

Appendix A  
NMR Spectra

## NMR Spectral Data.

### 2,4-D

300 MHz proton NMR, solvent  $\text{CDCl}_3$ . 7.43 ppm, 1 proton doublet, H3; 7.21 ppm 1 proton doublet of doublets H5; 6.84 ppm 2 proton doublet H6; 4.76 ppm 2 proton singlet  $\text{CH}_2$ . Consistent with expected spectrum of 2,4-D.

### Alachlor

300 MHz proton NMR, solvent  $\text{CDCl}_3$ . 7.31 ppm, 1 proton multiplet, H4; 7.21 ppm 1 proton doublet of doublets H5; 4.96 ppm 2 proton singlet  $\text{N-CH}_2\text{-O}$ ; 3.72 ppm 2 proton singlet  $\text{CH}_2\text{Cl}$ ; 3.51 ppm 3 proton singlet  $\text{OCH}_3$ ; 2.59 ppm, 4 proton multiplet 2 x  $\text{CH}_2\text{-CH}_3$ ; 2.59 ppm, 6 proton multiplet 2 x  $\text{CH}_2\text{-CH}_3$ . Consistent with expected spectrum of alachlor.

### Bensulide

300 MHz proton NMR, solvent  $\text{CD}_3\text{OD}$ . 7.85 ppm, 2 proton multiplet, H2, H6; 7.63 ppm 3 proton multiplet H3, H4, H5; 4.73 ppm 2 proton singlet 2 x  $\text{O-CH-(CH}_3)_2$ ; 3.13 ppm 2 proton multiplet  $\text{NH-CH}_2$ ; 2.90 ppm 2 proton multiplet  $\text{CH}_2\text{-S}$ ; 12 proton triplet, 2 x  $\text{O-CH-(CH}_3)_2$ . Consistent with expected spectrum of bensulide.

### Chloridazon

300 MHz proton NMR, solvent  $\text{CD}_3\text{OD}$ . 7.77 ppm, 1 proton singlet H4, 5 proton multiplet at 7.42 ppm, consistent with the 4 protons on the benzene ring and the single proton on the pyridazine ring. Consistent with the expected spectrum of chloridazon.

### Diazinon

300 MHz proton NMR, solvent  $\text{CDCl}_3$ . 6.77 ppm 1 proton singlet, ring; 4.35 ppm 4 proton multiplet 2 x  $\text{CH}_2$ ; 2.55 ppm 3 proton singlet  $\text{CH}_3$ ; 3.23 ppm 1 proton singlet  $\text{CH}$ ; 12 proton multiplets at approximately 1.4 ppm 4 x  $\text{CH}_3$ . Consistent with the expected spectrum of chloridazon.

### Flufenacet

300 MHz proton NMR, solvent  $\text{CDCl}_3$ . 7.24 ppm, 4 proton multiplet ring protons; 4.97 ppm, 1 proton multiplet  $\text{CH}$ ; 4.75 ppm, 2 proton multiplet  $\text{CH}_2$ ; 1.10 ppm, 6 protons, 2 x  $\text{CH}_3$ . Consistent with the expected NMR spectrum for flufenacet.

### Propyzamide

300 MHz proton NMR, solvent  $\text{CDCl}_3$ . 7.64 ppm, 2 proton multiplet, H2, H5; 7.54 ppm 1 proton multiplet H4; 6.13 ppm 1 proton singlet  $\text{NH}$ ; 2.43 ppm 1 proton singlet  $\text{CH}$ ; 1.80 ppm, 6 proton singlet 2 x  $\text{CH}_3$ . Consistent with expected spectrum of propyzamide

### Pyrithiobac

300 MHz proton NMR, solvent  $\text{CD}_3\text{OD}$ . 7.62 ppm, 1 proton multiplet H4, 2 x 1 proton multiplet at 7.4 and 7.26 ppm, consistent with the H3 and H5 protons on the benzene ring; 5.74 ppm, 1 proton singlet

for the pyrimidine ring proton; 3.73 ppm 6 proton singlet 2 x CH<sub>3</sub>. Consistent with expected spectrum of Pyriithiobac

