

**Intramolecular functional group differentiation as a strategy for synthesis of bridged bicyclic
β-amino acids**

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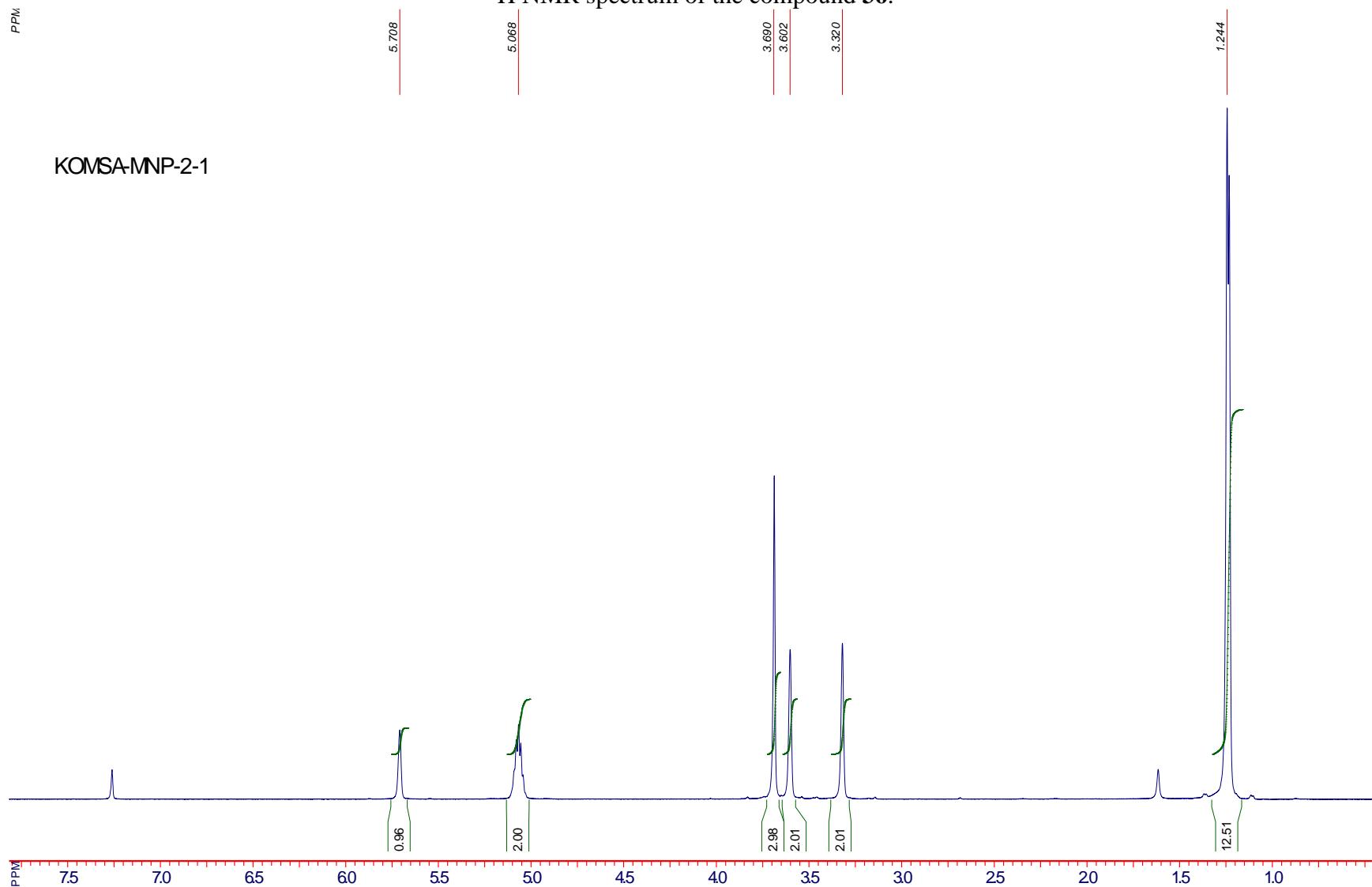
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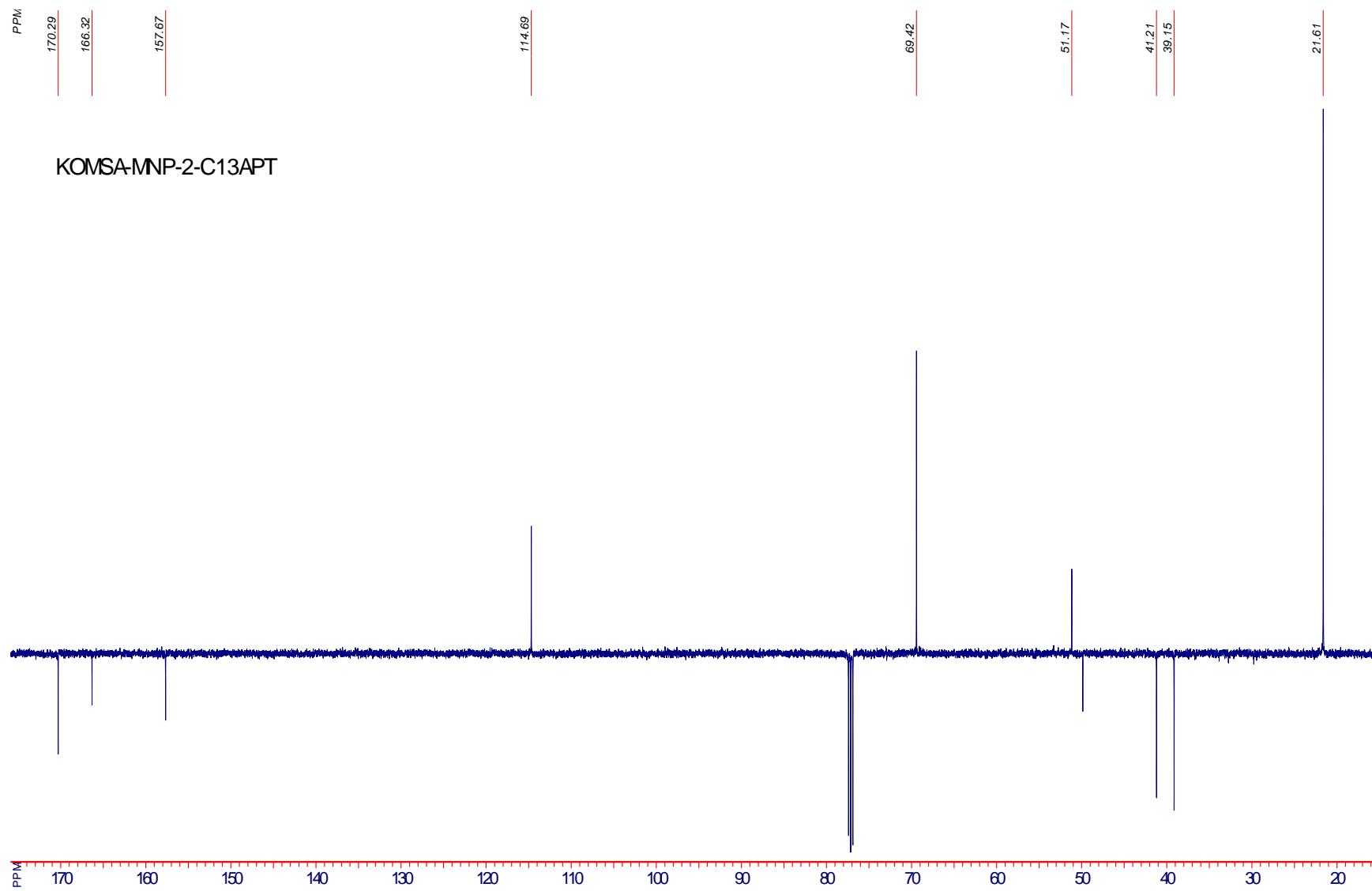
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PPM

¹H NMR spectrum of the compound 36.

