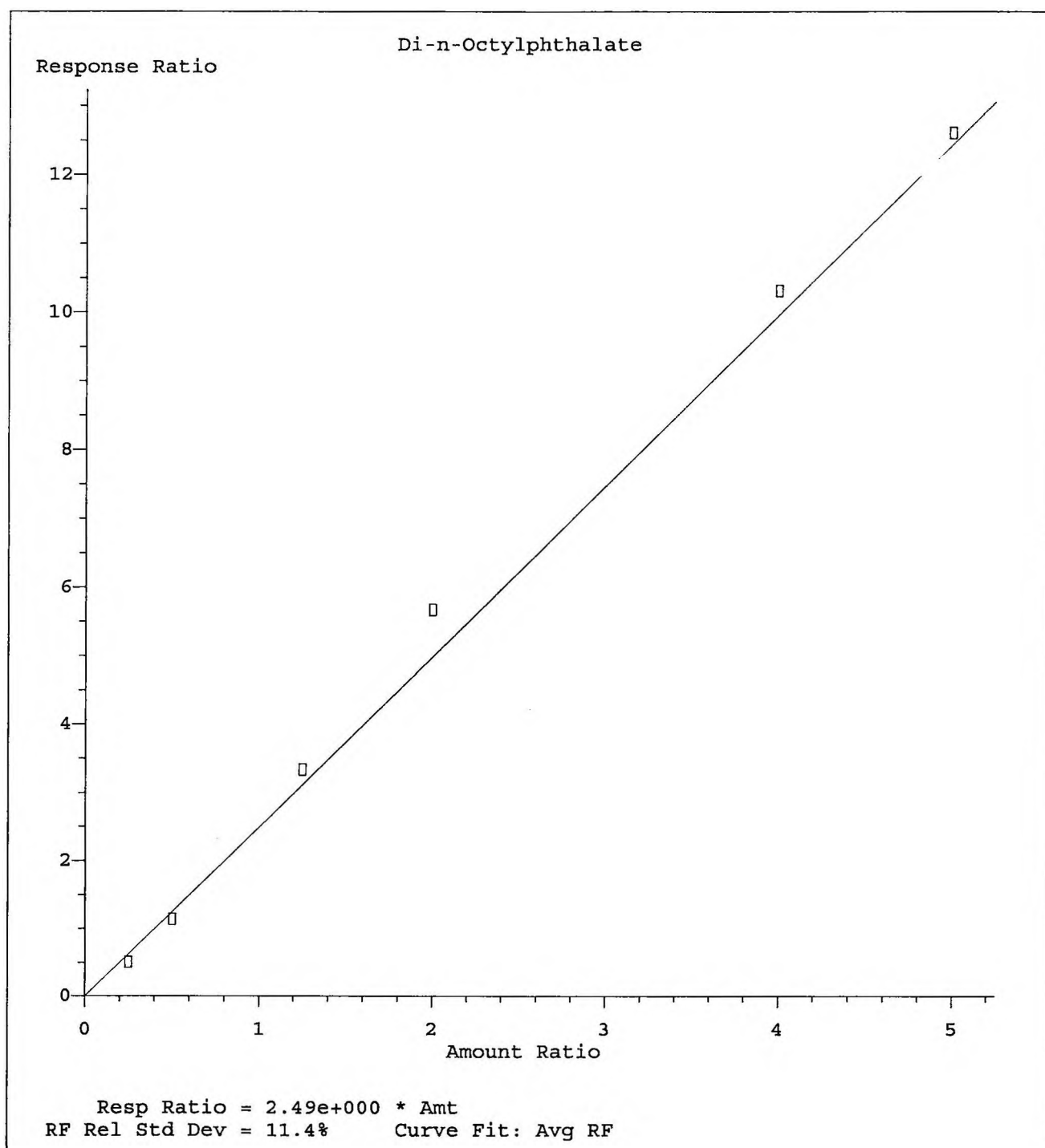
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