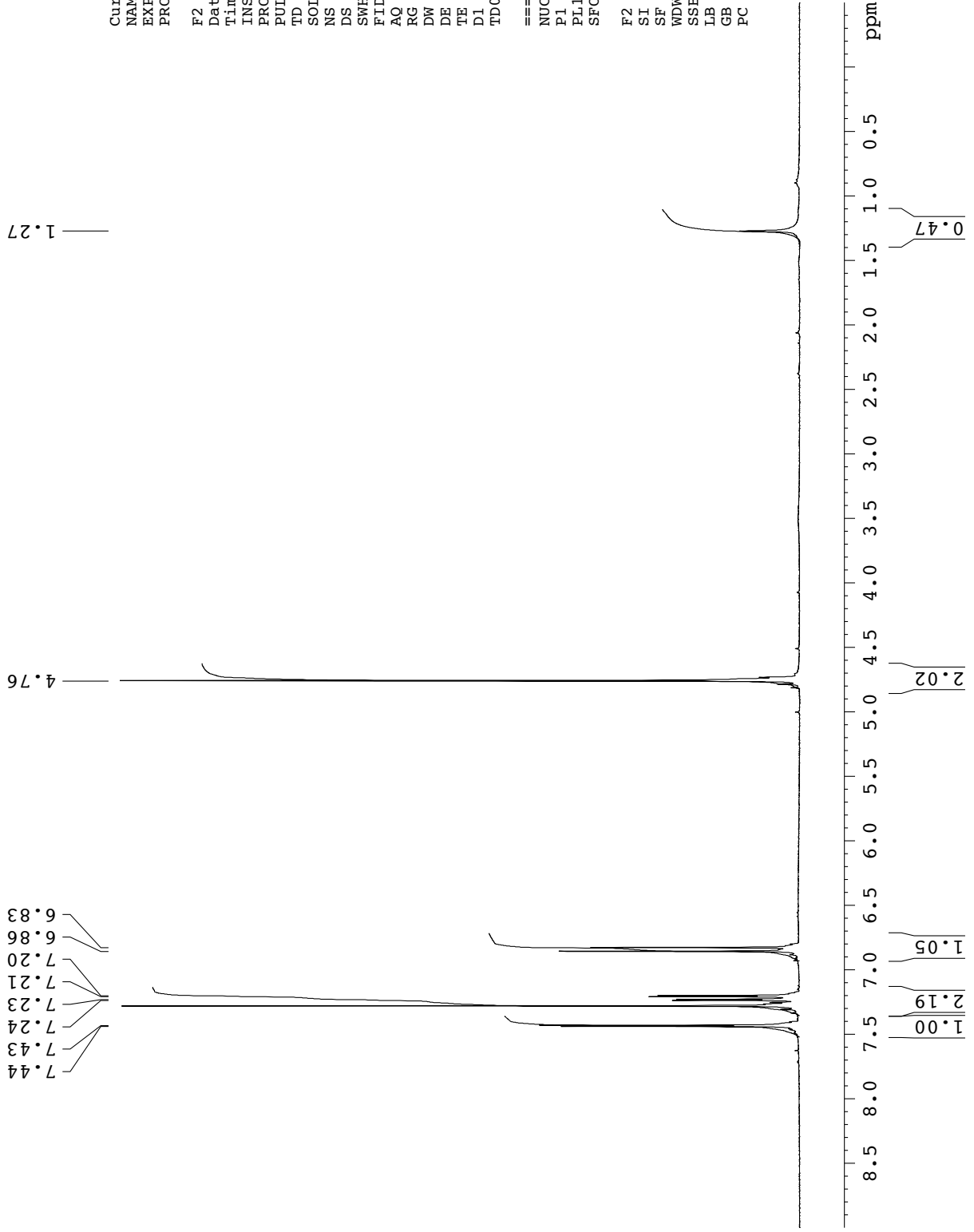
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 NS 64  
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 FIDRES 0.094190 Hz  
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 SSB 0  
 LB 0.30 Hz  
 GB 0  
 PC 1.00