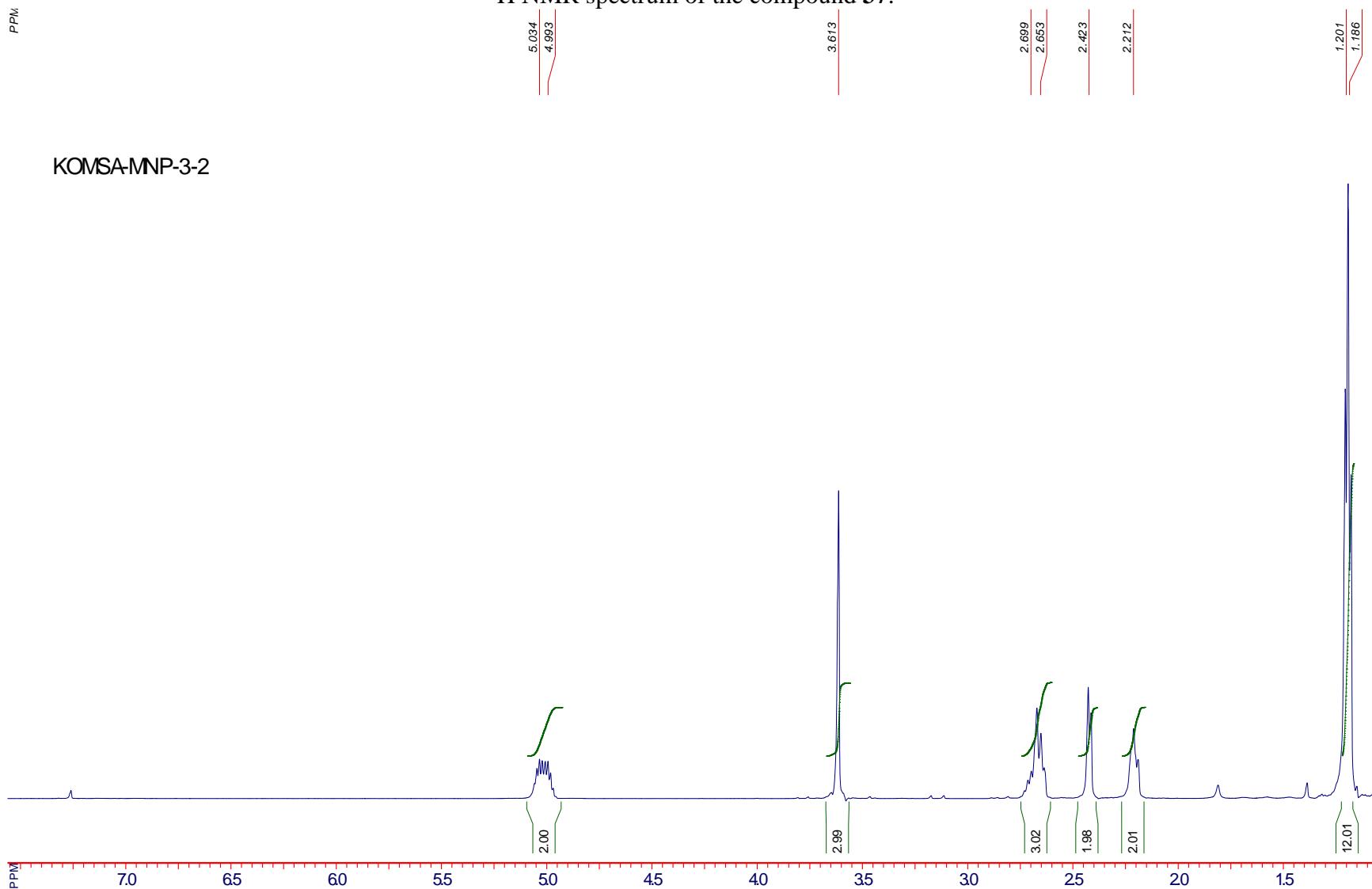


^{13}C NMR spectrum of the compound **36**.

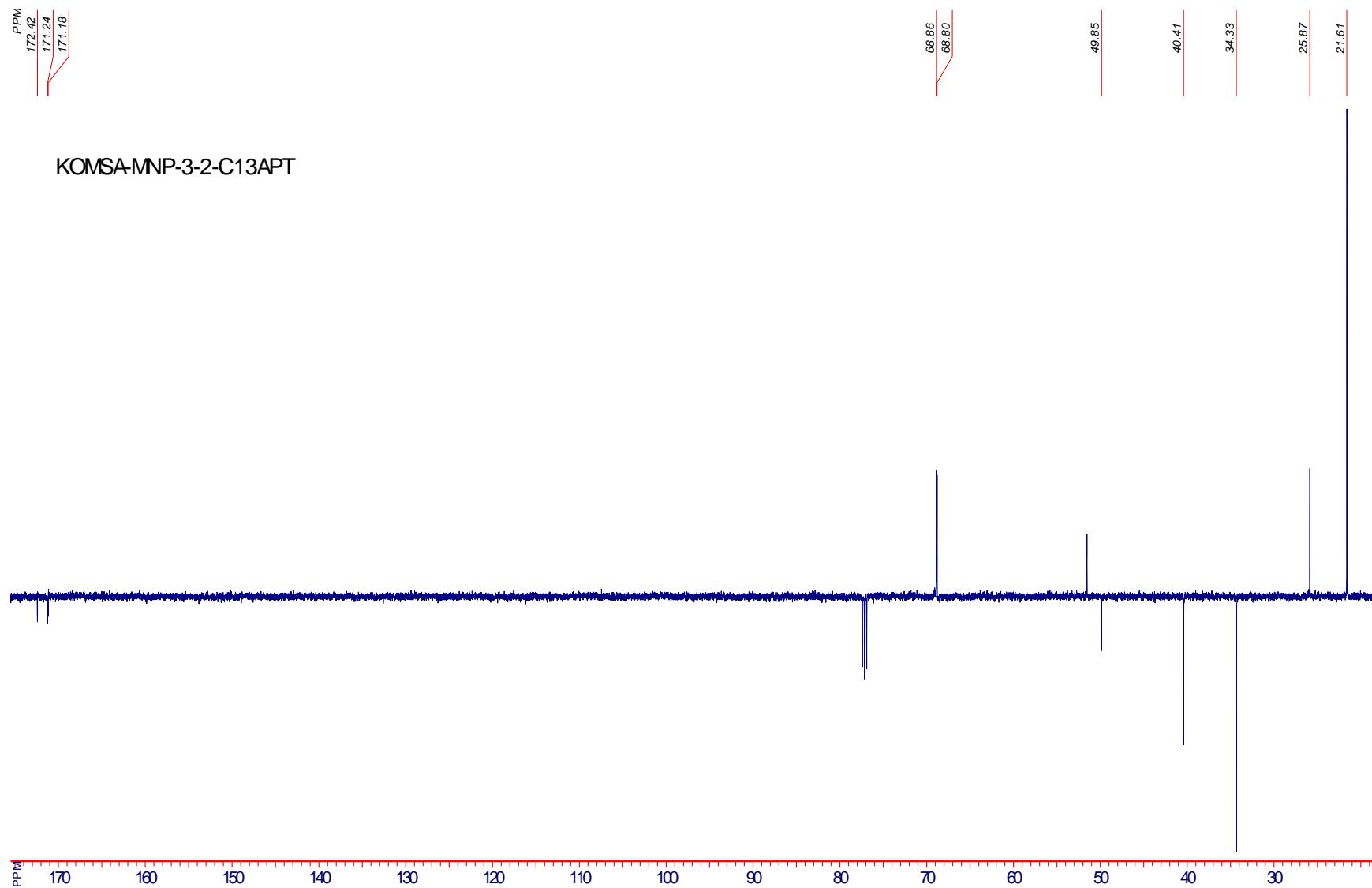


PPM

^1H NMR spectrum of the compound **37**.

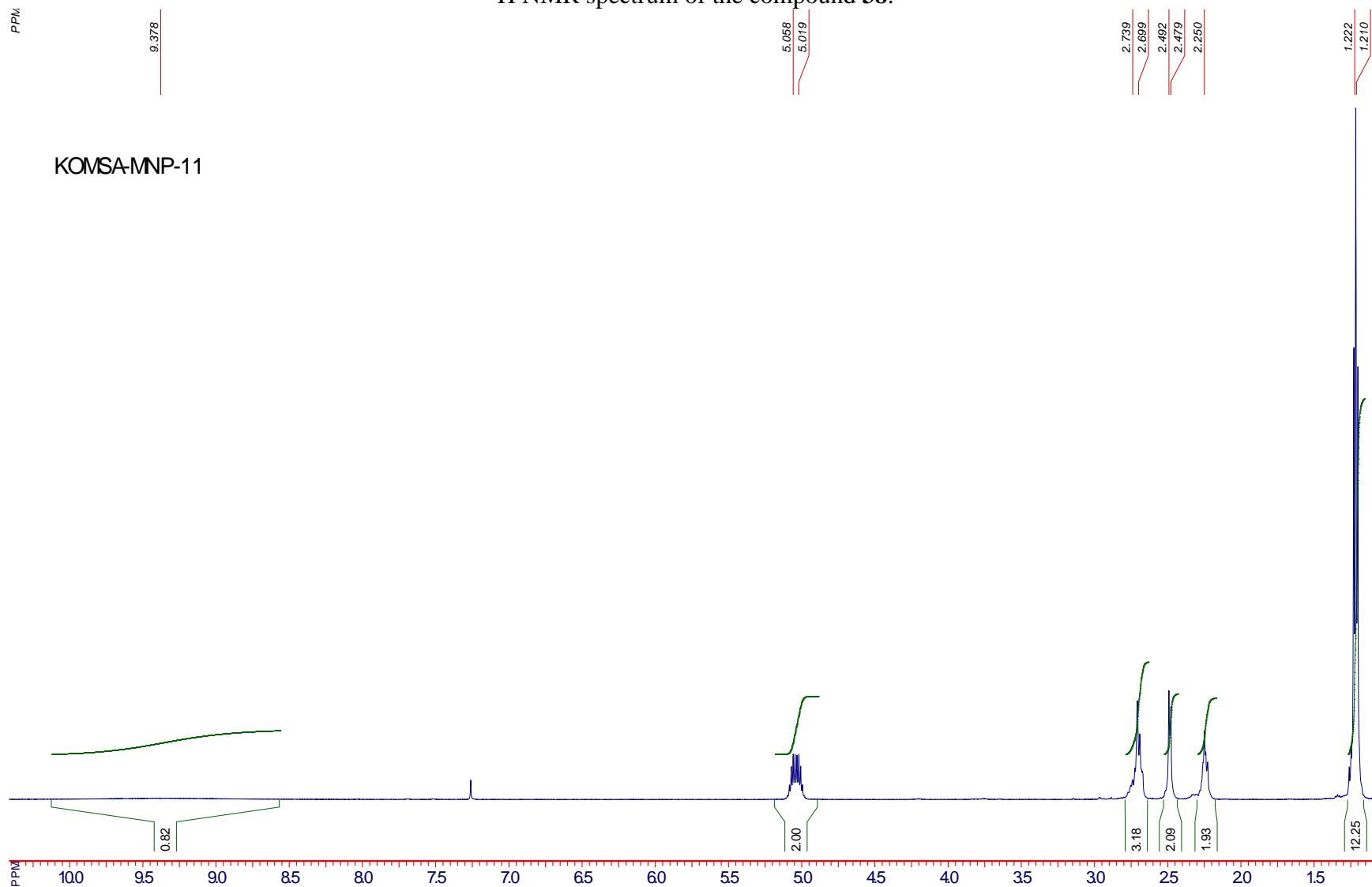


^{13}C NMR spectrum of the compound **37**.

