

Ionic Liquid Phase Organic Synthesis. Preparation of a Parallel Array of Isoxazolines

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Supporting Information

Spectra of compounds **3(A1-D4)**

¹HNMR (200 MHz, CDCl₃) ¹³C NMR (50 MHz, CDCl₃)

Compound 3A1	¹ HNMR	pag 2
Compound 3A2	¹ HNMR	pag 3
	¹³ CNMR	pag 4
Compound 3A3	¹ HNMR	pag 5
	¹³ CNMR	pag 6
Compound 3A4	¹ HNMR	pag 7
Compound 3B1	¹ HNMR	pag 8
Compound 3B2	¹ HNMR	pag 9
Compound 3B3	¹ HNMR	pag 10
Compound 3B4	¹ HNMR	pag 11
Compound 3C1	¹ HNMR	pag 12
Compound 3C2	¹ HNMR	pag 13
	¹³ CNMR	pag 14
Compound 3C3	¹ HNMR	pag 15
Compound 3C4	¹ HNMR	pag 16
Compound 3D1	¹ HNMR	pag 17
	¹³ CNMR	pag 18
Compound 3D2	¹ HNMR	pag 19
	¹³ CNMR	pag 20
Compound 3D3	¹ HNMR	pag 21
	¹³ CNMR	pag 22
Compound 3D4	¹ HNMR	pag 23

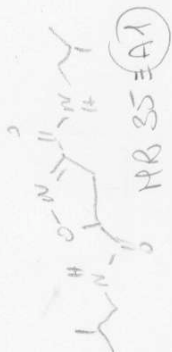
CC(C)NCC(=O)N1C(=O)C(=N1)C(=O)N(C)C

HR 35 = 41

2.68 1.00 1.63 2.93 0.57 5.18 28.64

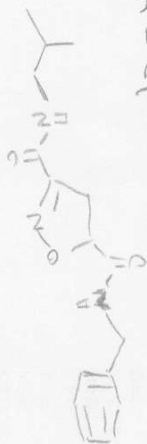
6.50 6.00 5.50 5.00 4.50 4.00 3.50 3.00 2.50 2.00 1.50 1.00 .50

PPM

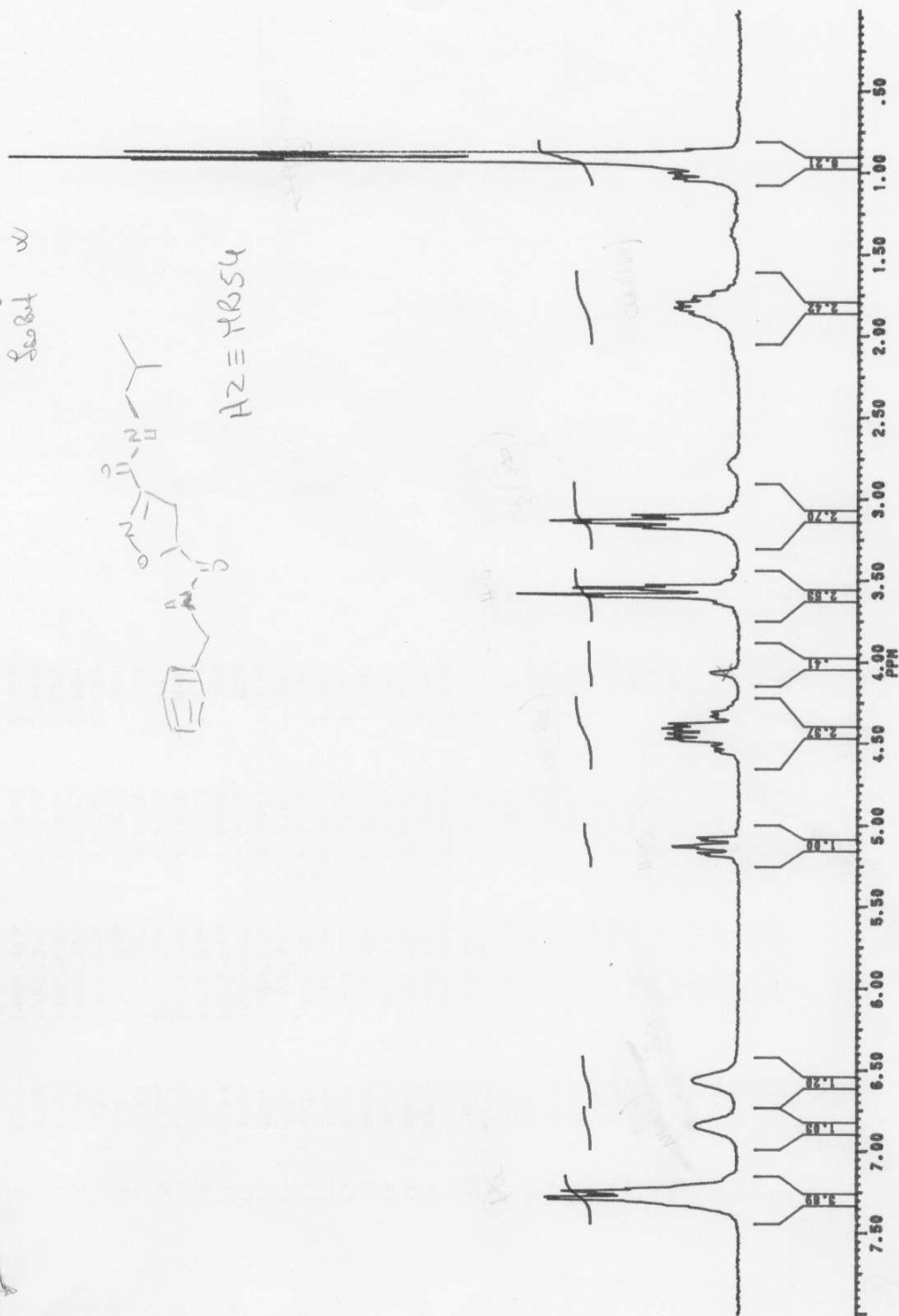


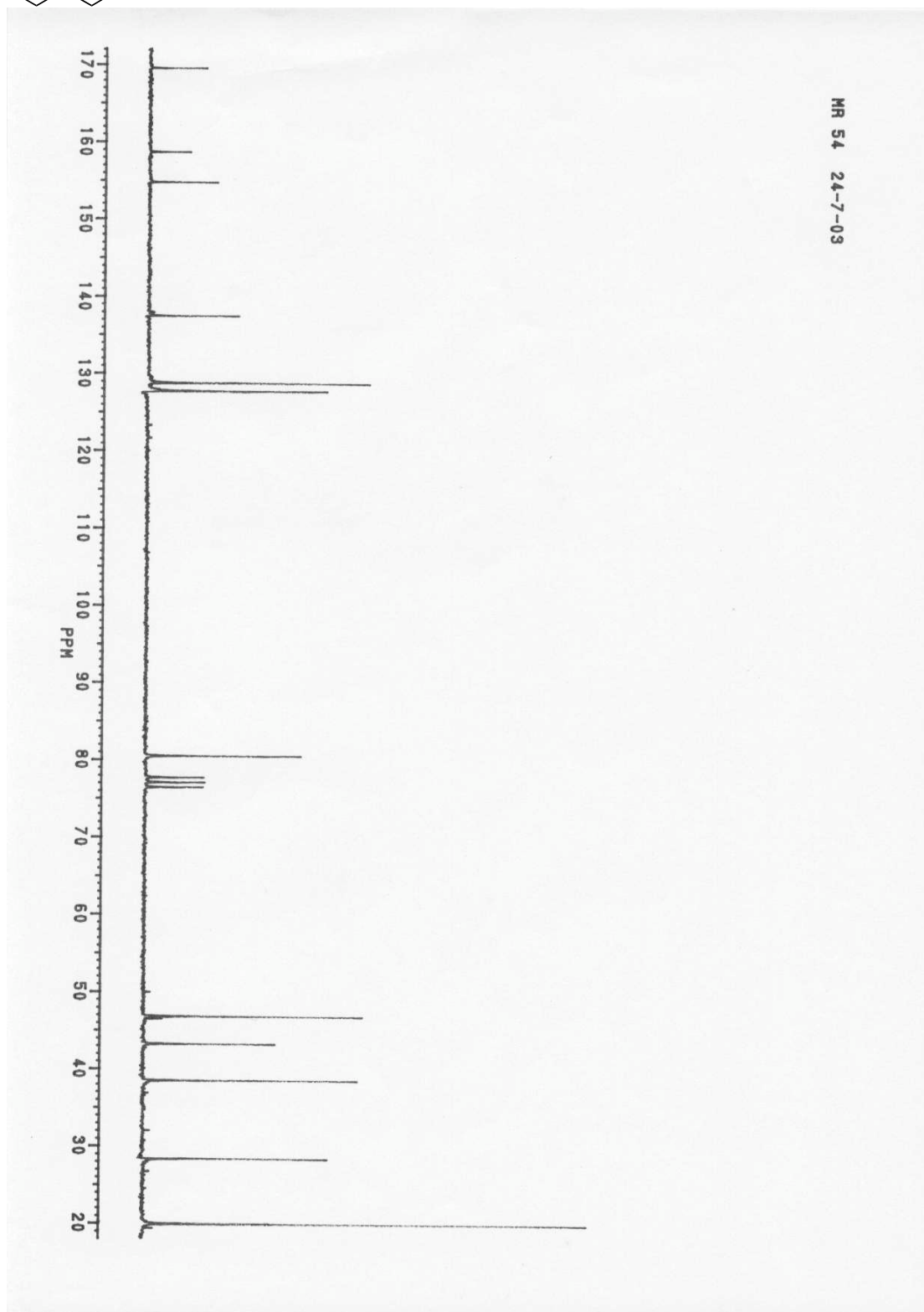
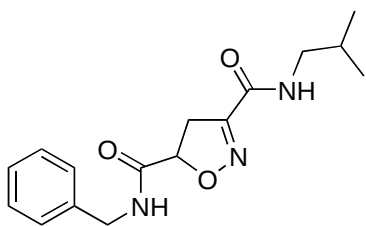
A2