# Alachlor PESTANAL



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7.36  
7.34  
7.31  
7.28  
7.25  
7.22

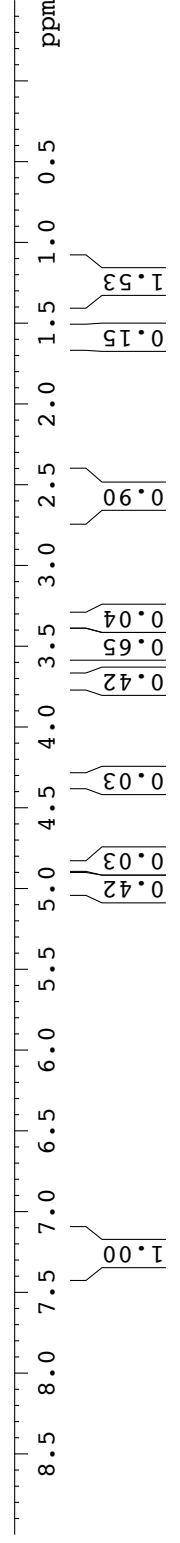
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2.59  
2.59  
2.57  
2.56  
2.54  
2.51  
2.49  
1.59  
1.29  
1.26  
1.24

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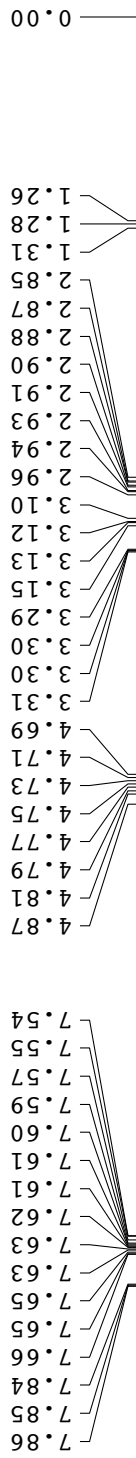
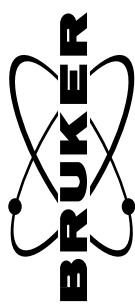
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SOLVENT CDCl3  
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DS 2  
SWH 6172.839 Hz  
FIDRES 0.094190 Hz  
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TD0 1

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Bensulide PESTANAL



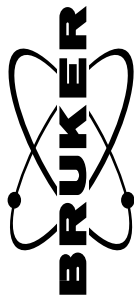
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TD 65536  
SOLVENT MeOD  
NS 64  
DS 2  
SWH 6172.839 Hz  
FIDRES 0.094190 Hz  
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# Chloridazon PESTANAL