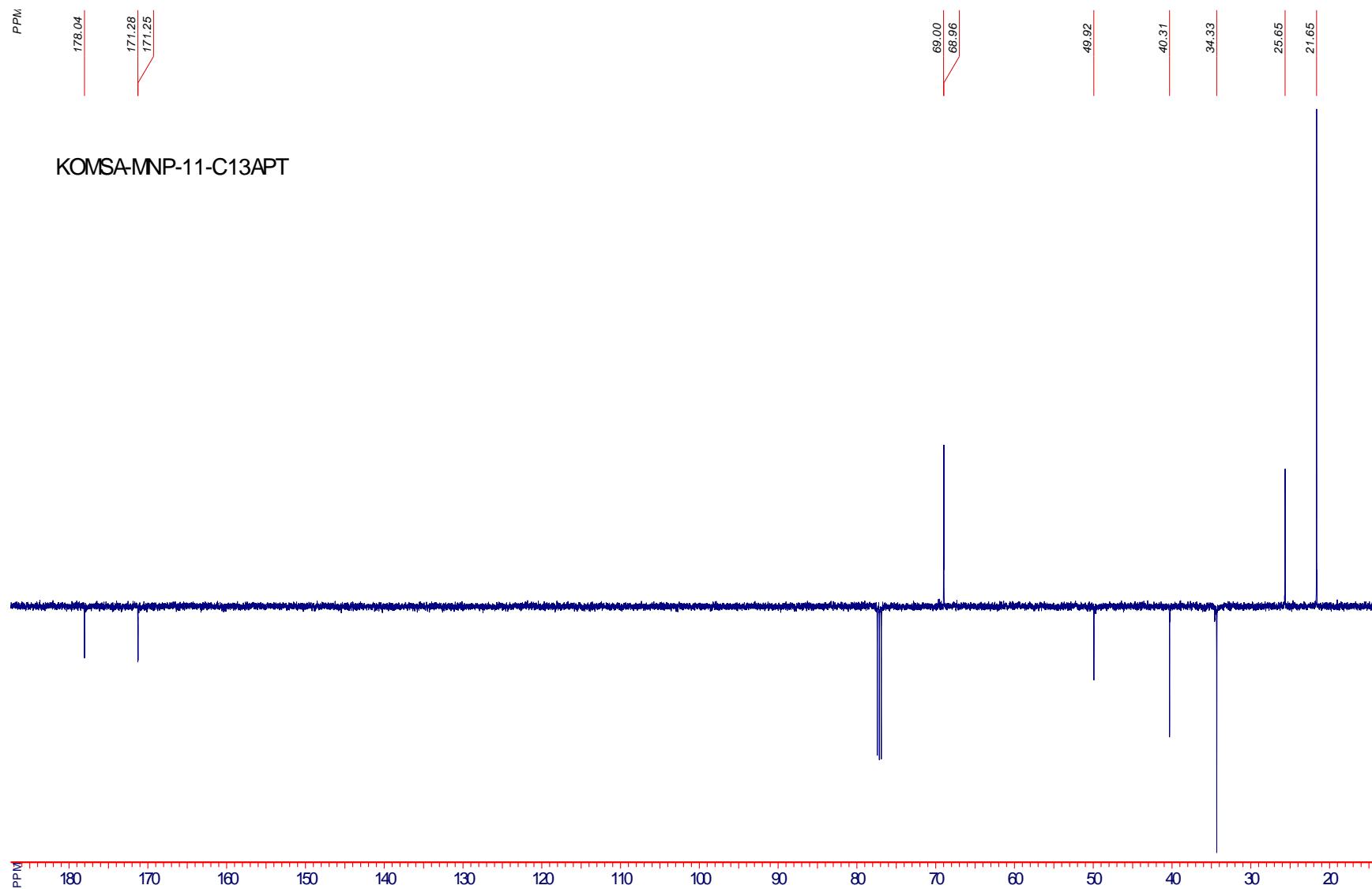


PPM

¹H NMR spectrum of the compound 38.

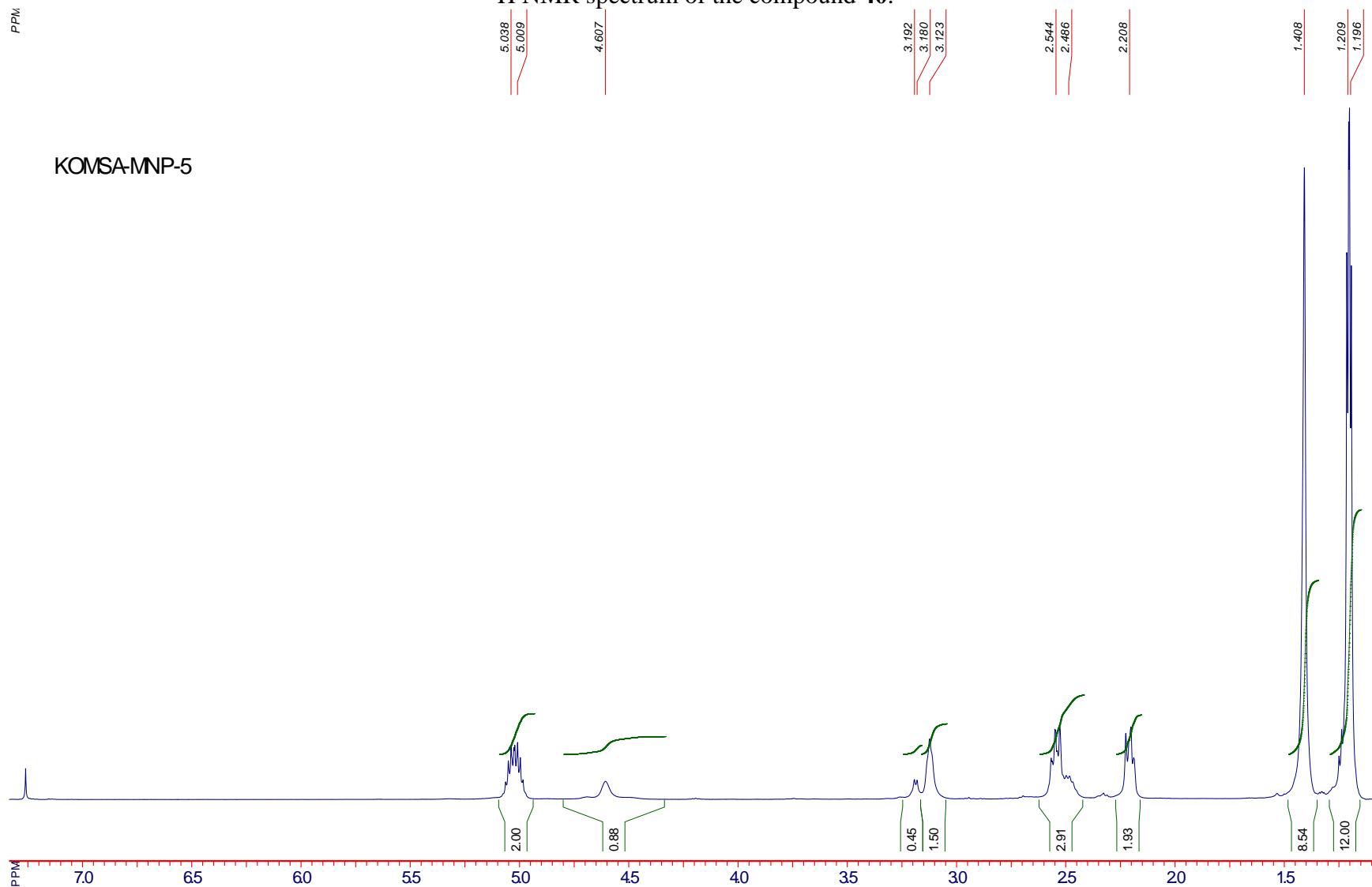


^{13}C NMR spectrum of the compound **38**.