BuNH₂ 5.83
 Sec But 4.0

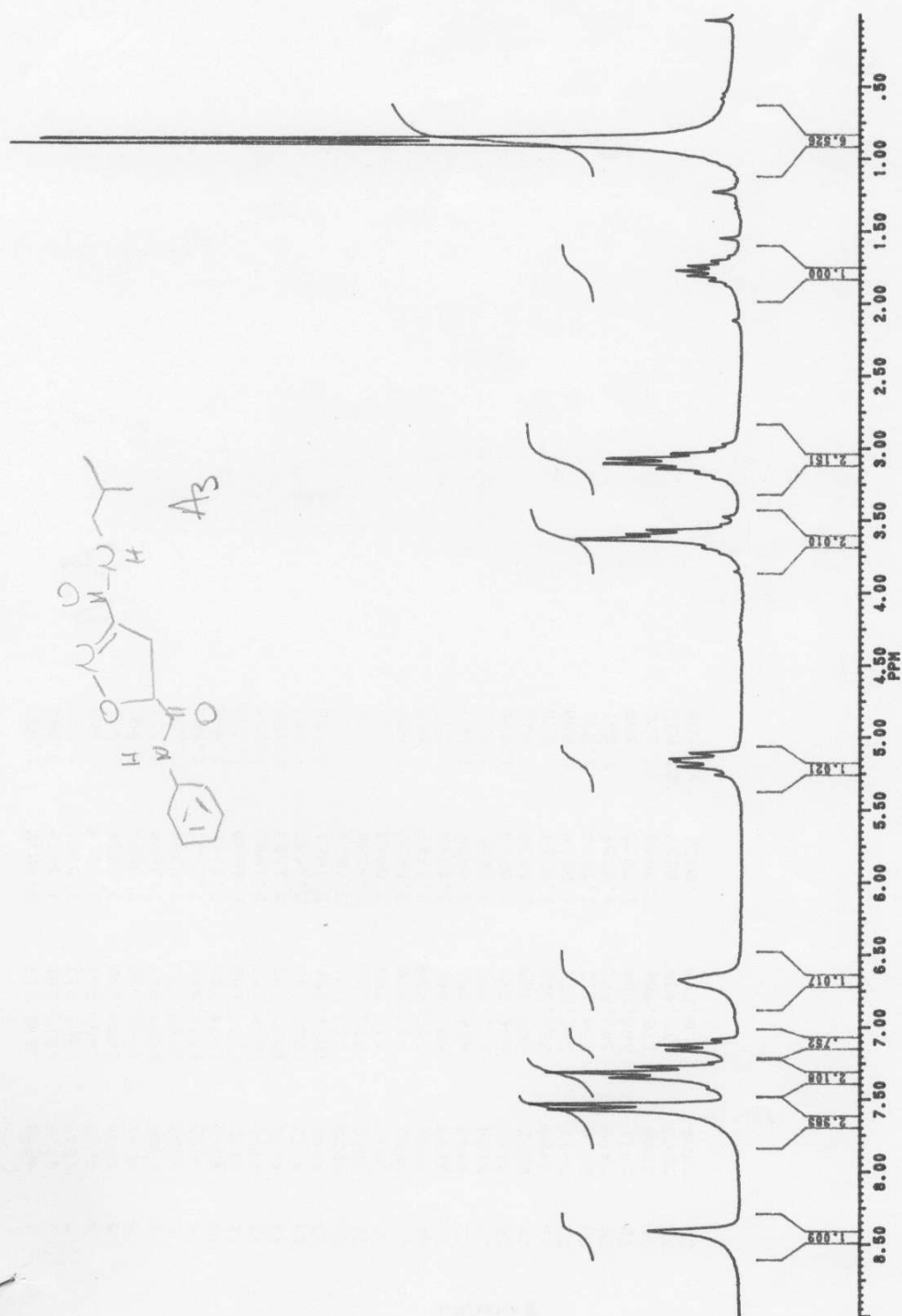
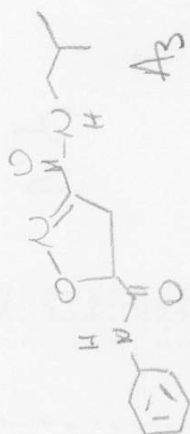


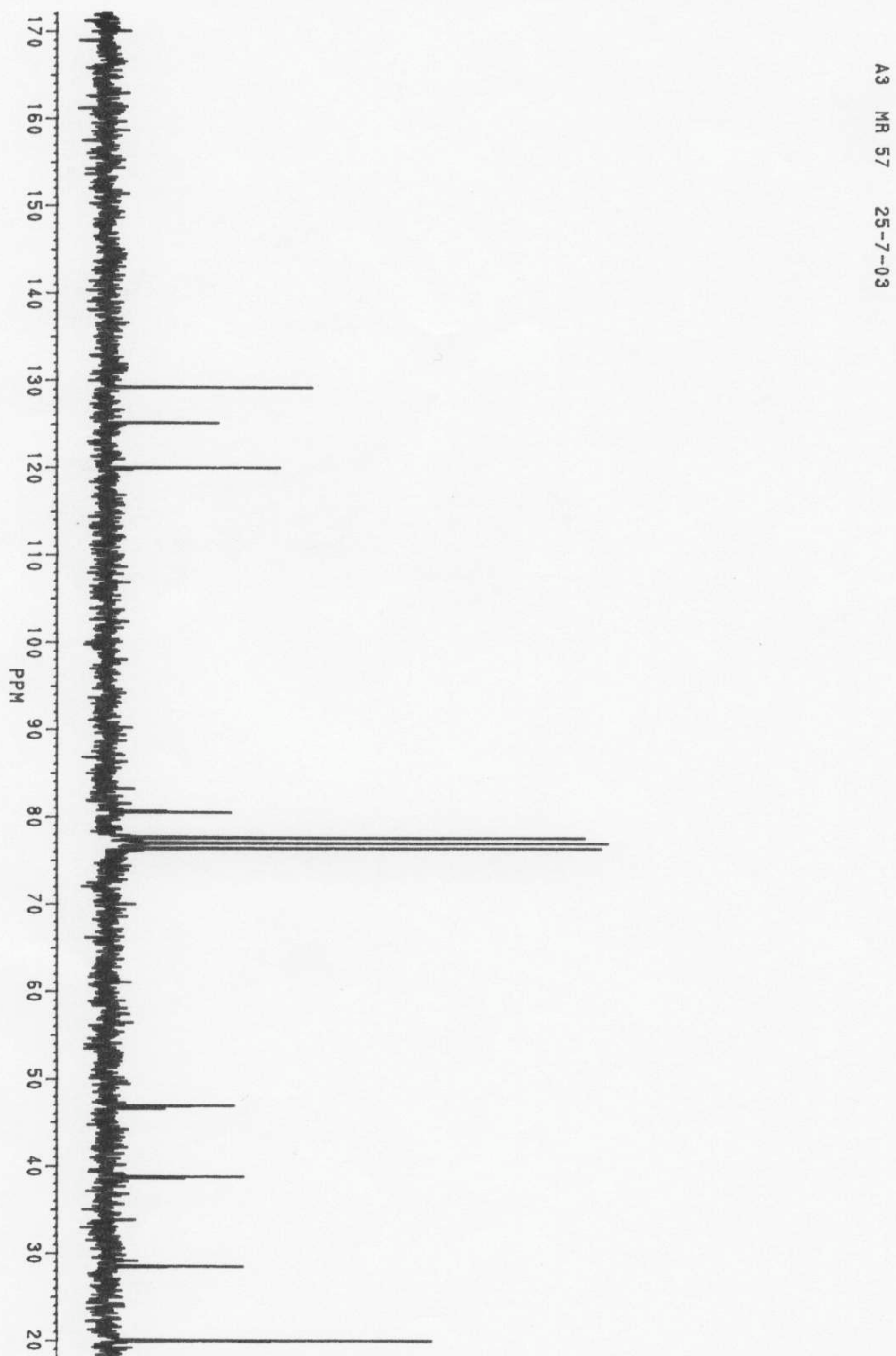
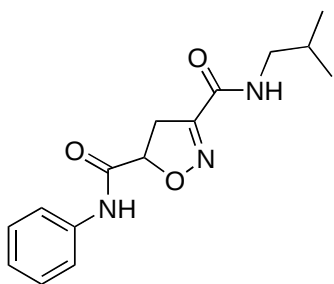
AZ = 4854