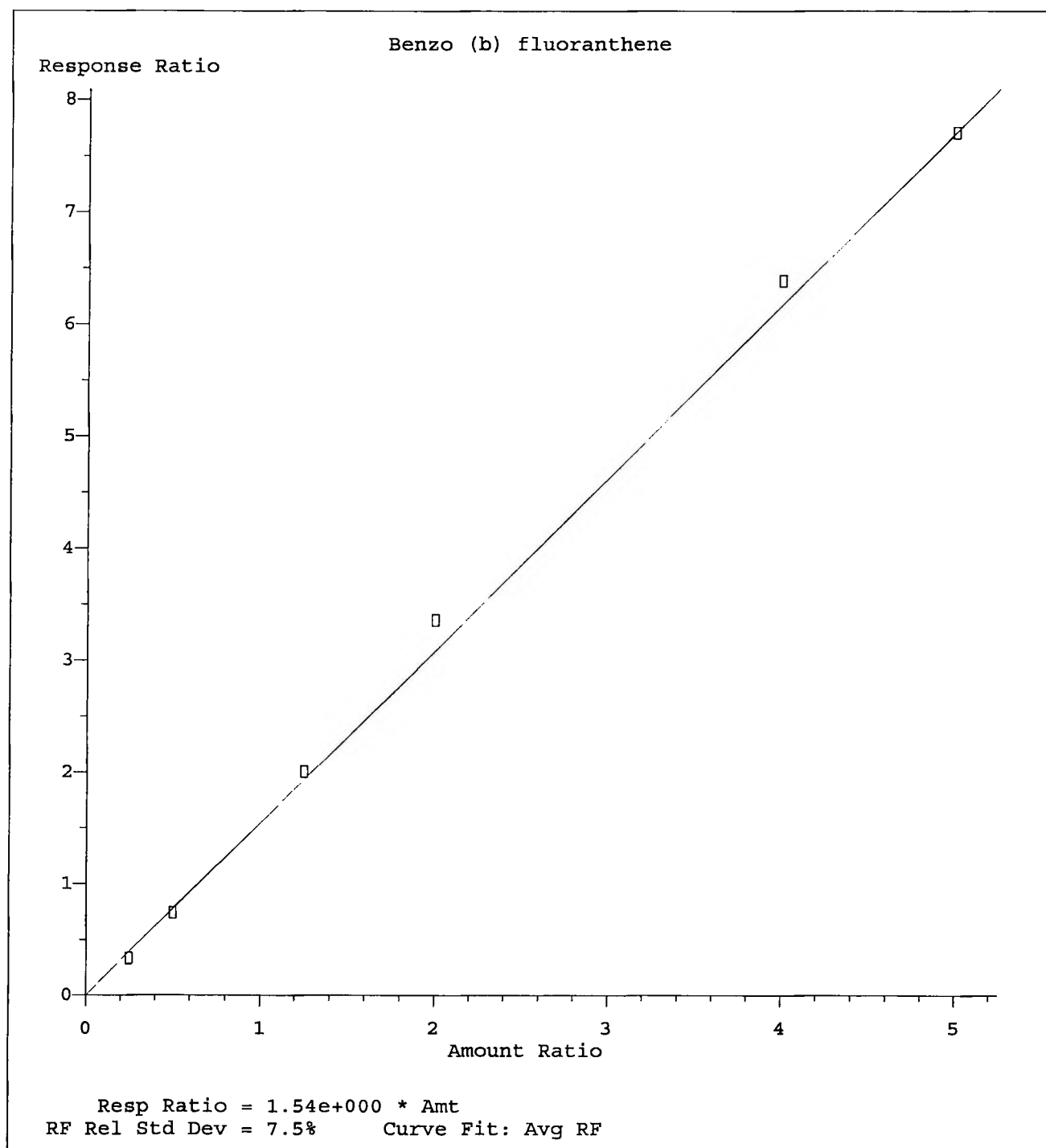
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Calibration Table Last Updated: Mon