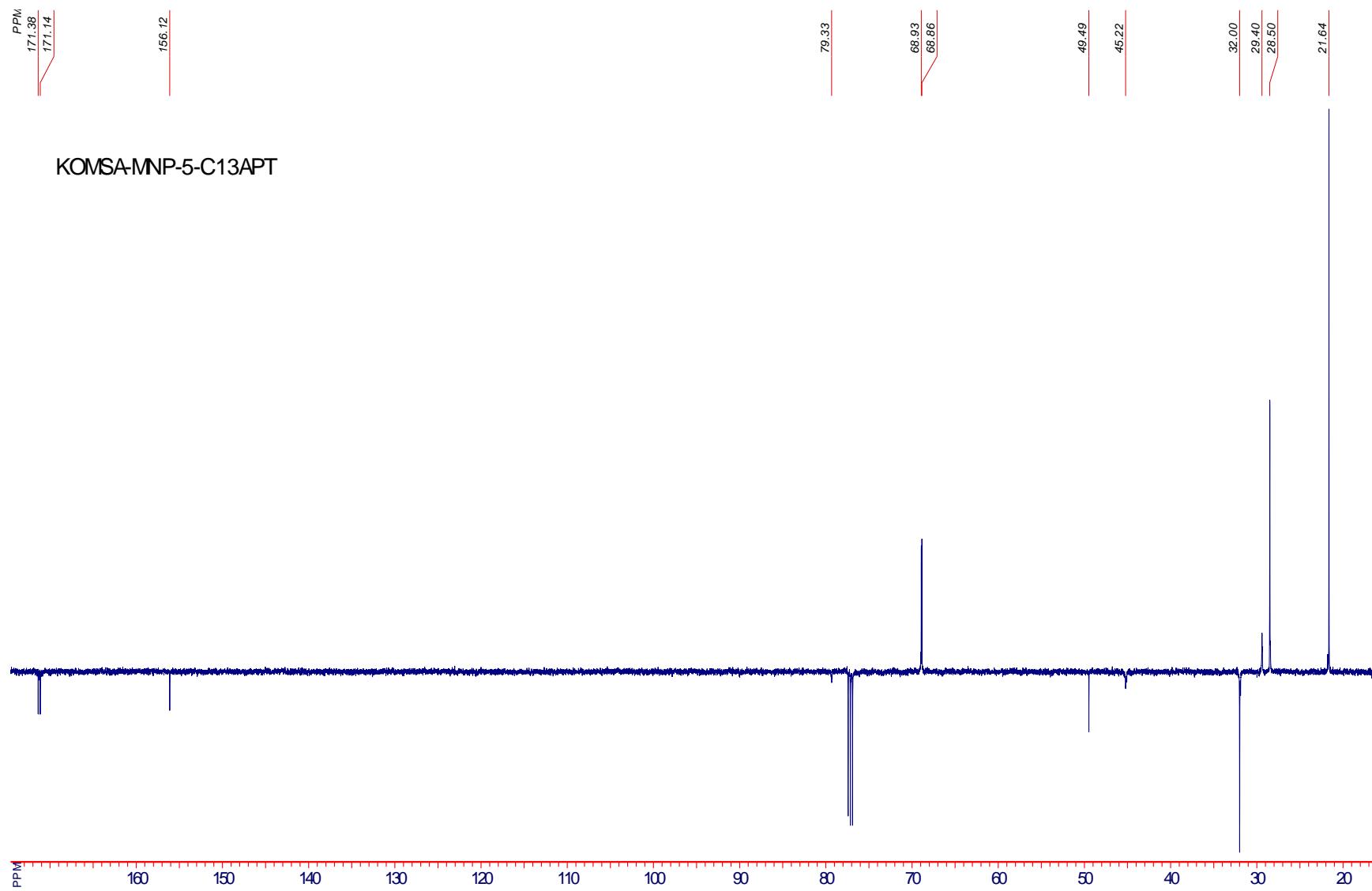


PPM

^1H NMR spectrum of the compound **40**.



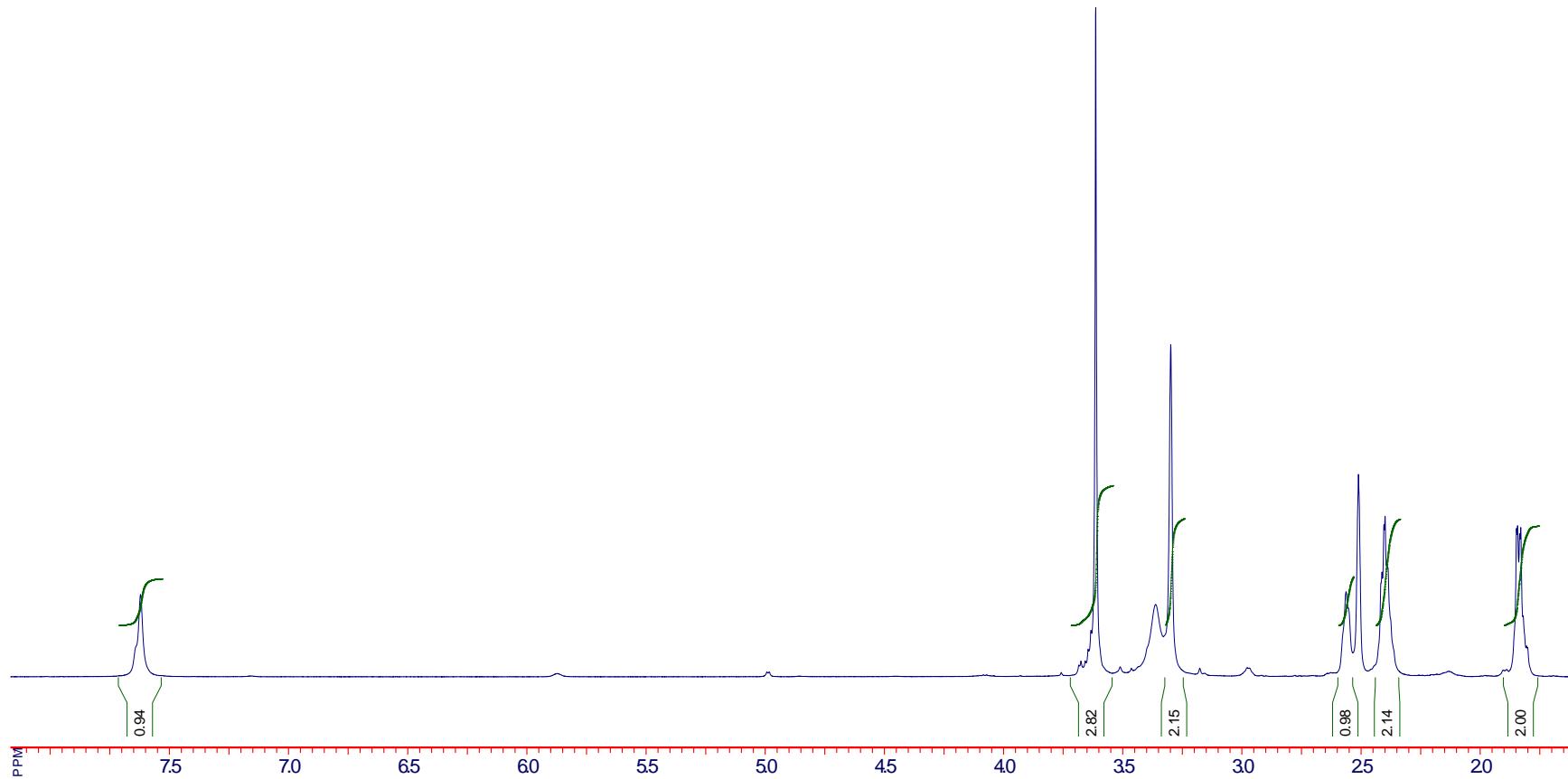
^{13}C NMR spectrum of the compound **40**.



PPM

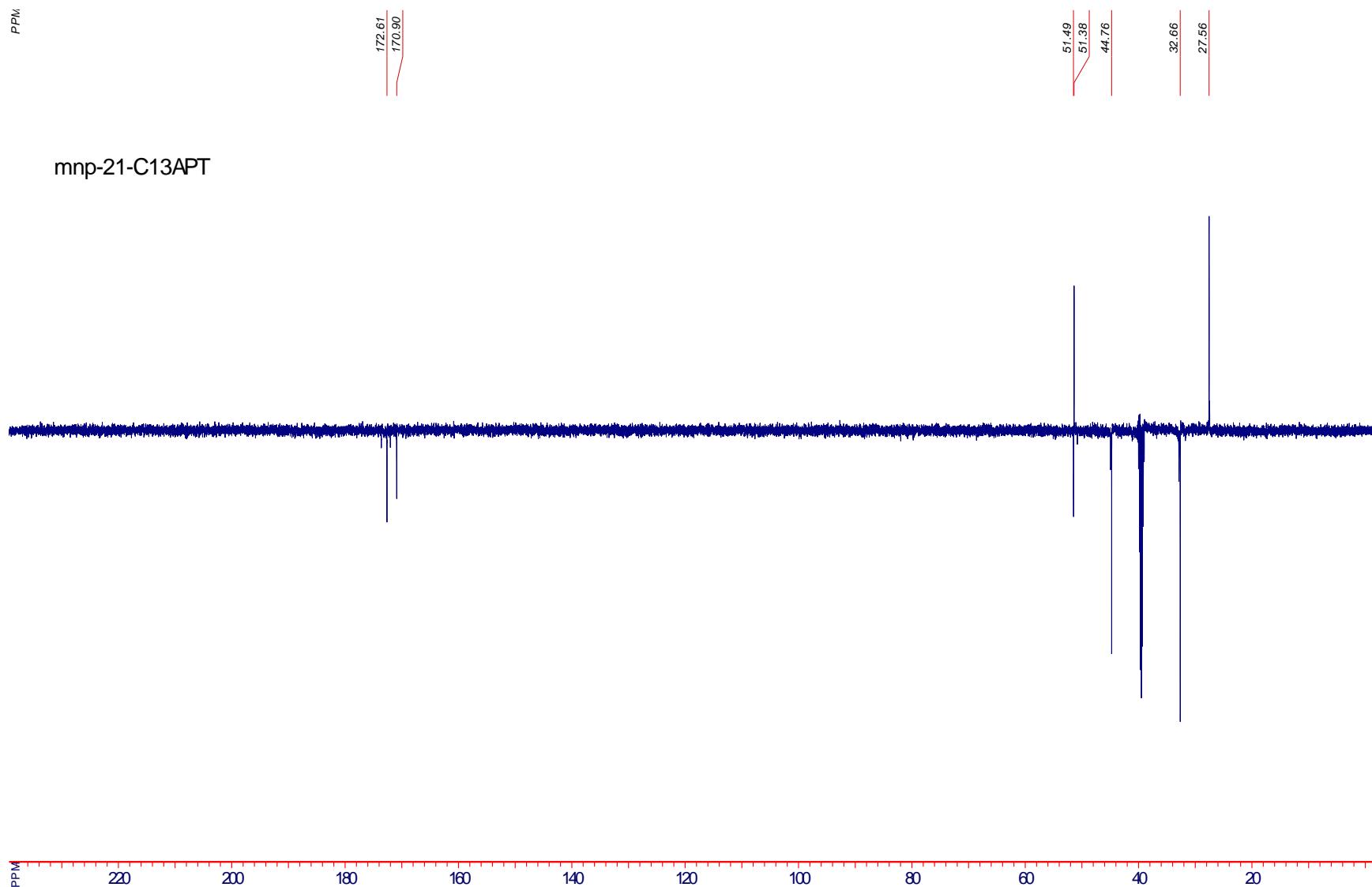
^1H NMR spectrum of the compound **41**.