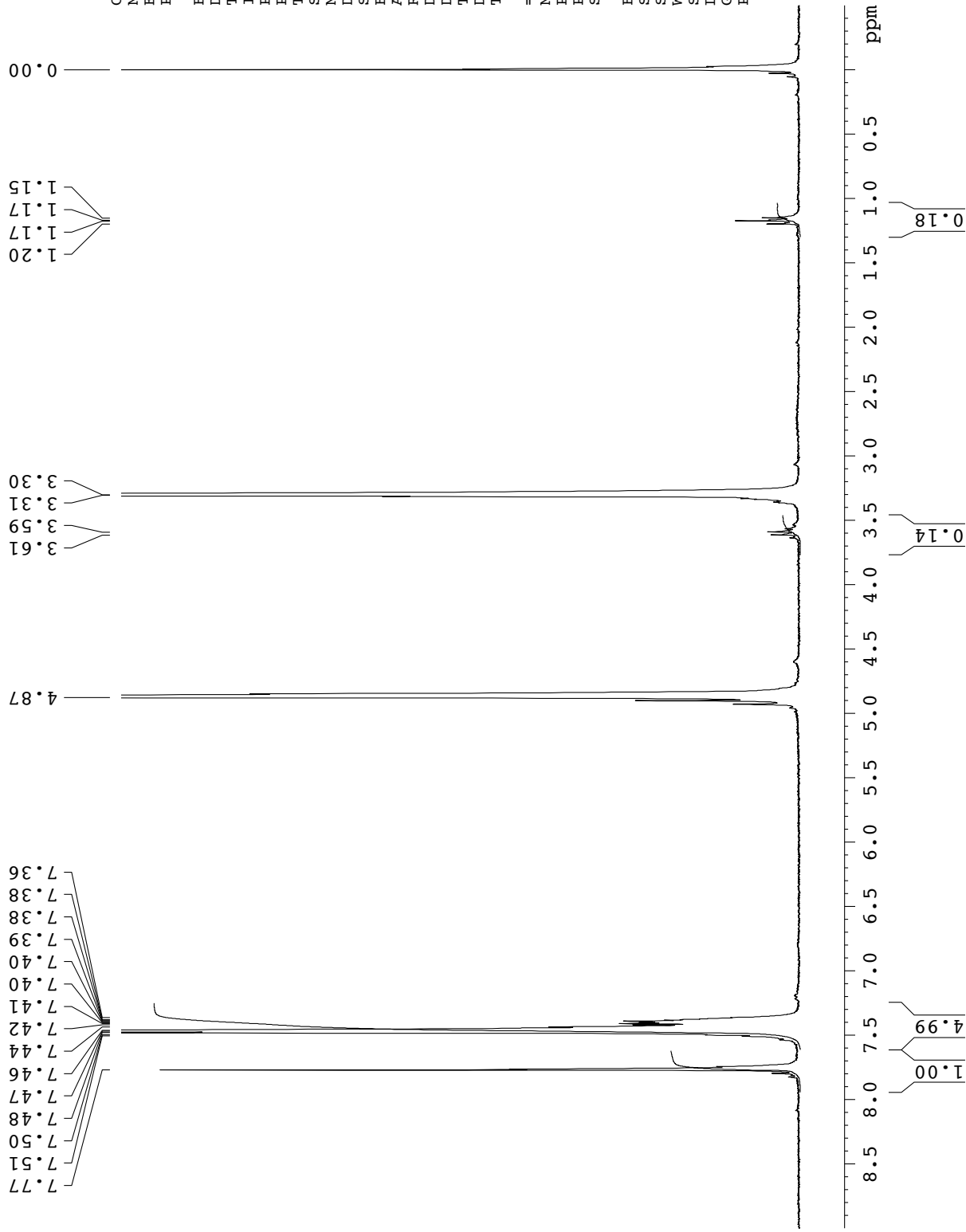
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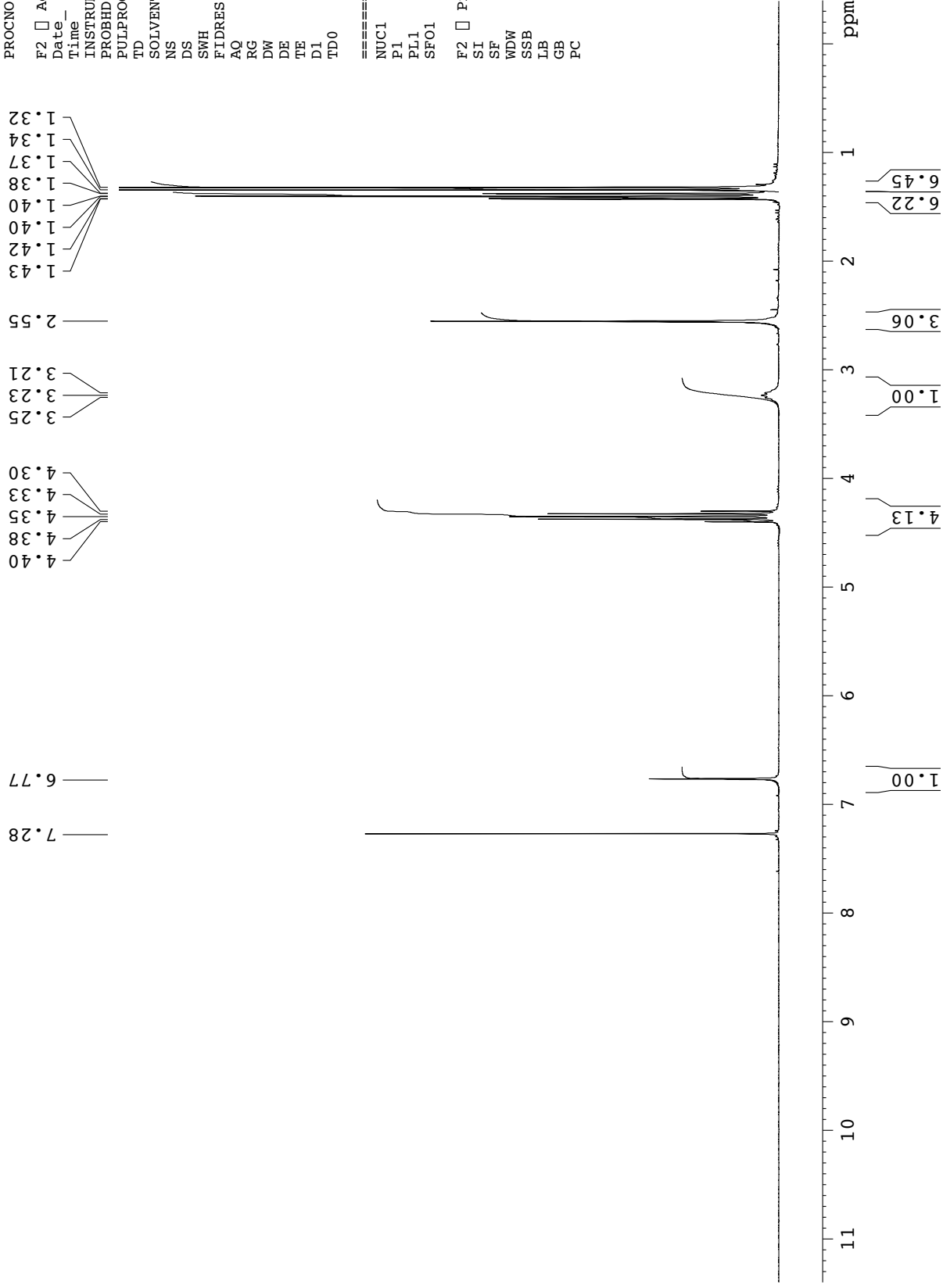
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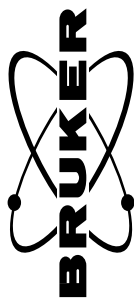
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Current Data Parameters  
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 TD 65536  
 SOLVENT CDCl<sub>3</sub>  
 NS 16  
 DS 2  
 SWH 6172.839 Hz  
 FIDRES 0.094190 Hz  
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 PL1 6.00 dB  
 SFO1 300.1318534 MHz  
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# Flufenacet PESTANAL