Apr 15 14:48:10 2002



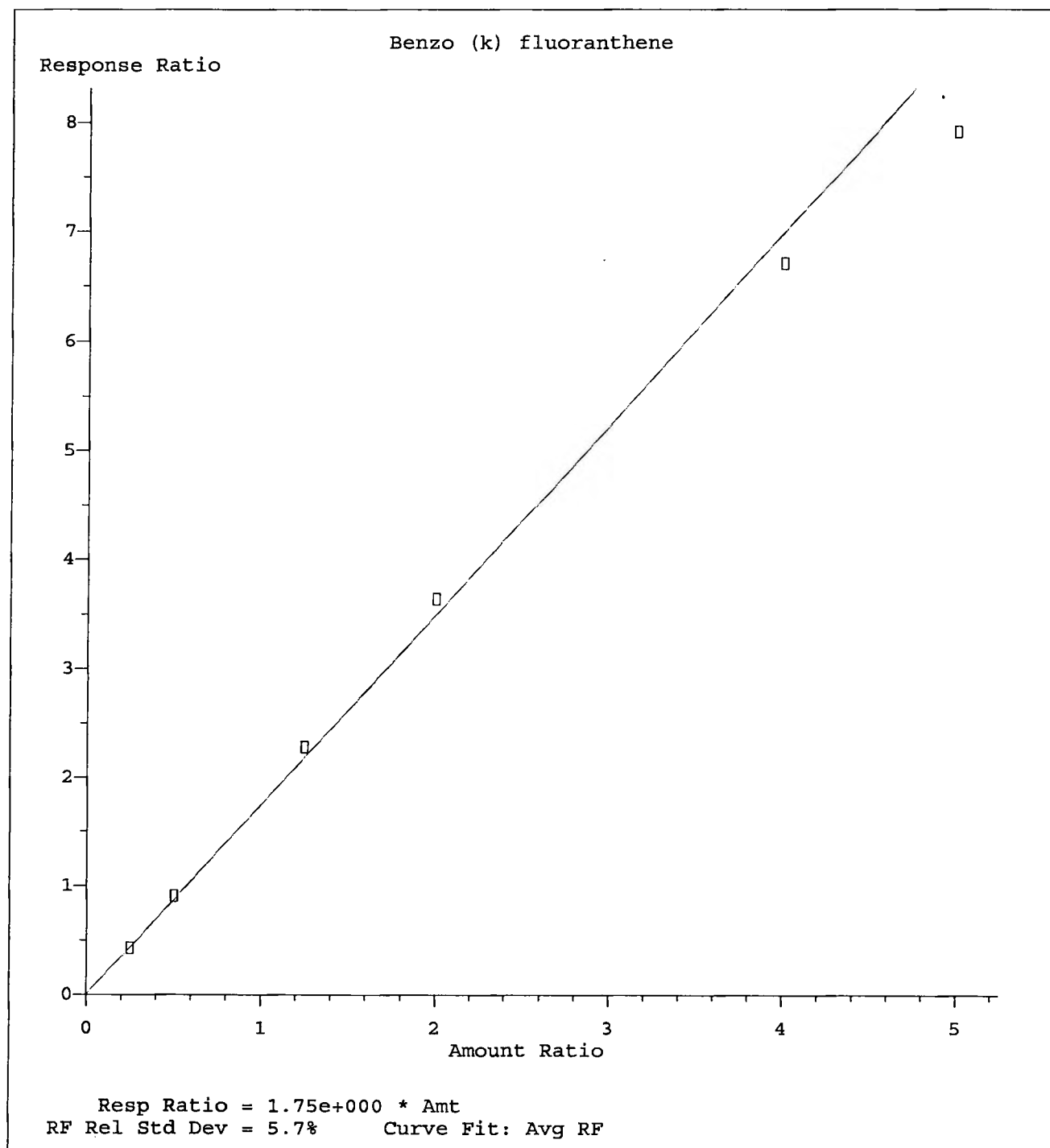
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 