mnp-21



^{13}C NMR spectrum of the compound **41**.

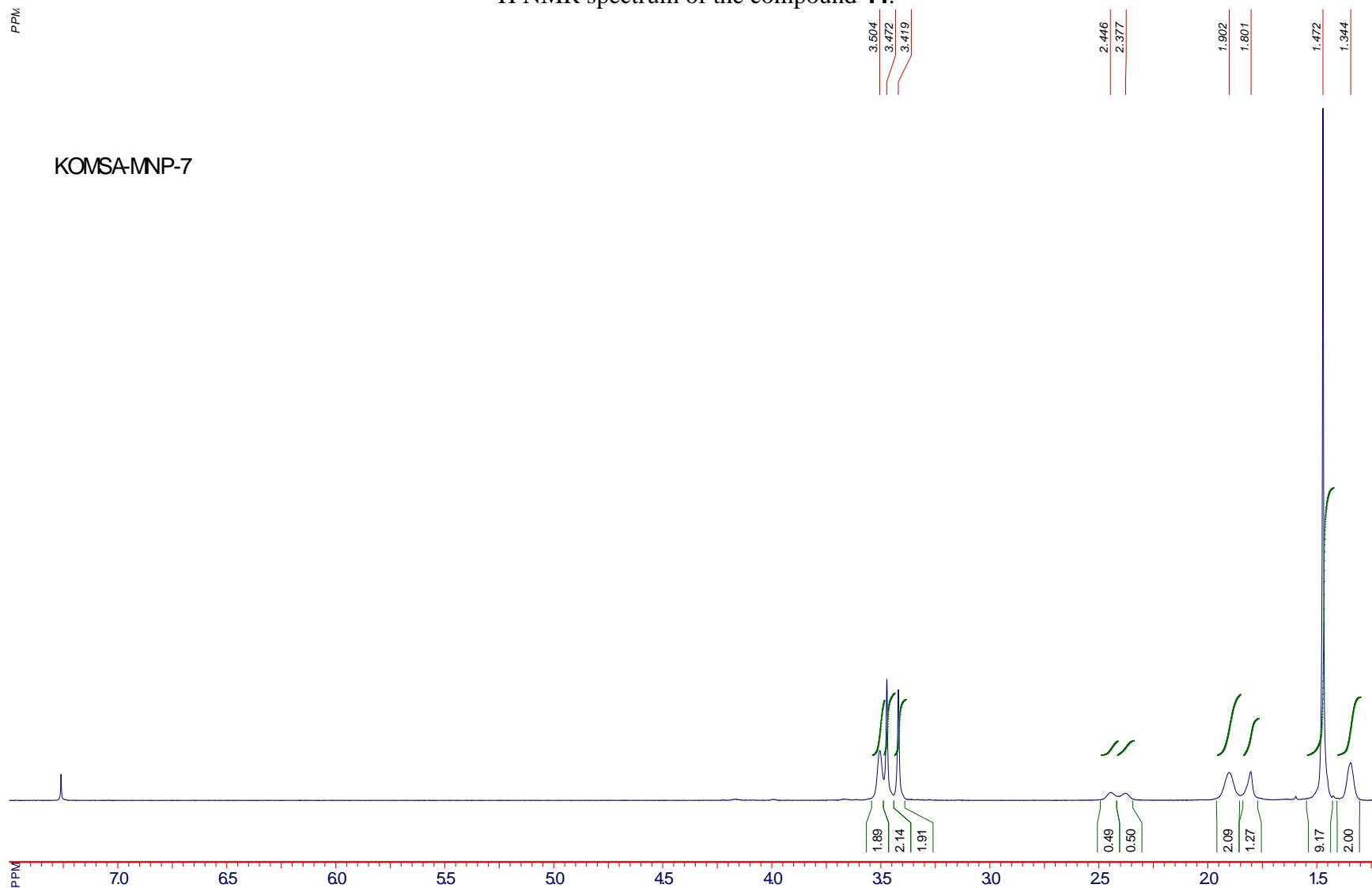
PPM



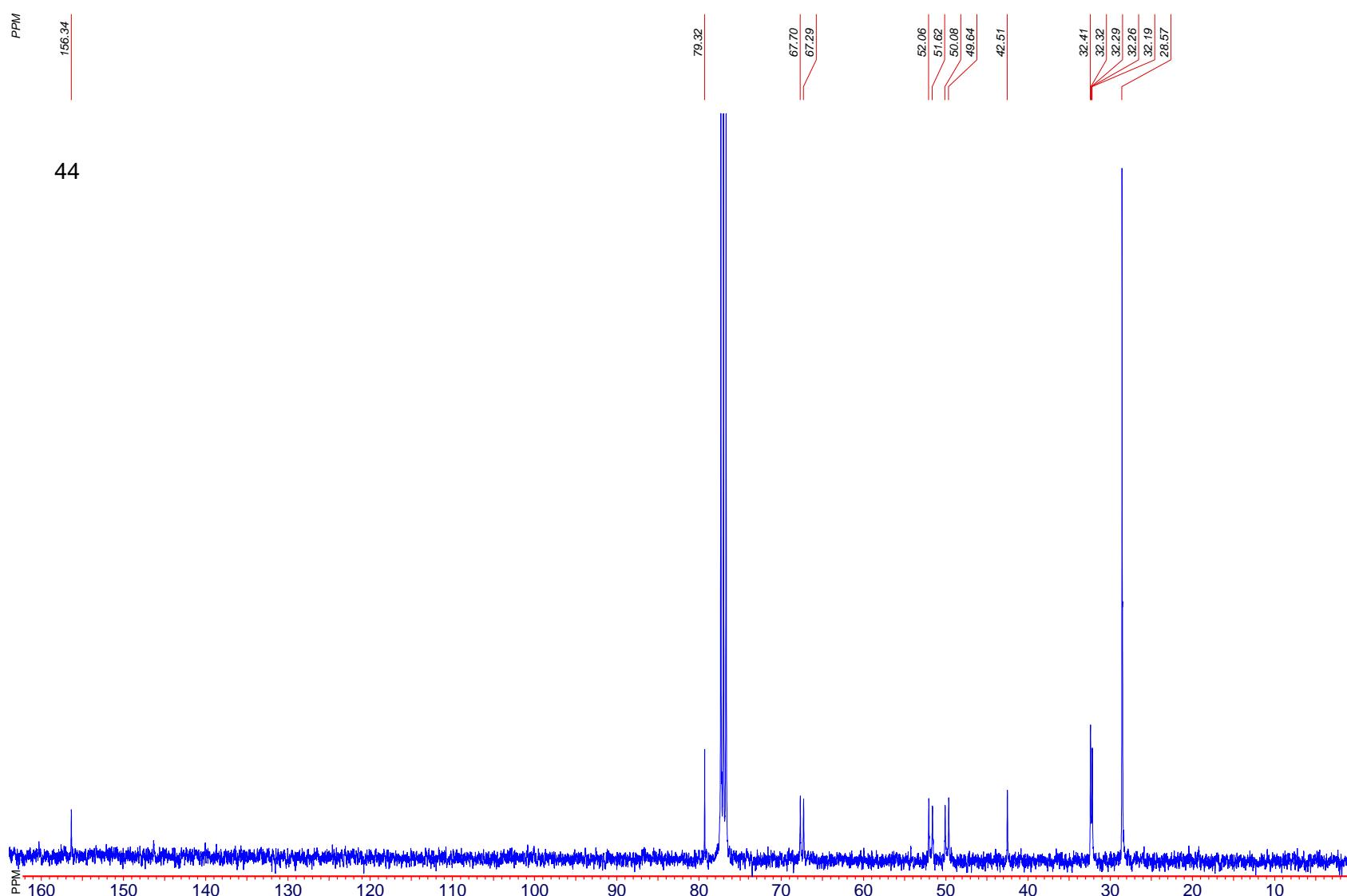
mnp-21-C13APT

PPM

^1H NMR spectrum of the compound **44**.