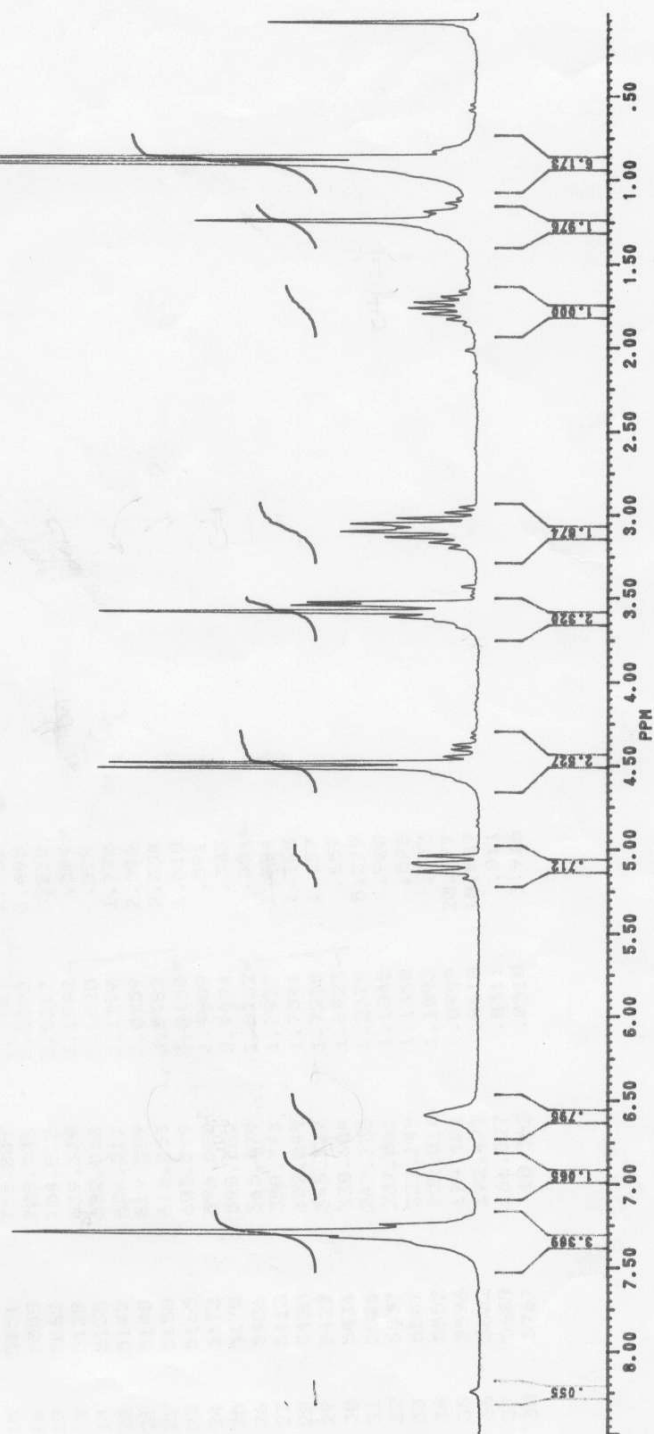
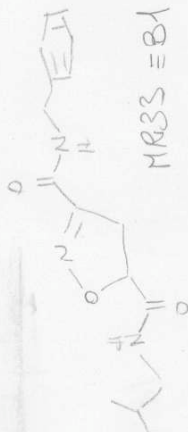
A3



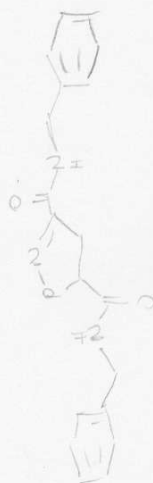


55% pure

MR 33 CDCL3 18/04/03

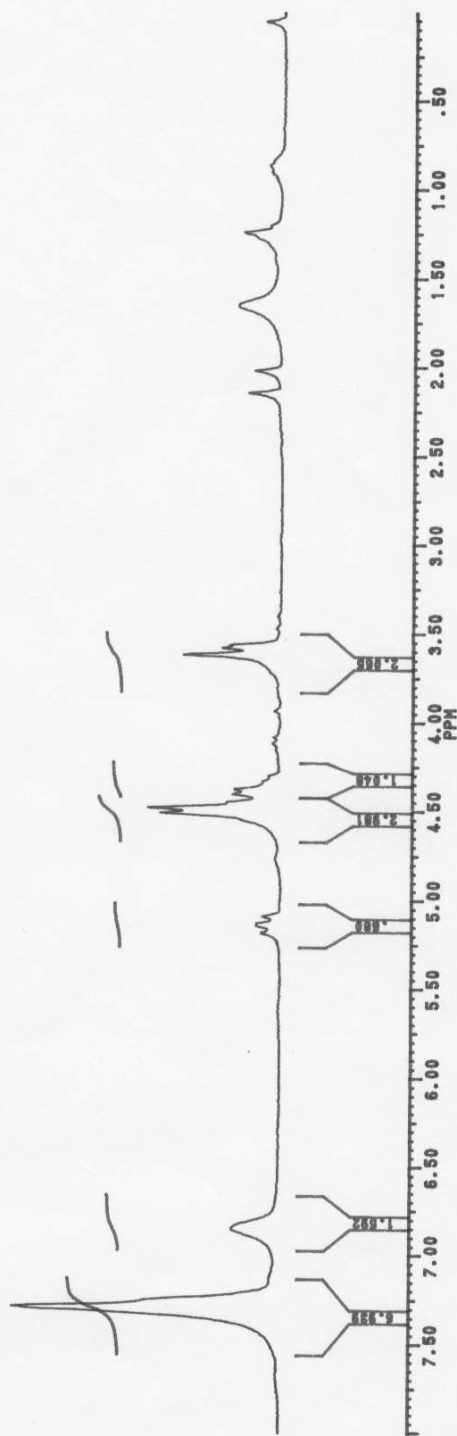


A2 36/52 CDCL3 27/05/03



B2

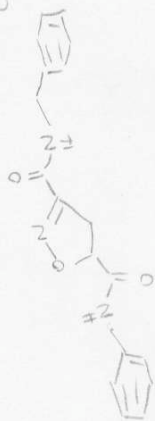
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1	4296	1458.815	7.2892	5.010
2	4385	1370.816	6.8495	.875
3	4724	1037.043	5.1818	.358
4	4733	1028.813	5.1406	.446
5	4861	902.583	4.5099	2.268
6	4867	896.907	4.4815	2.541
7	4882	881.540	4.4048	.906
8	4888	876.292	4.3785	.884
9	5042	724.062	3.6179	1.846
10	5050	716.339	3.5793	1.115
11	5053	713.423	3.5647	.982
12	5342	425.254	2.1448	.663
13	5367	404.212	2.0197	.526
14	5442	330.653	1.6522	.844
15	5526	248.073	1.2395	.757
16	5767	10.837	.0542	.362