Calibration Table Last Updated: Mon Apr 15 14:48:10 2002



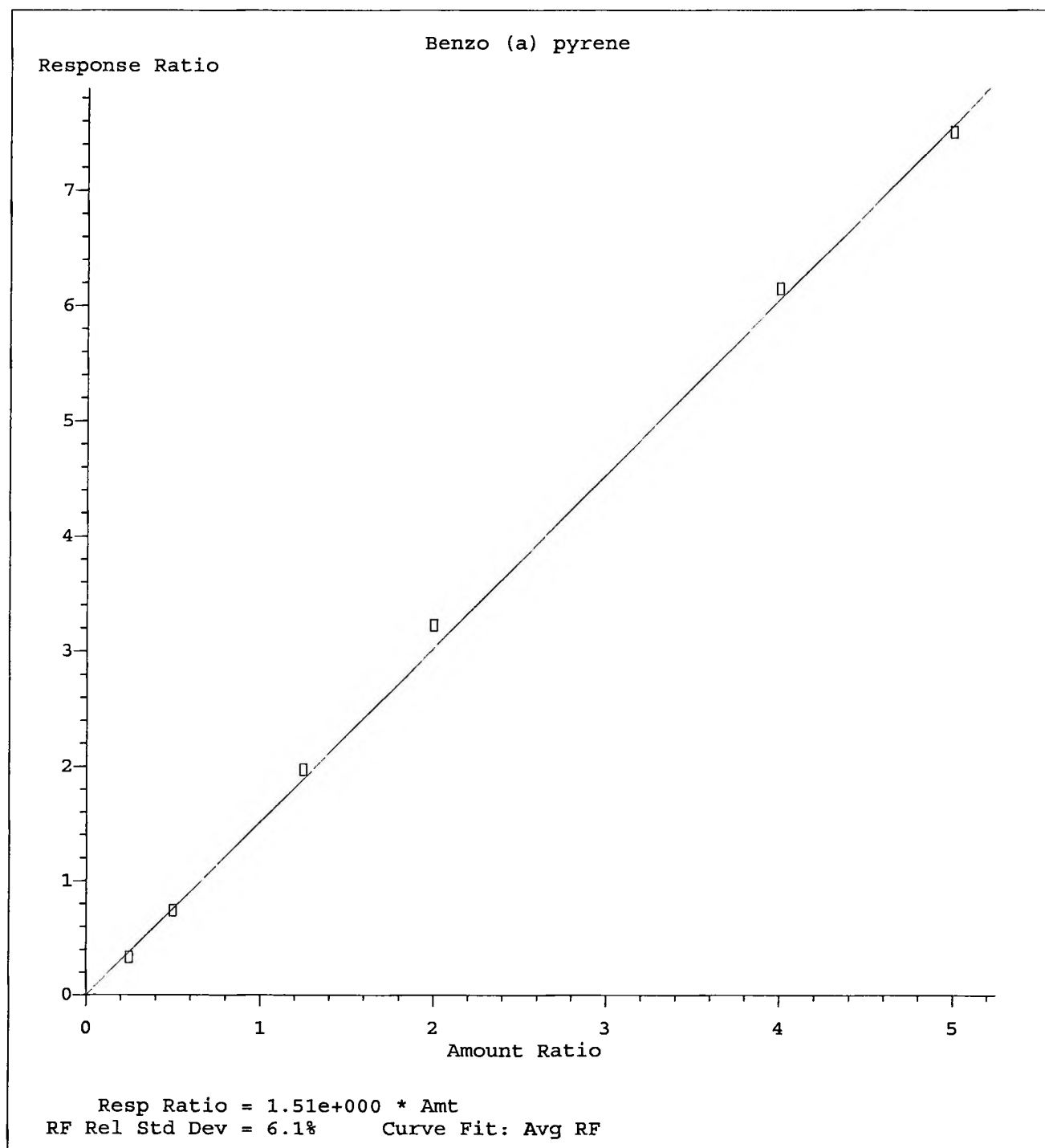