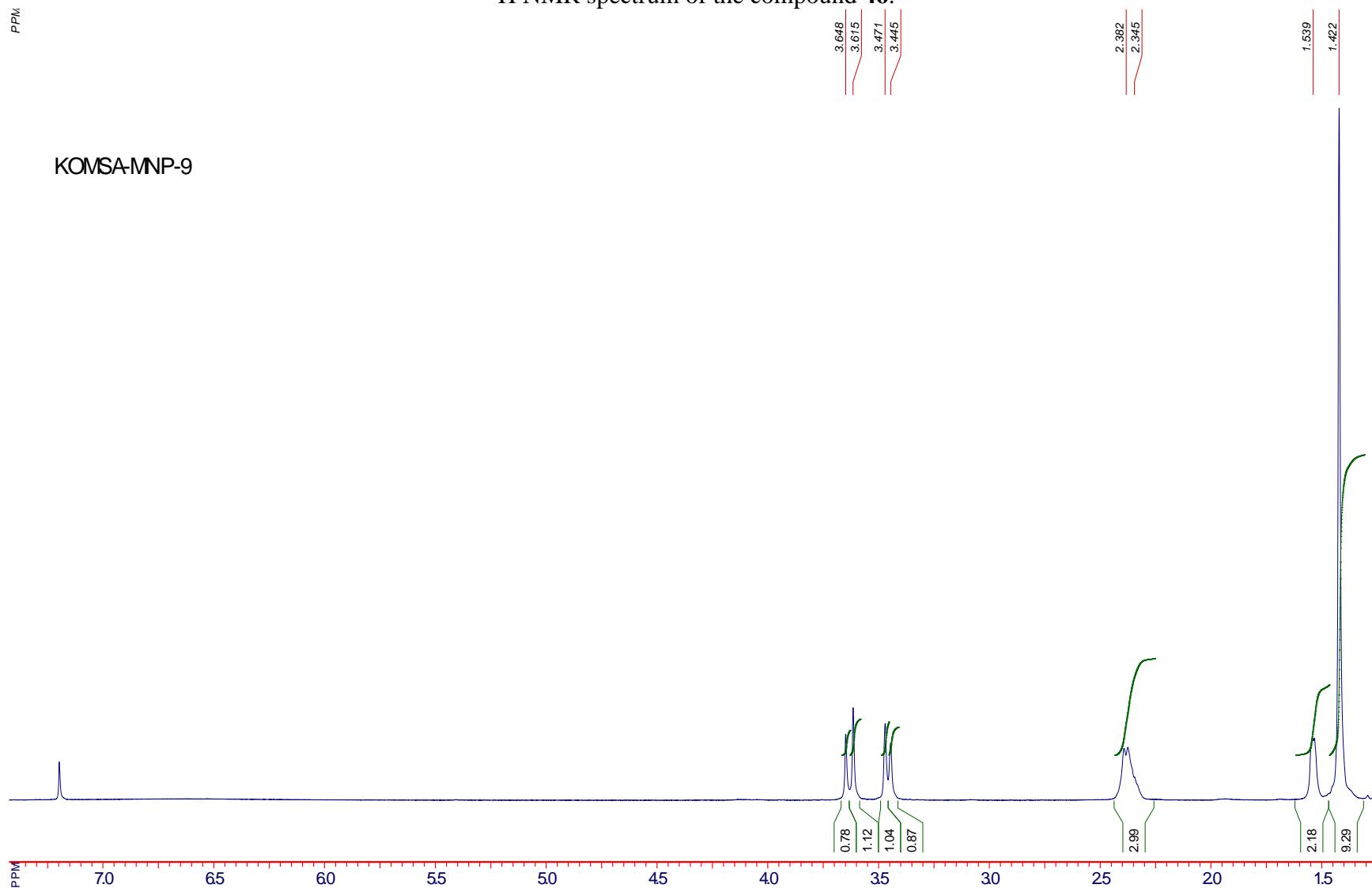


^{13}C NMR spectrum of the compound **44**.

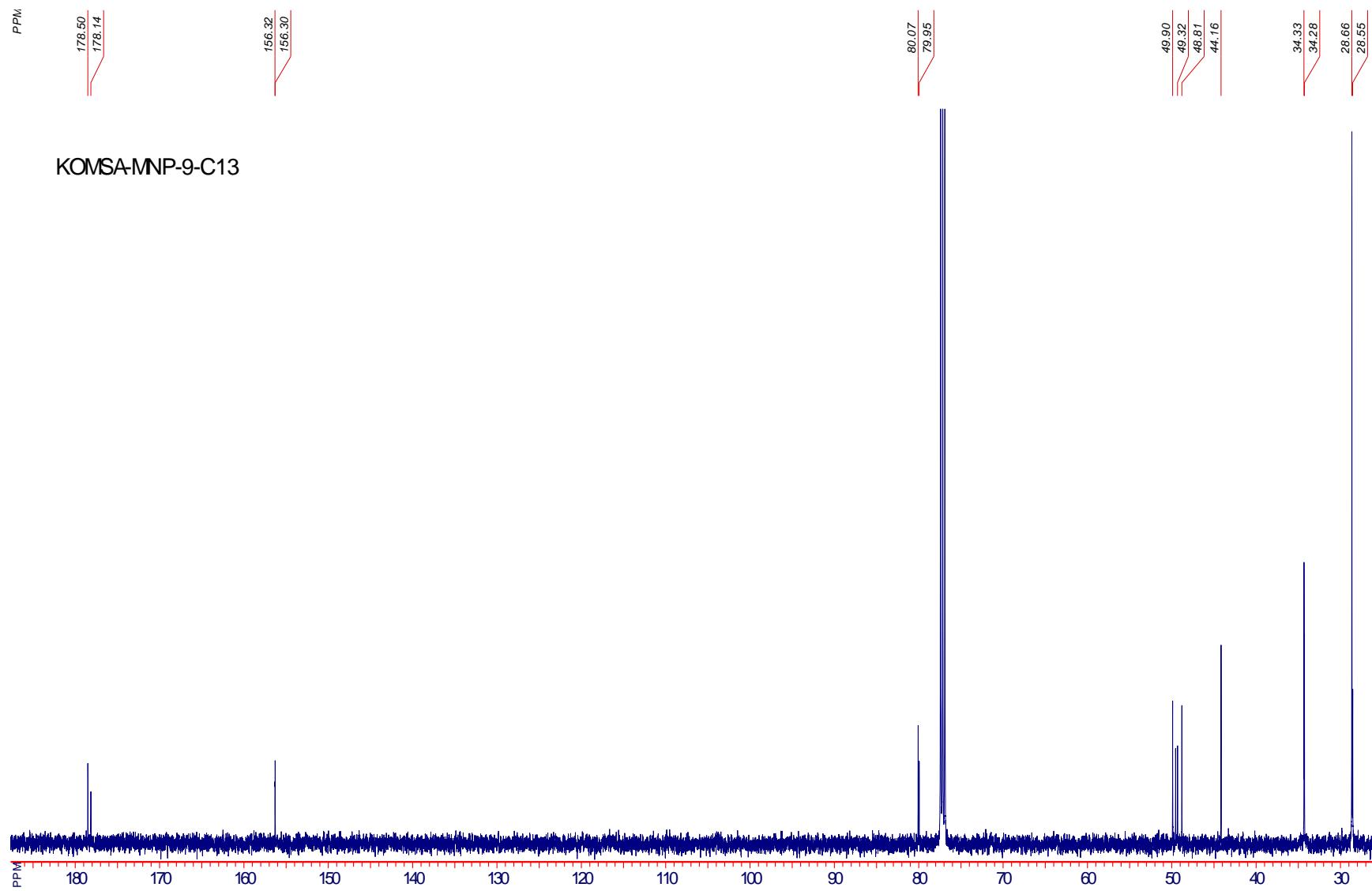


PPM

^1H NMR spectrum of the compound **46**.