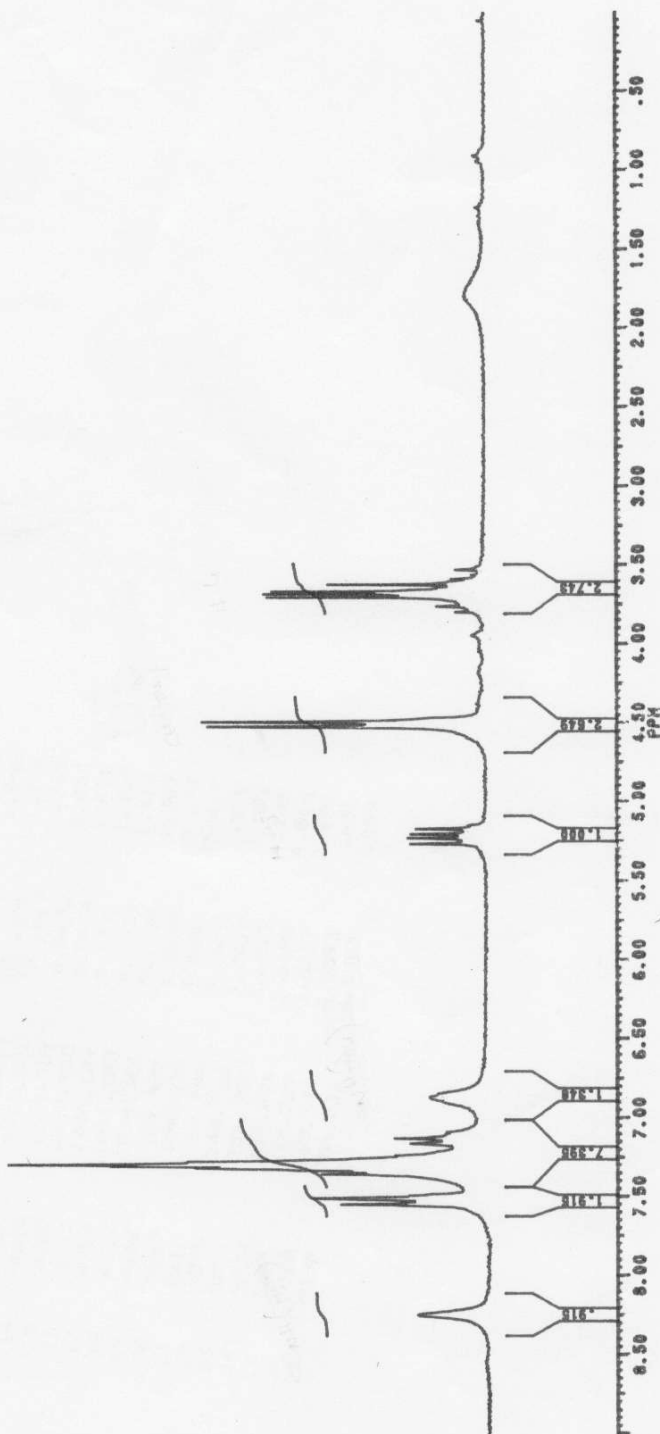
SAVINI 2

↓ B3

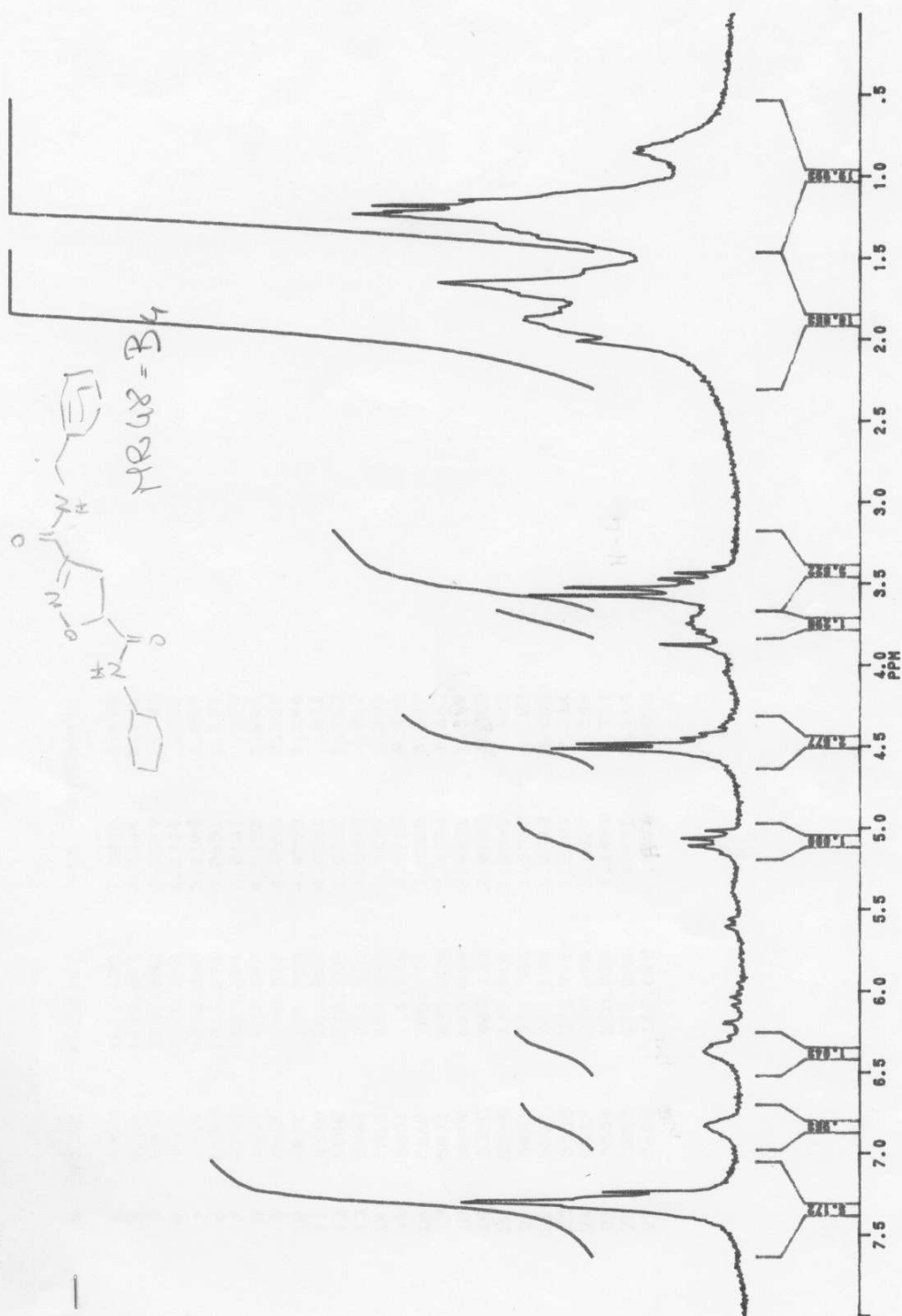
Ami S2B
Bund IX



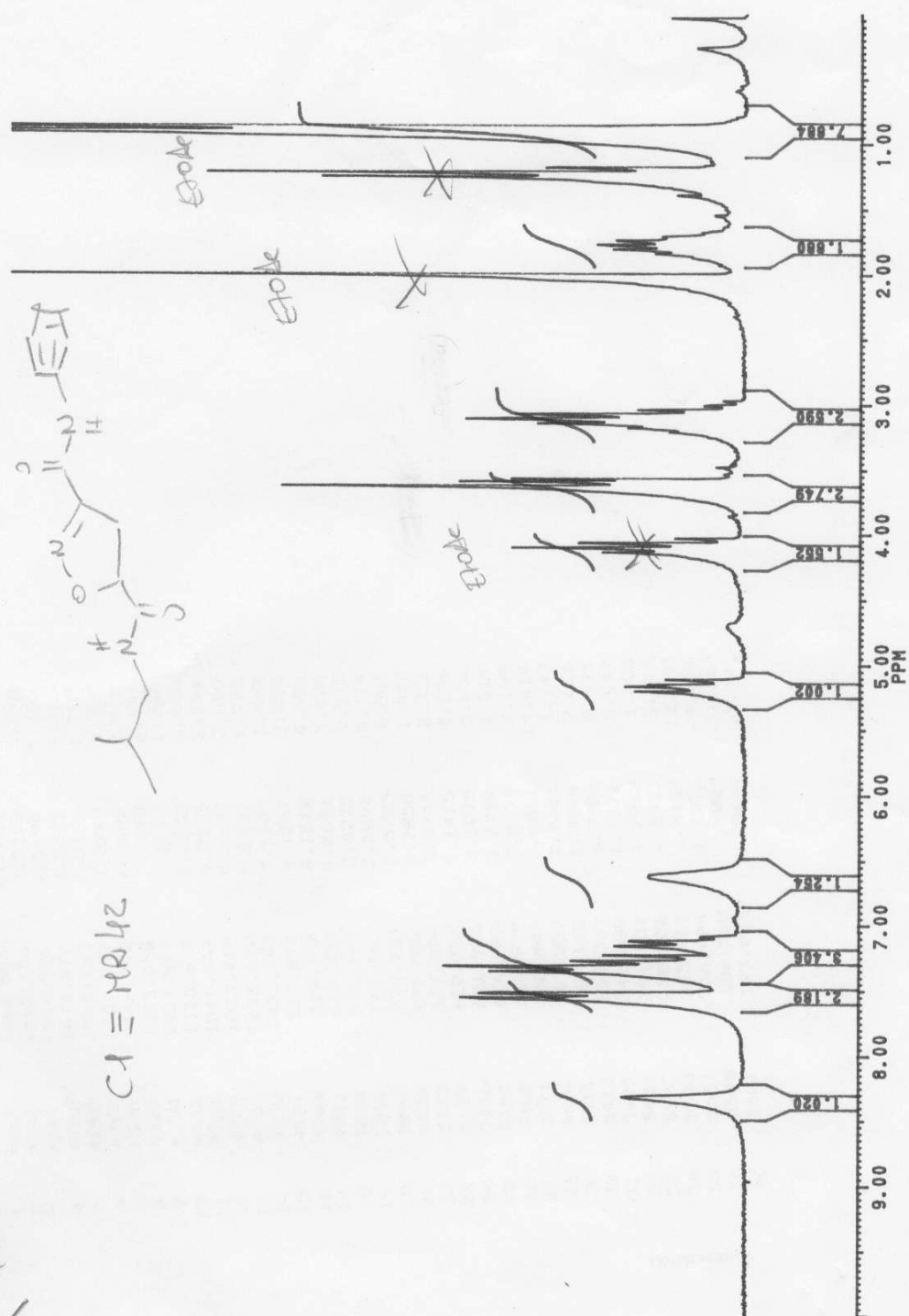
B3 = HR 56