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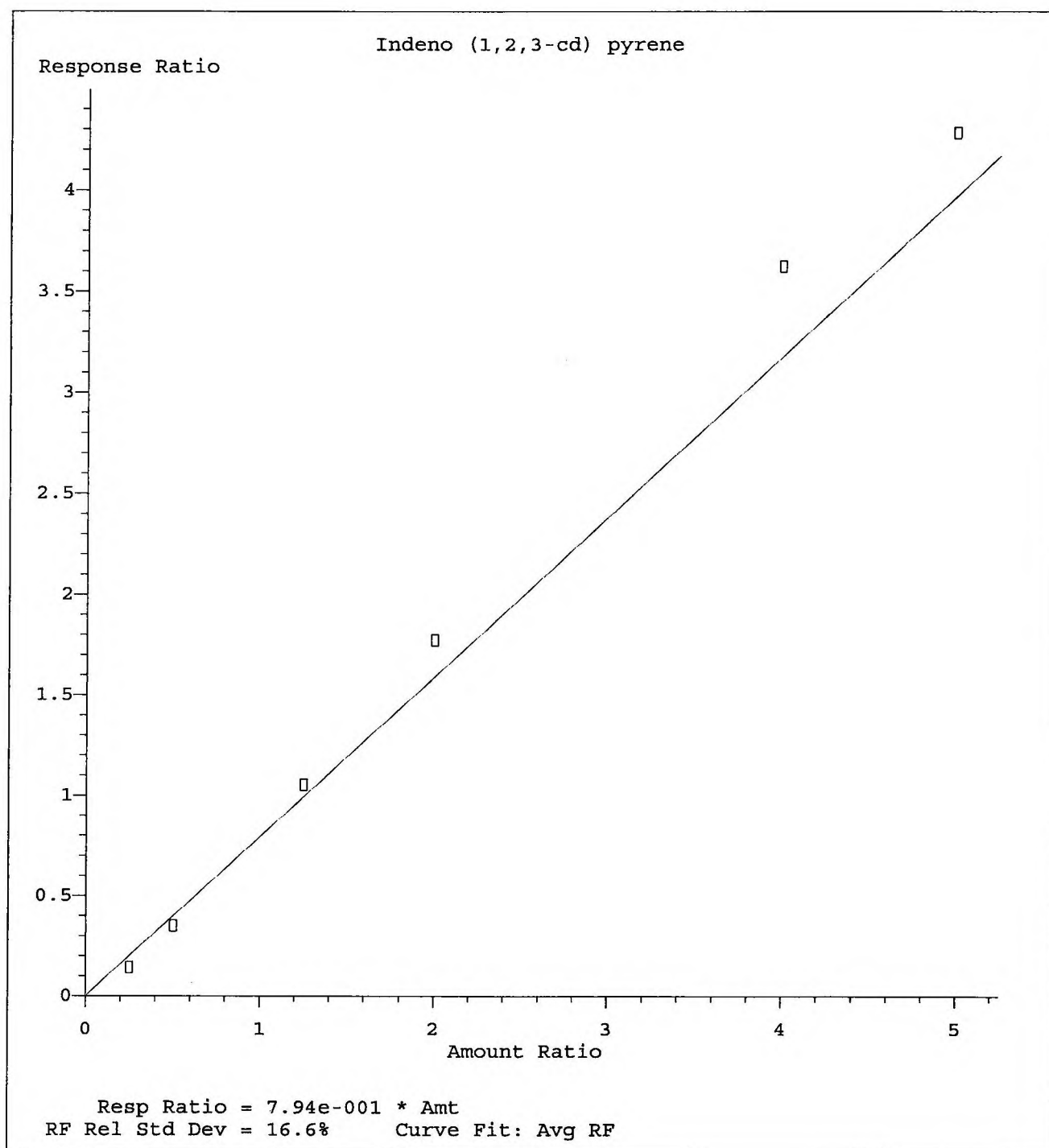