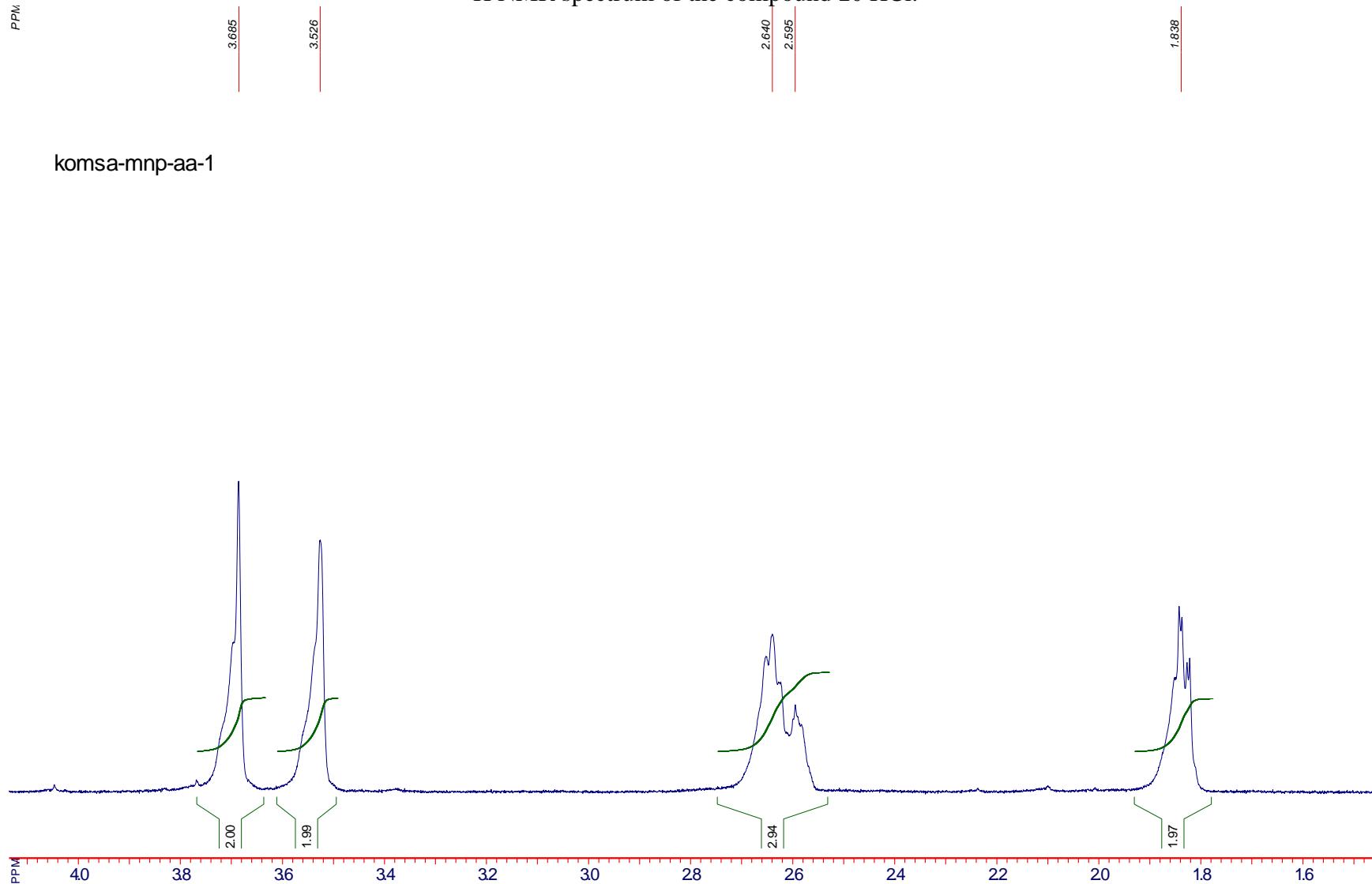


^{13}C NMR spectrum of the compound **46**.

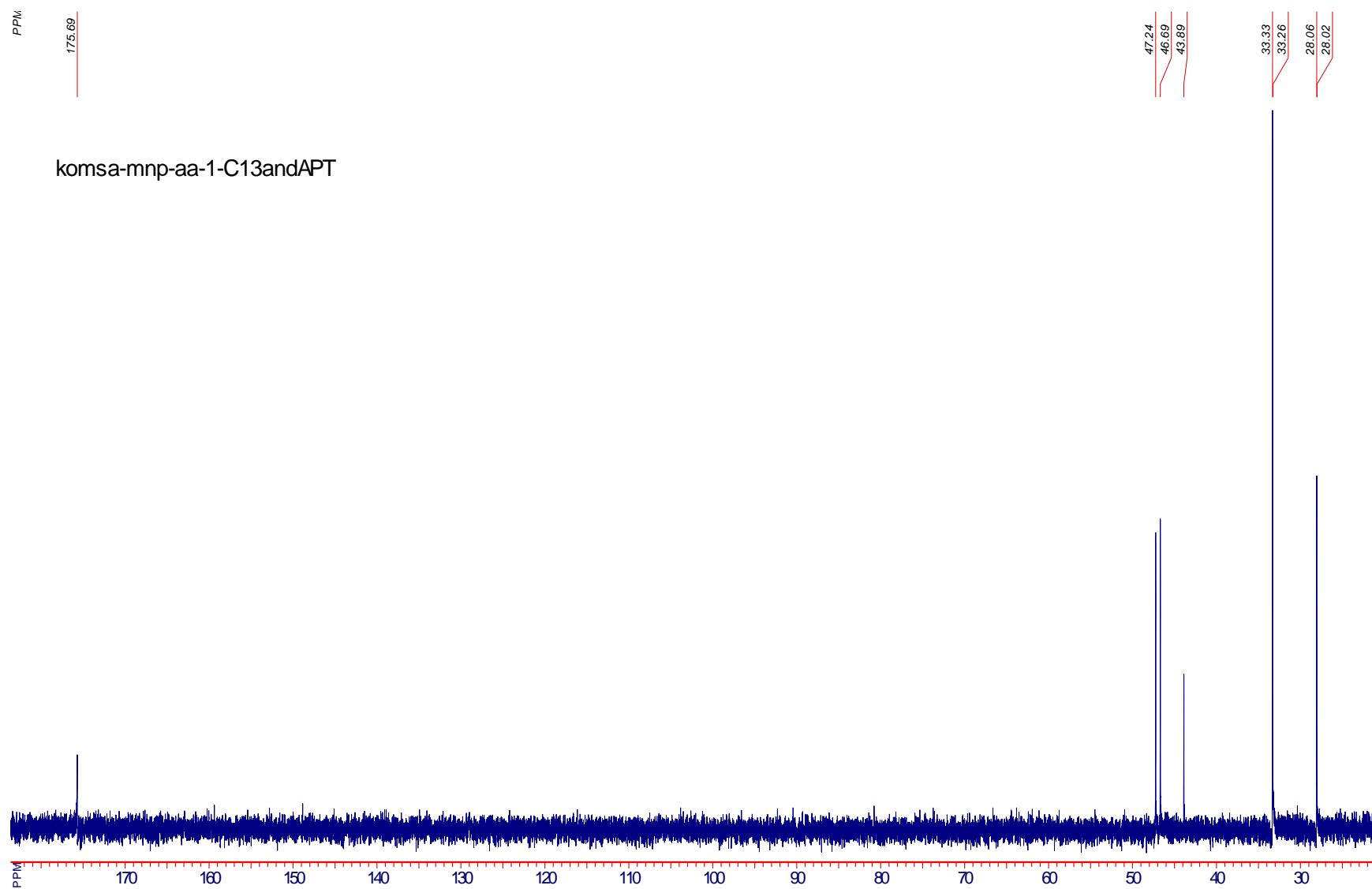


PPM

¹H NMR spectrum of the compound **10**·HCl.



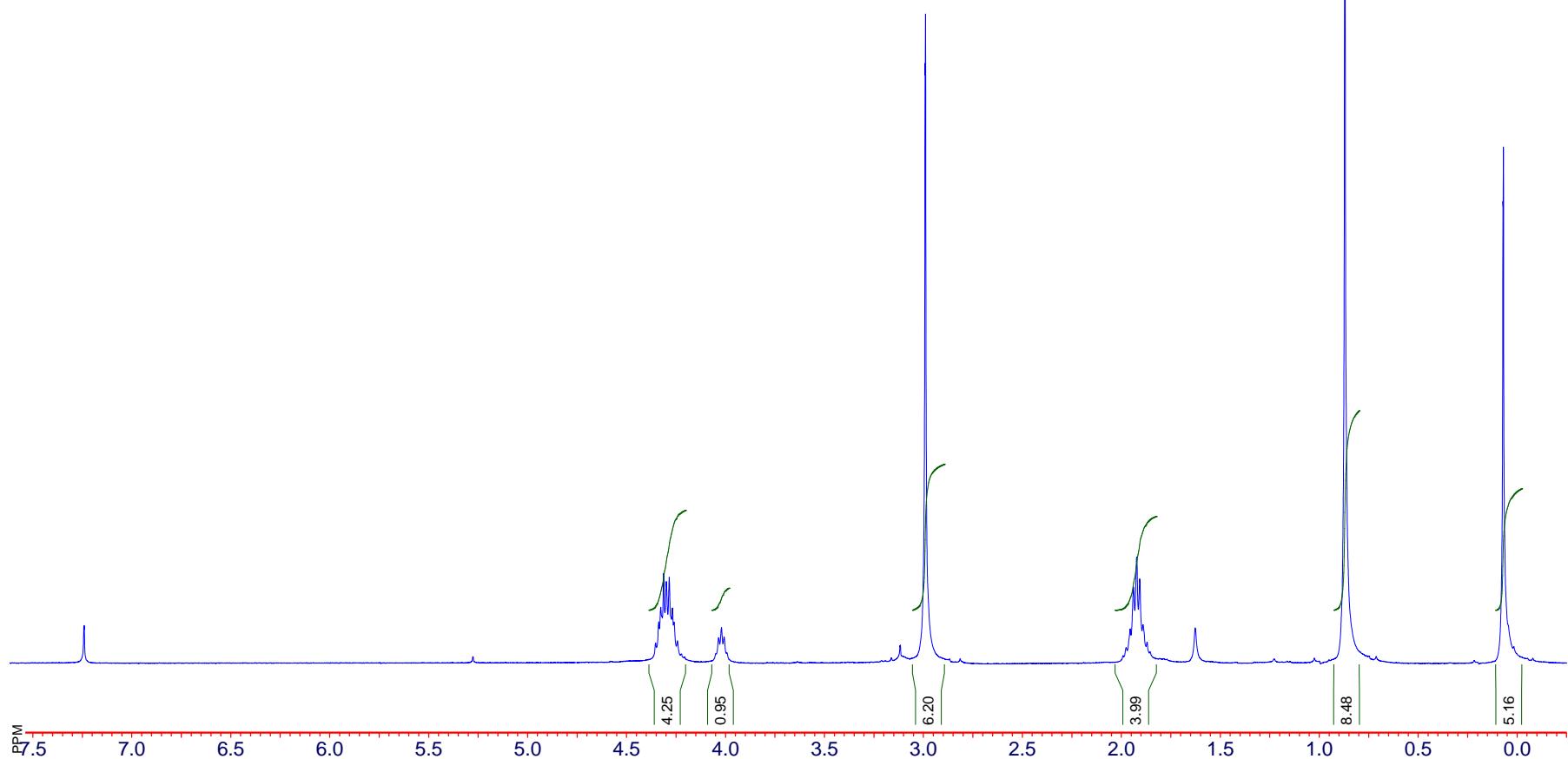
^{13}C NMR spectrum of the compound **10**·HCl.