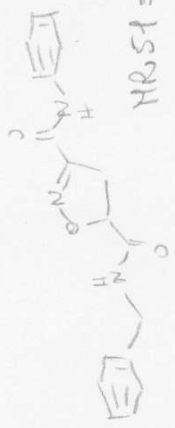
3 B4



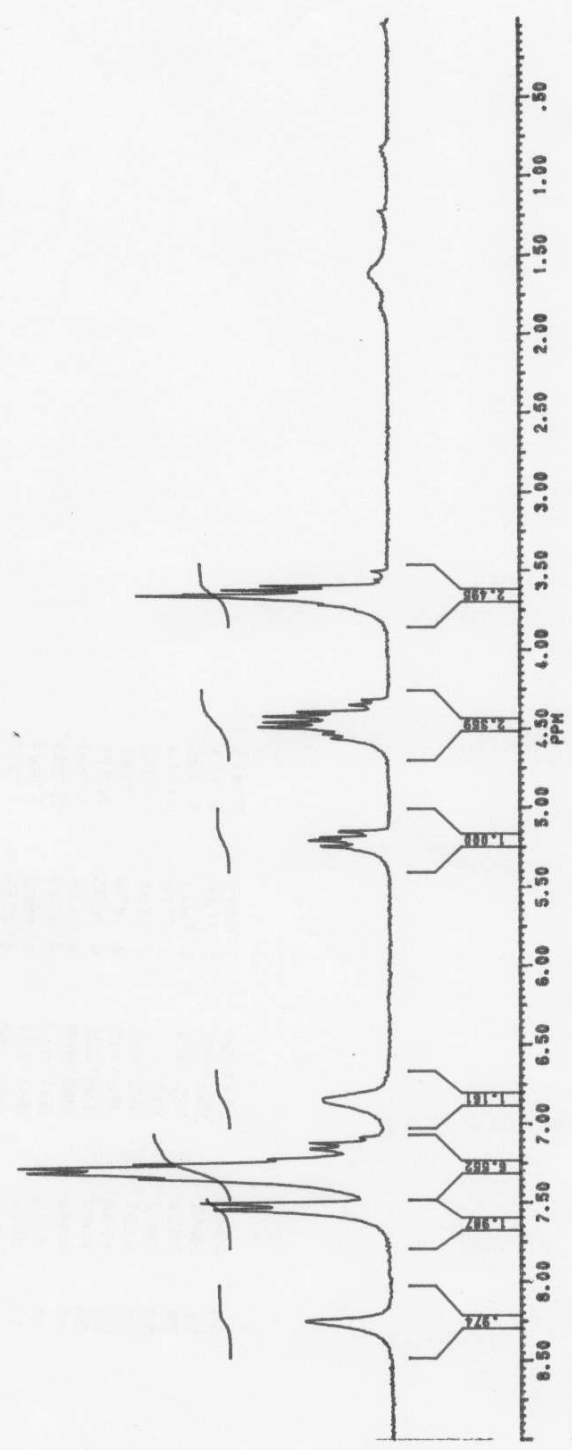
34 C1

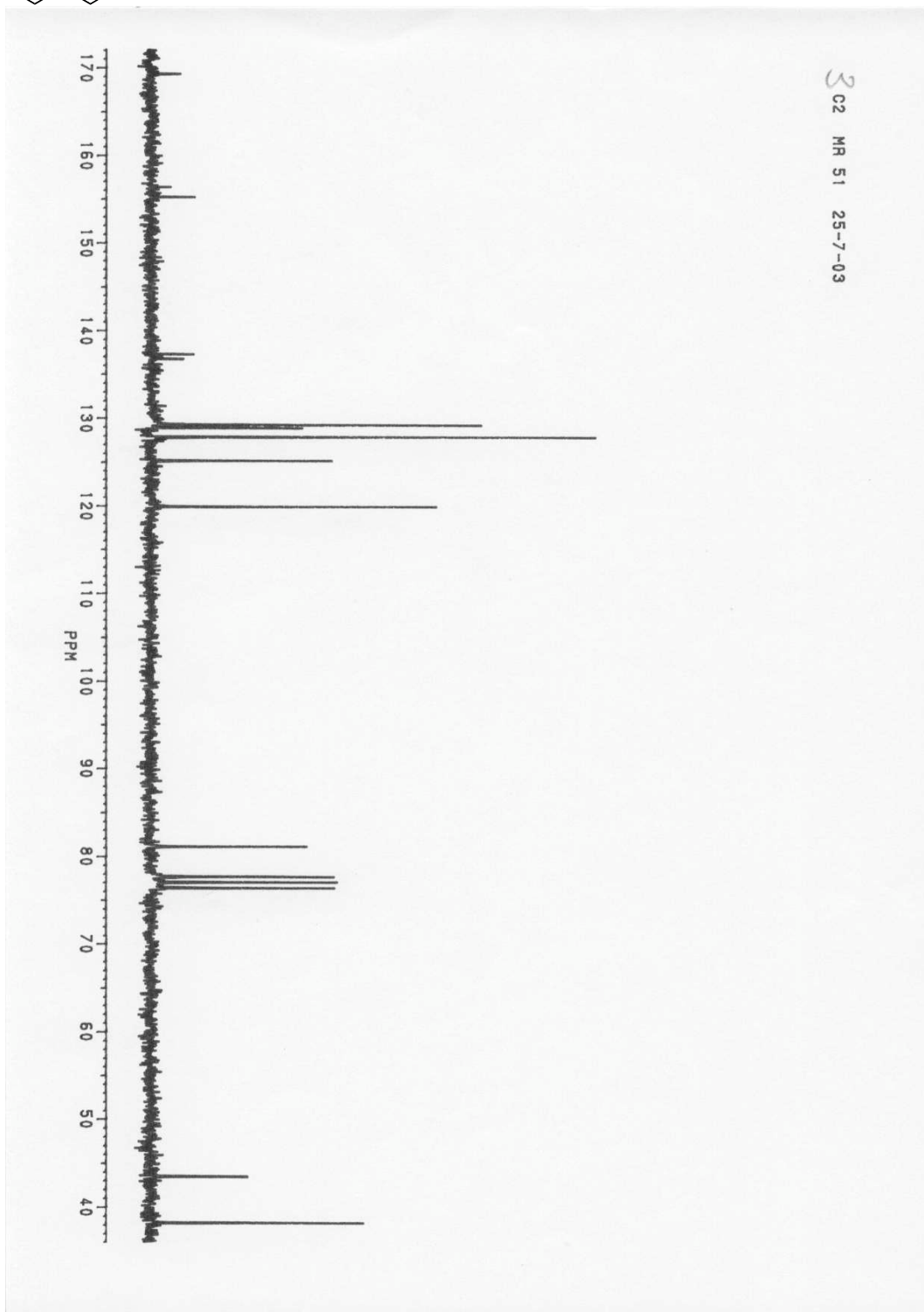
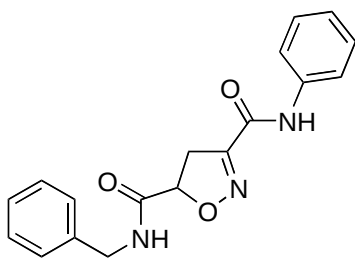