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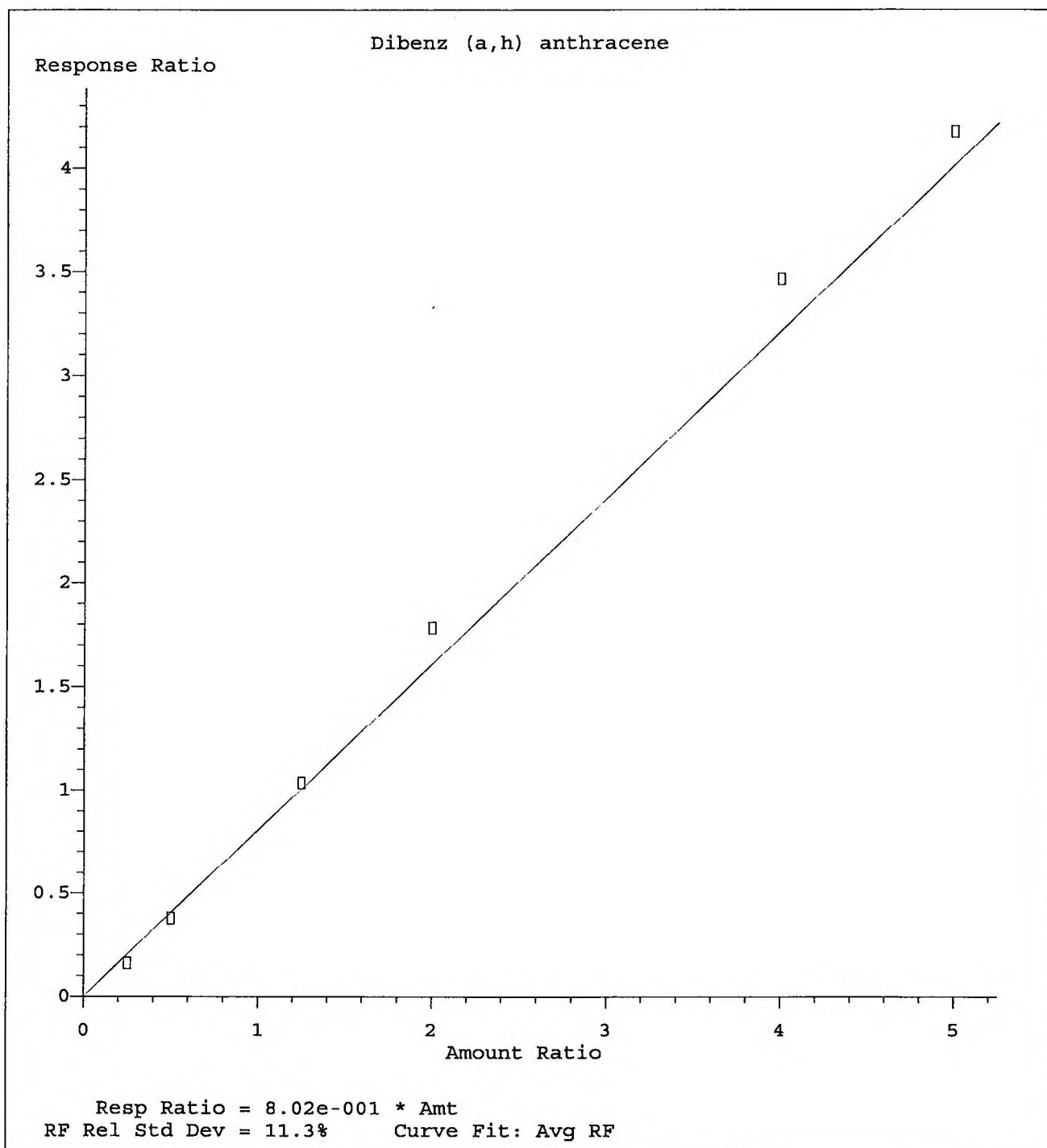
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