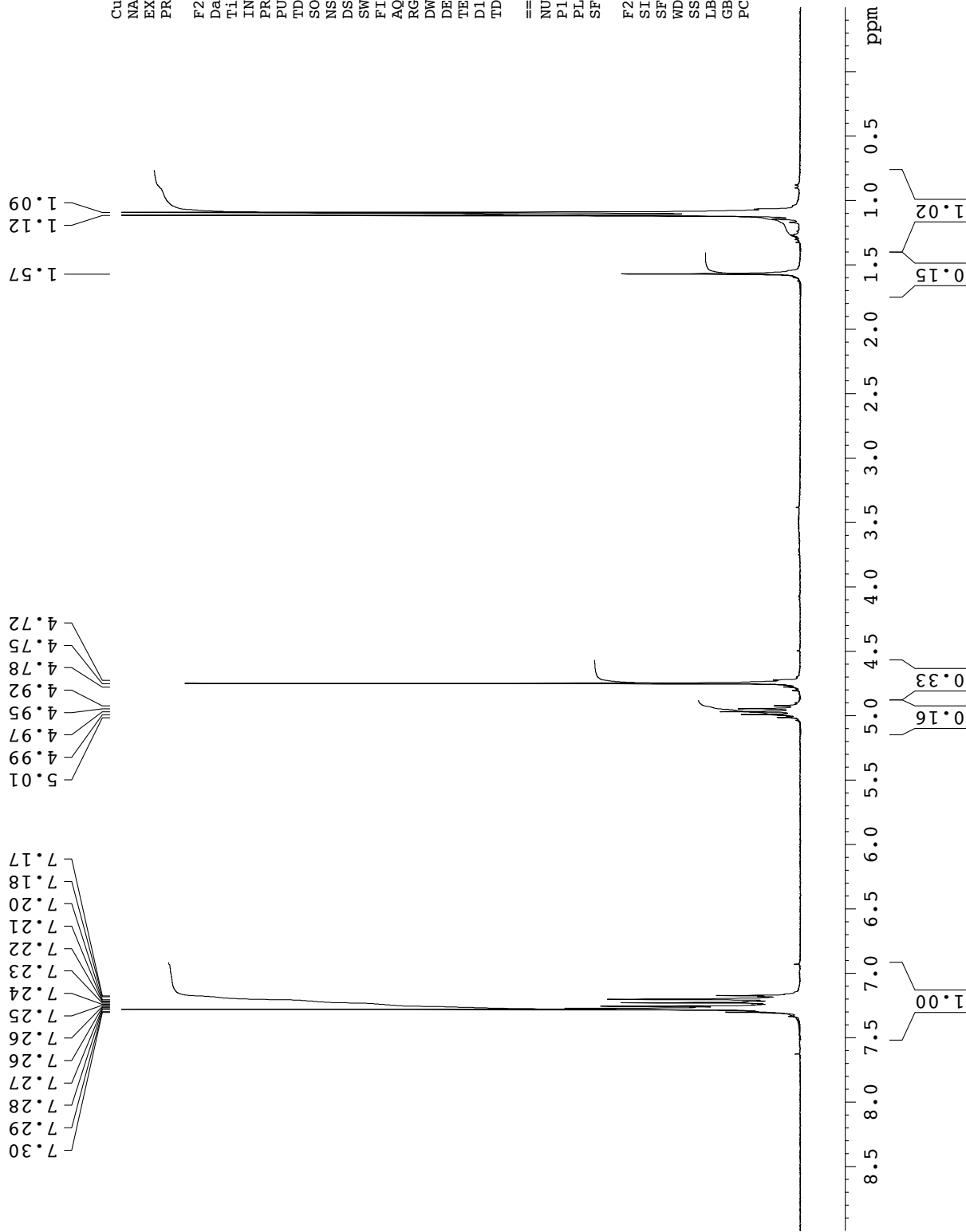


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 SOLVENT CDC13  
 NS 64  
 DS 2  
 SWH 6172.839 Hz  
 FIDRES 0.094190 Hz  
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F2 Processing parameters  
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# Propyzamid PESTANAL



8.18  
8.18  
7.94  
7.93  
7.92  
7.91  
7.85  
7.84  
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6.22  
6.13