Pou NH₂ S&B
 Ani W



HR 51 = C2

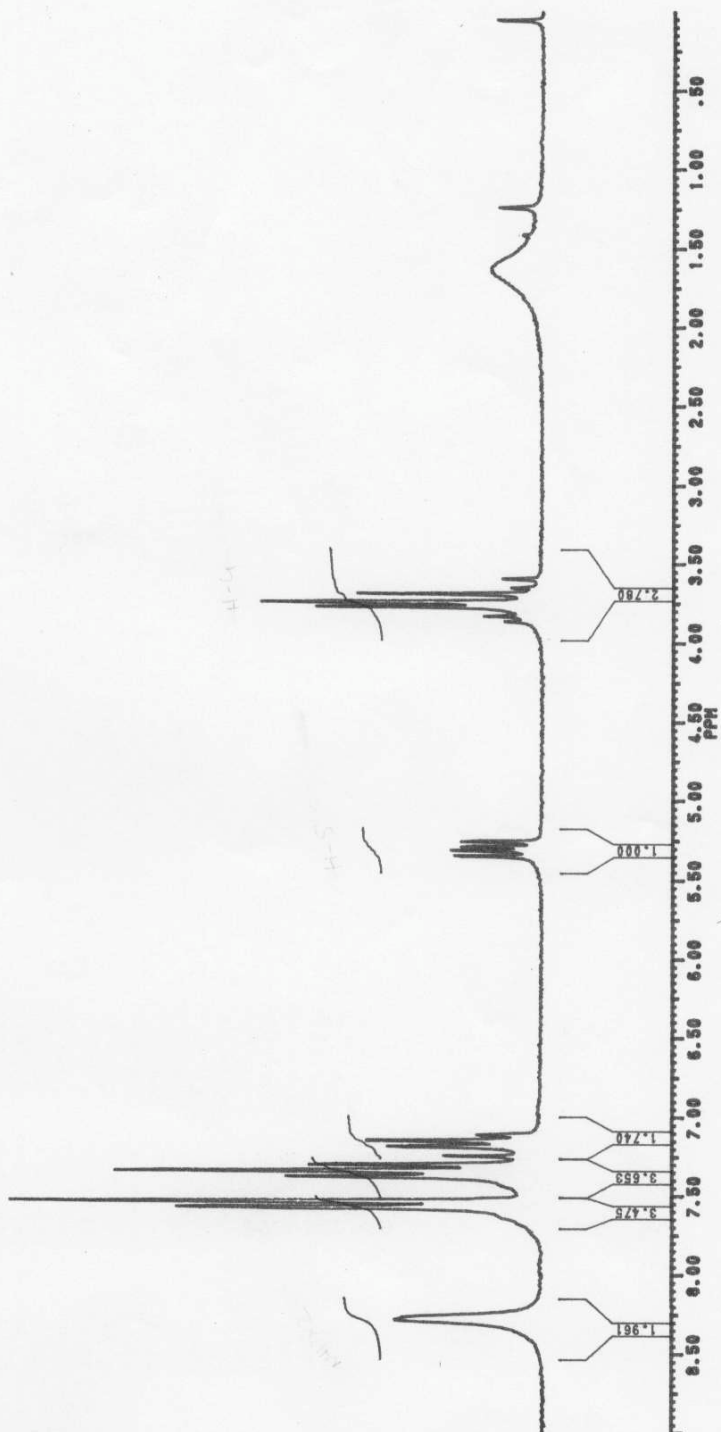
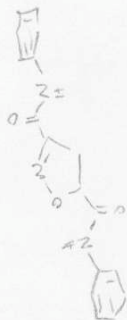




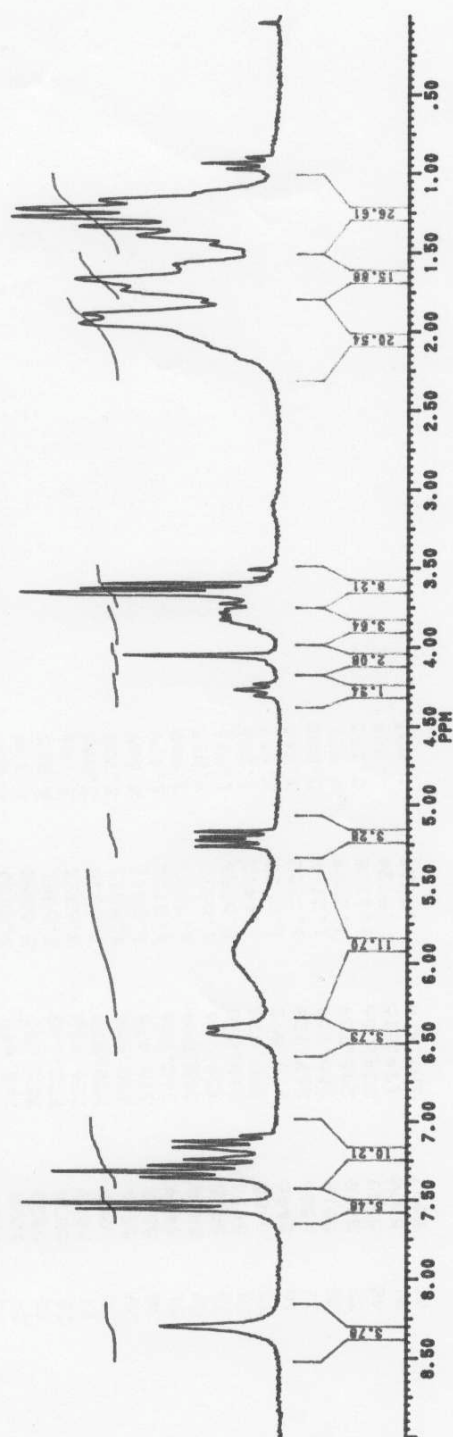
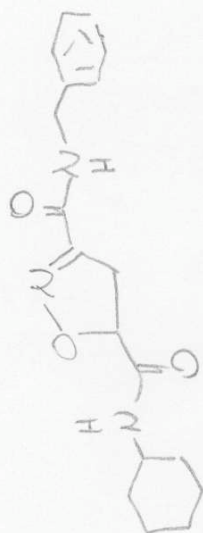
SPOT ALTO C3

HR 52 = C3

Anti S2B
Anti W

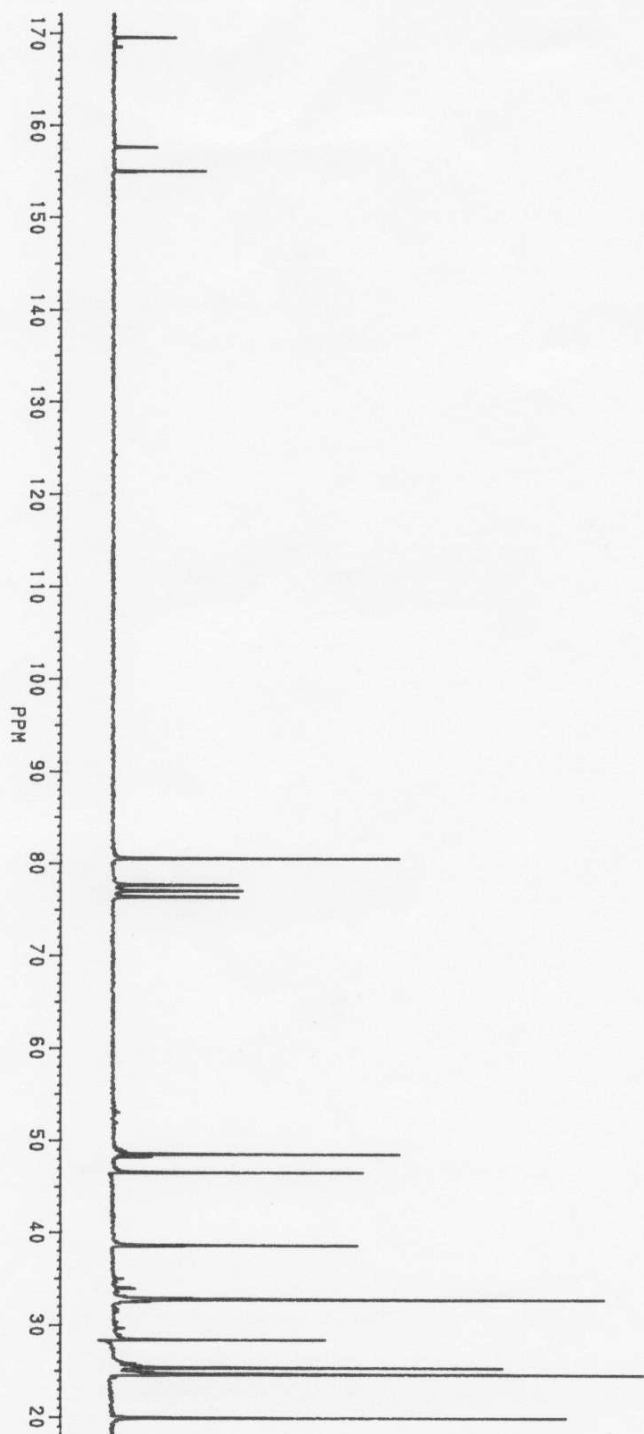
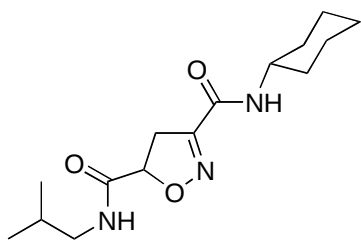