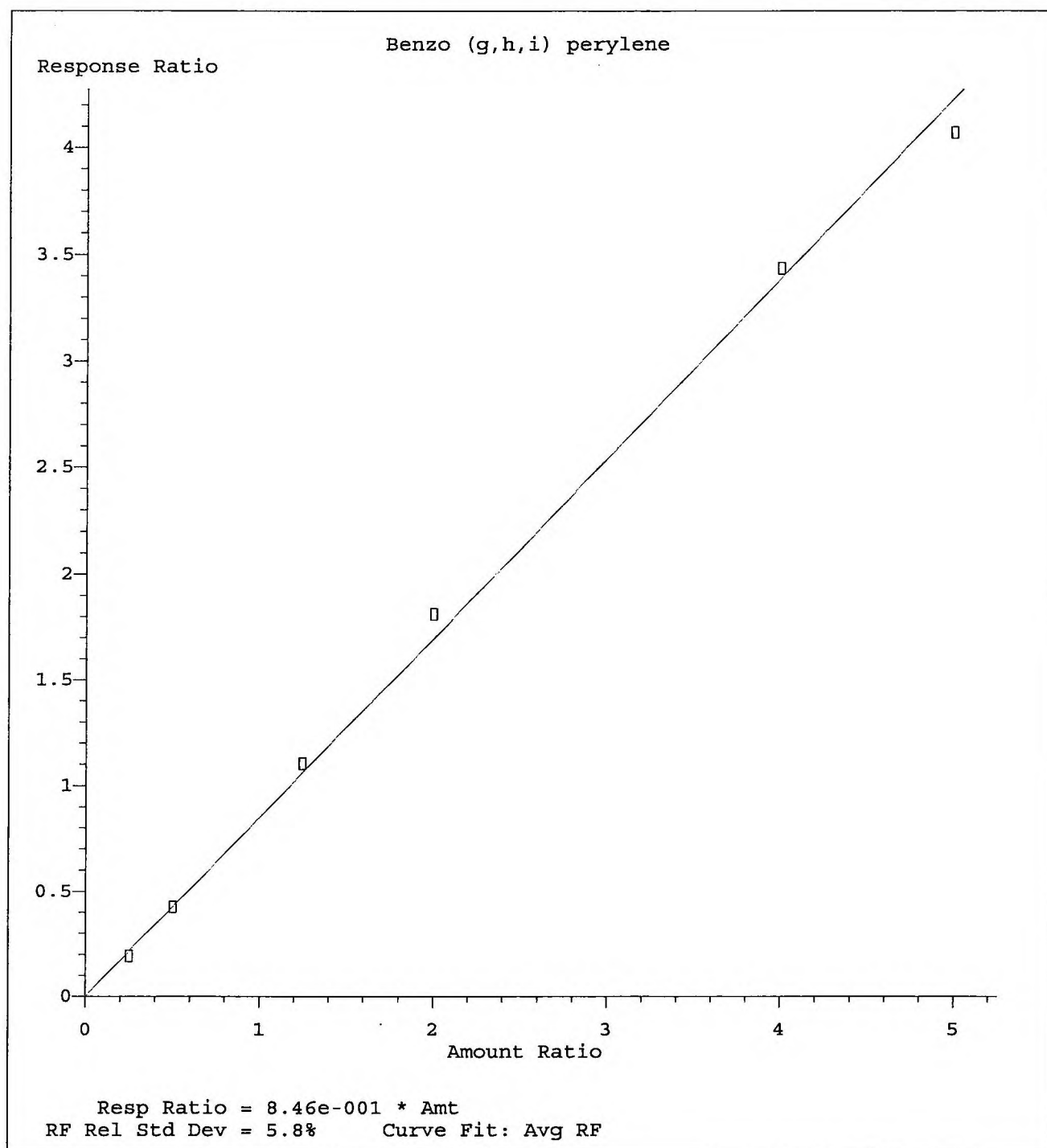
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