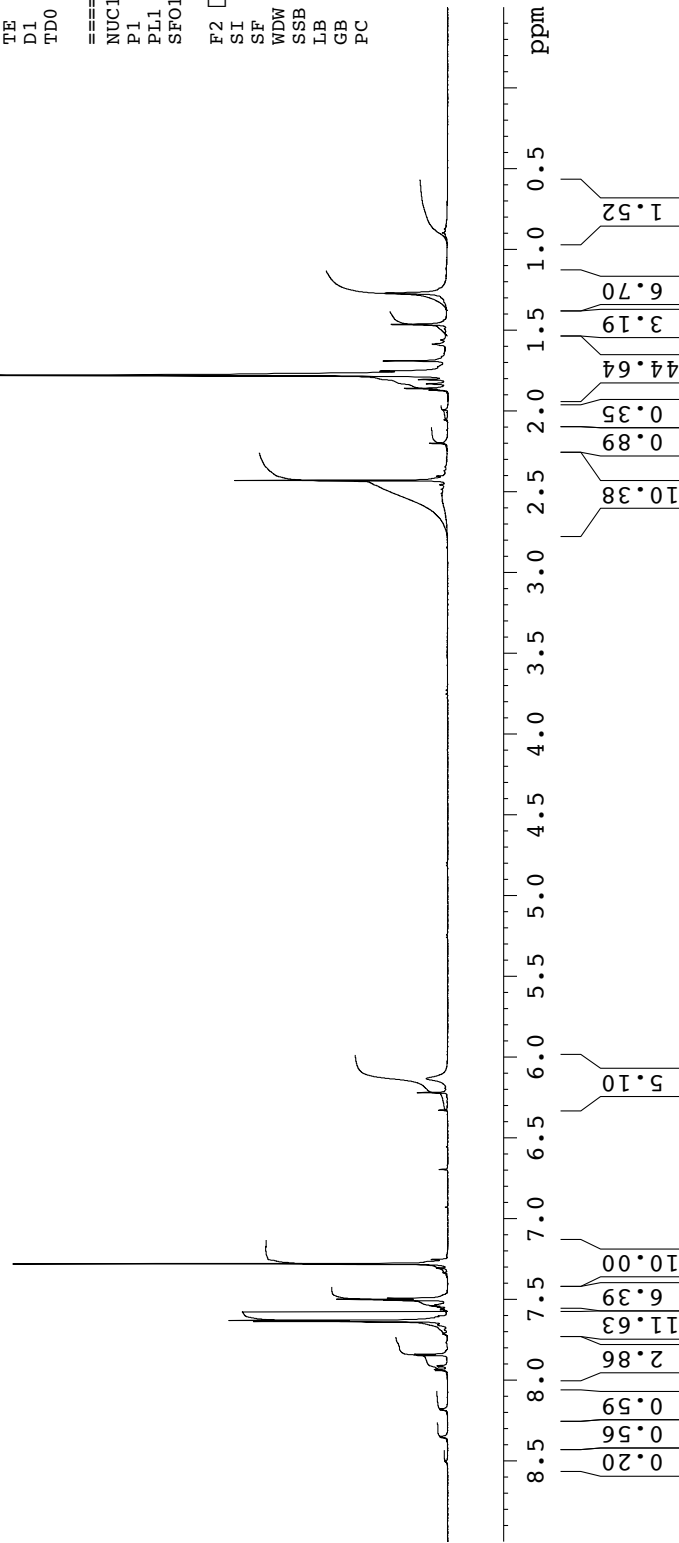
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1.27

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PROCNO 1

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SOLVENT CDCl3  
NS 64  
DS 2  
SWH 6172.839 Hz  
FIDRES 0.094190 Hz  
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PC 1.00



# Pyrethiobac Sodium

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PROCNO 1

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TD 65536  
SOLVENT MeOD  
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DS 2  
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FIDRES 0.094190 Hz  
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TE 300.0 K  
D1 1.00000000 sec  
TD0 1

==== CHANNEL f1 =====  
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PL1 6.00 dB  
SFO1 300.1318534 MHz

F2 Processing parameters  
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GB 0  
PC 1.00