C4



DI = YR 41

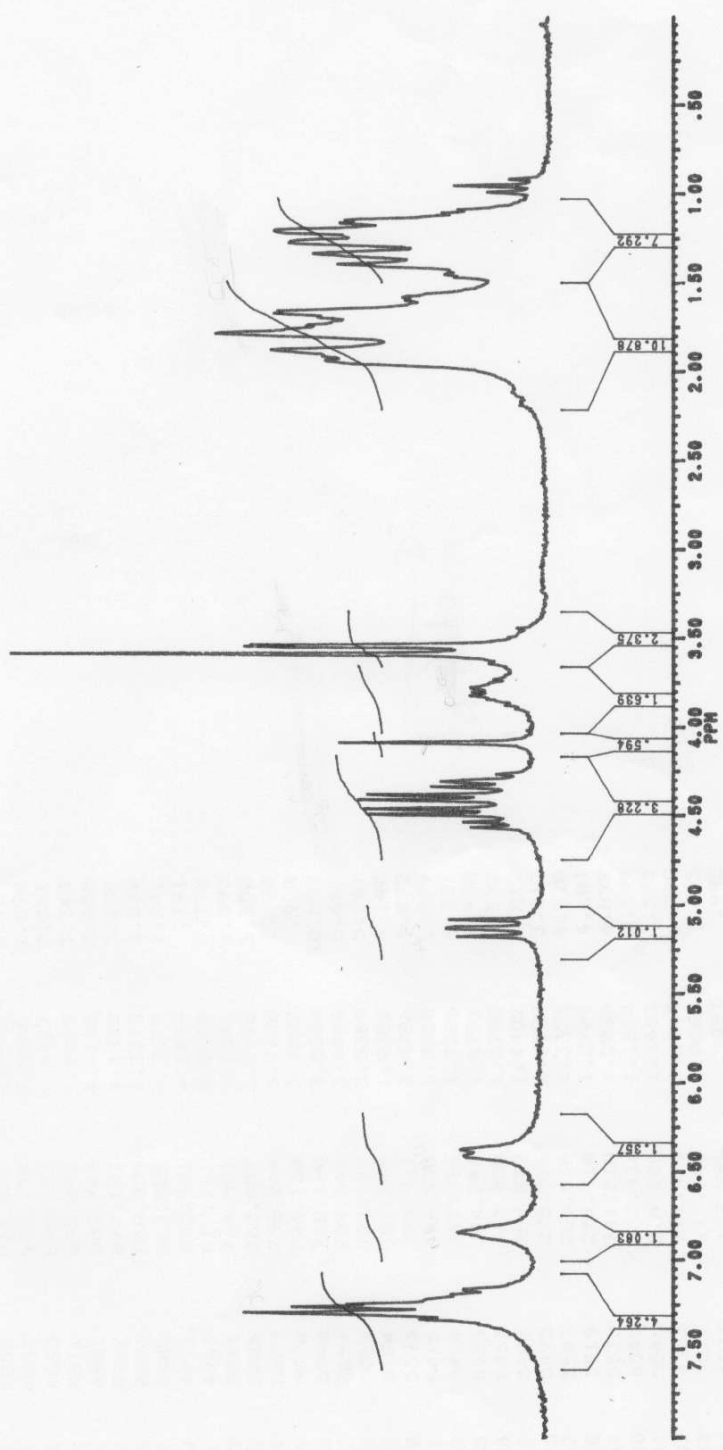
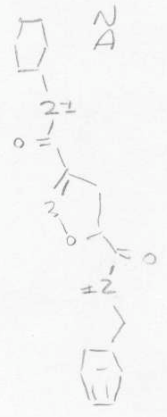
Chemical structure of compound 41 is shown in the top right corner. The structure is a substituted cyclohexane ring with a carboxylic acid group, a methyl group, and a side chain containing a double bond and a methyl group.

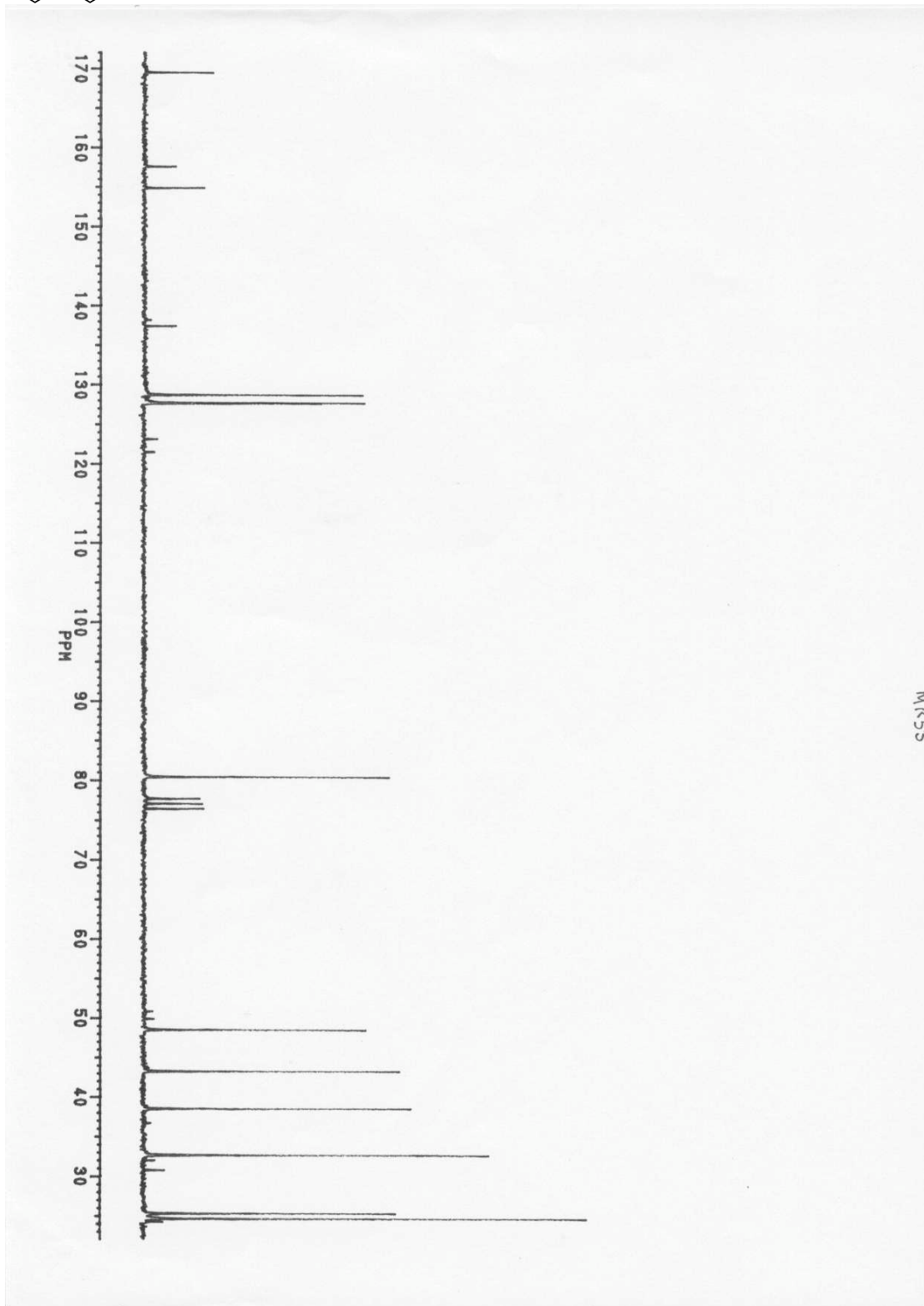
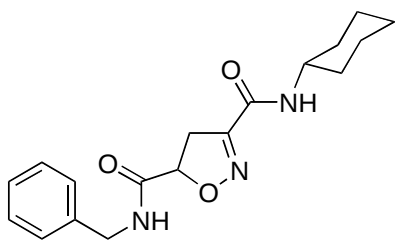


MR 4.1 25-7-03 3D1

SPOT ALTO PRECIPITATO D2

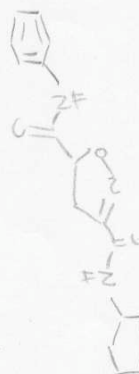
BuNH₂ S.B. acidomilung W



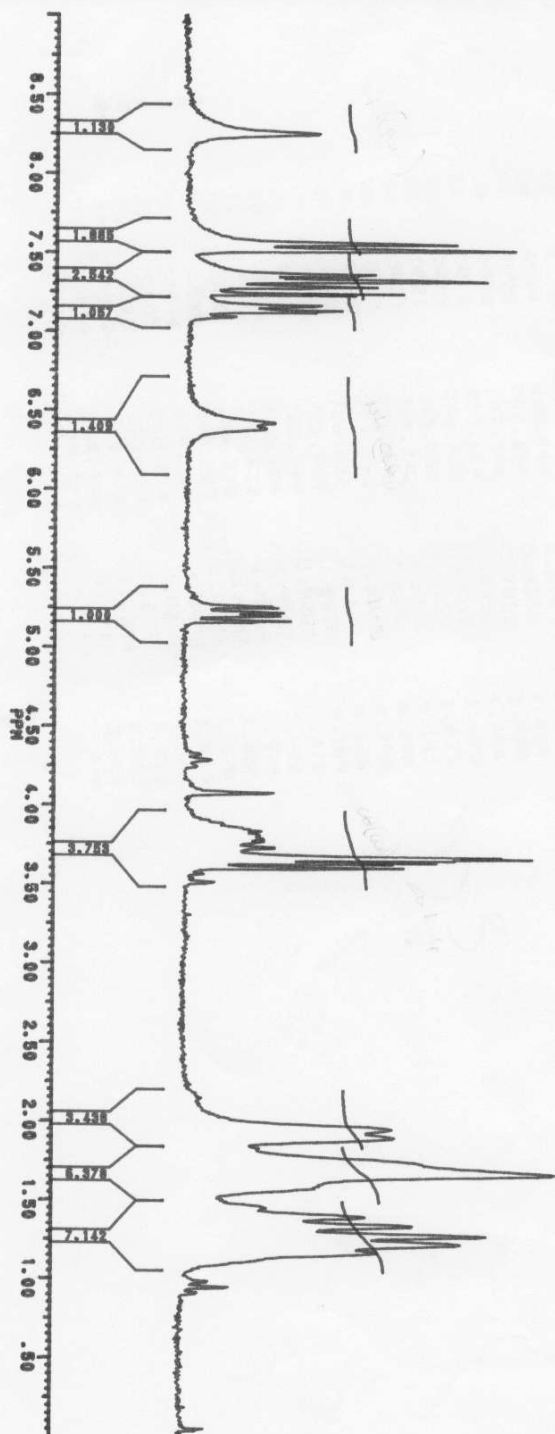


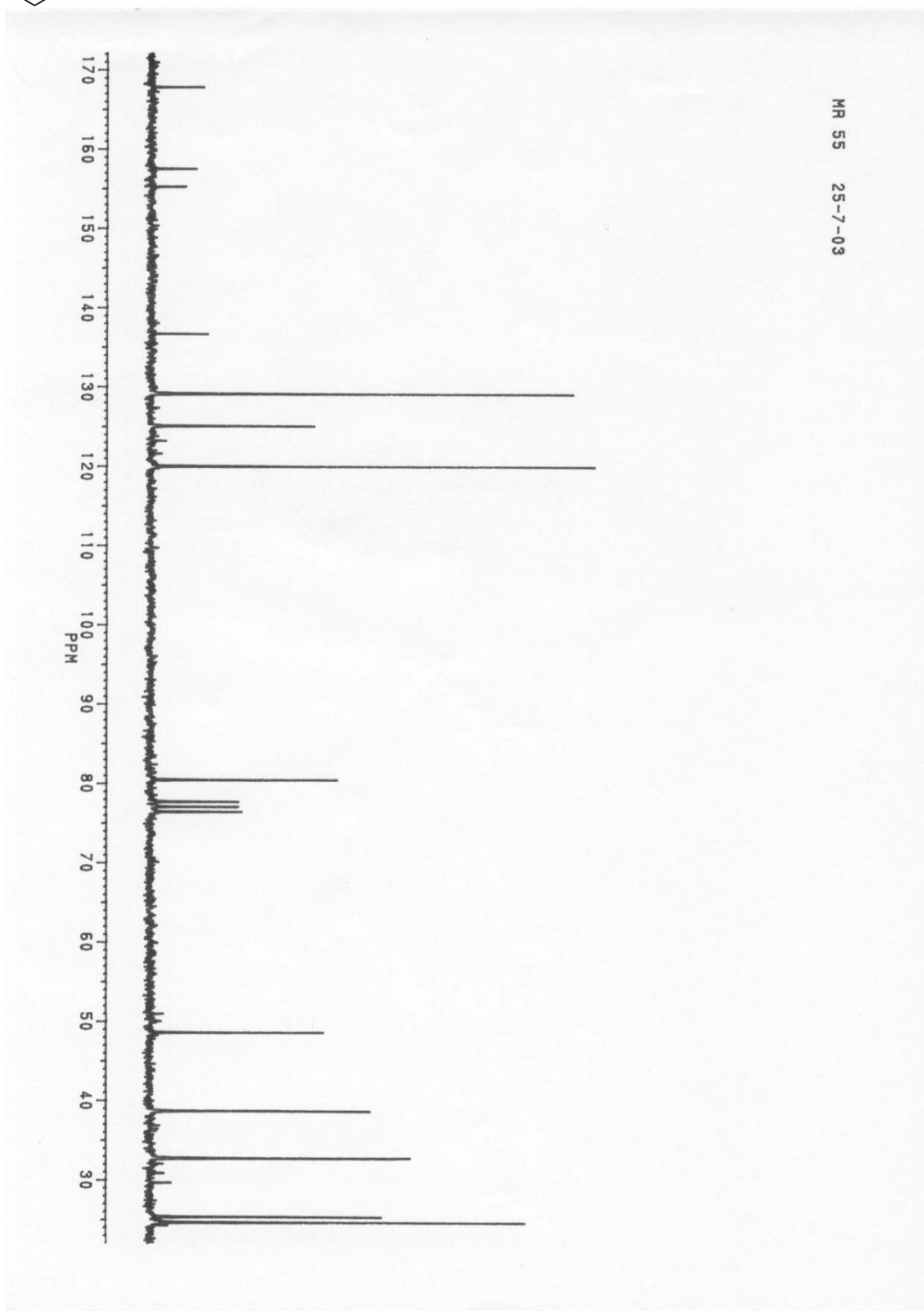
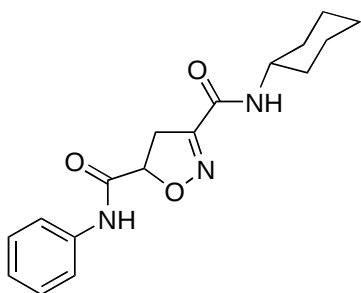
D3

D3 = NR SS



Ami S & B
CEA W





26/10 D4 CDCL3 29/05/03

#	CURSOR	FREQUENCY	PPM	INTENSITY
1	4477	1280.541	6.3984	.907
2	4744	1017.555	5.0844	.499
3	4754	1008.067	5.0370	.698
4	4763	998.564	4.9895	.323
5	4944	820.858	4.1016	.464
6	4951	813.881	4.0667	.477
7	5007	768.396	3.7895	.833
8	5018	748.089	3.7380	.936
9	5022	744.049	3.7178	.859
10	5026	740.191	3.6985	.852
11	5059	707.287	3.5341	2.477
12	5069	697.806	3.4867	1.943
13	5345	426.211	2.1296	.456
14	5370	401.100	2.0042	2.914
15	5394	377.601	1.8867	4.652
16	5440	332.358	1.6607	5.032
17	5512	261.799	1.3081	6.571
18	5523	250.946	1.2539	7.913
19	5530	244.107	1.2197	.959
20	5598	176.901	.8839	.975
21	5605	169.628	.8476	.769
22	5771	6.654	.0332	