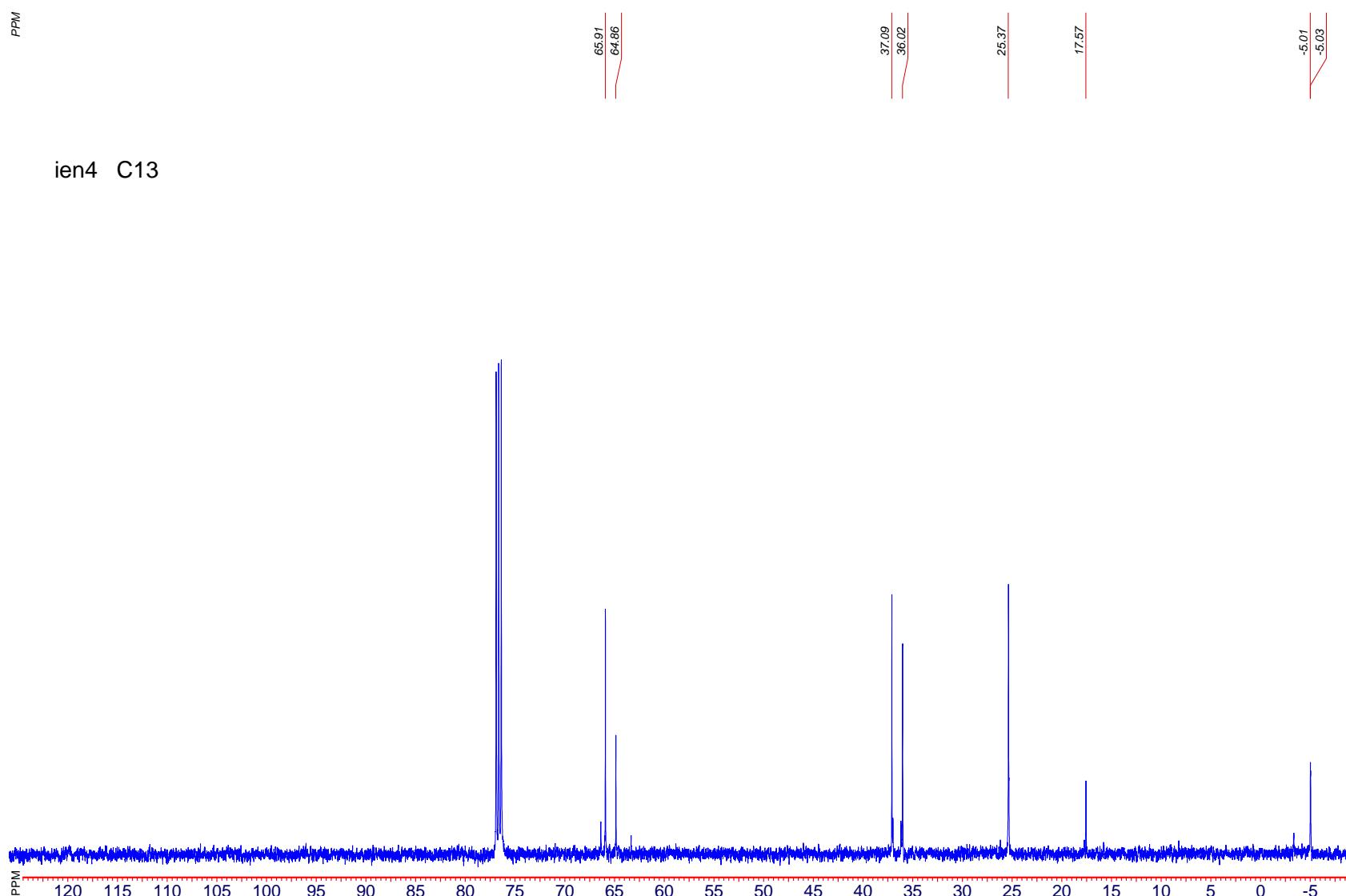
PPM

^1H NMR spectrum of the compound **23**.

ien-721-1

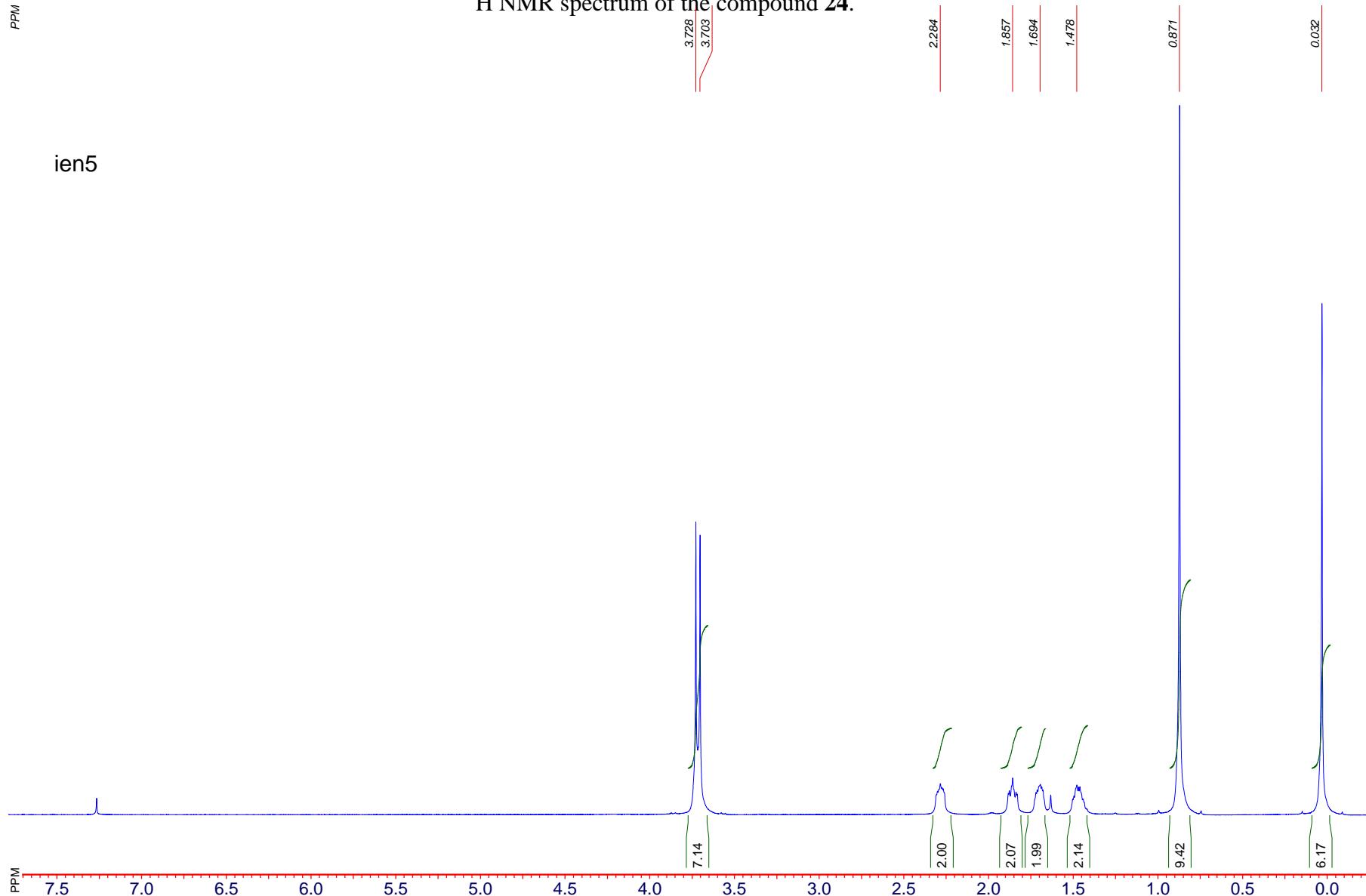


^{13}C NMR spectrum of the compound **23**.

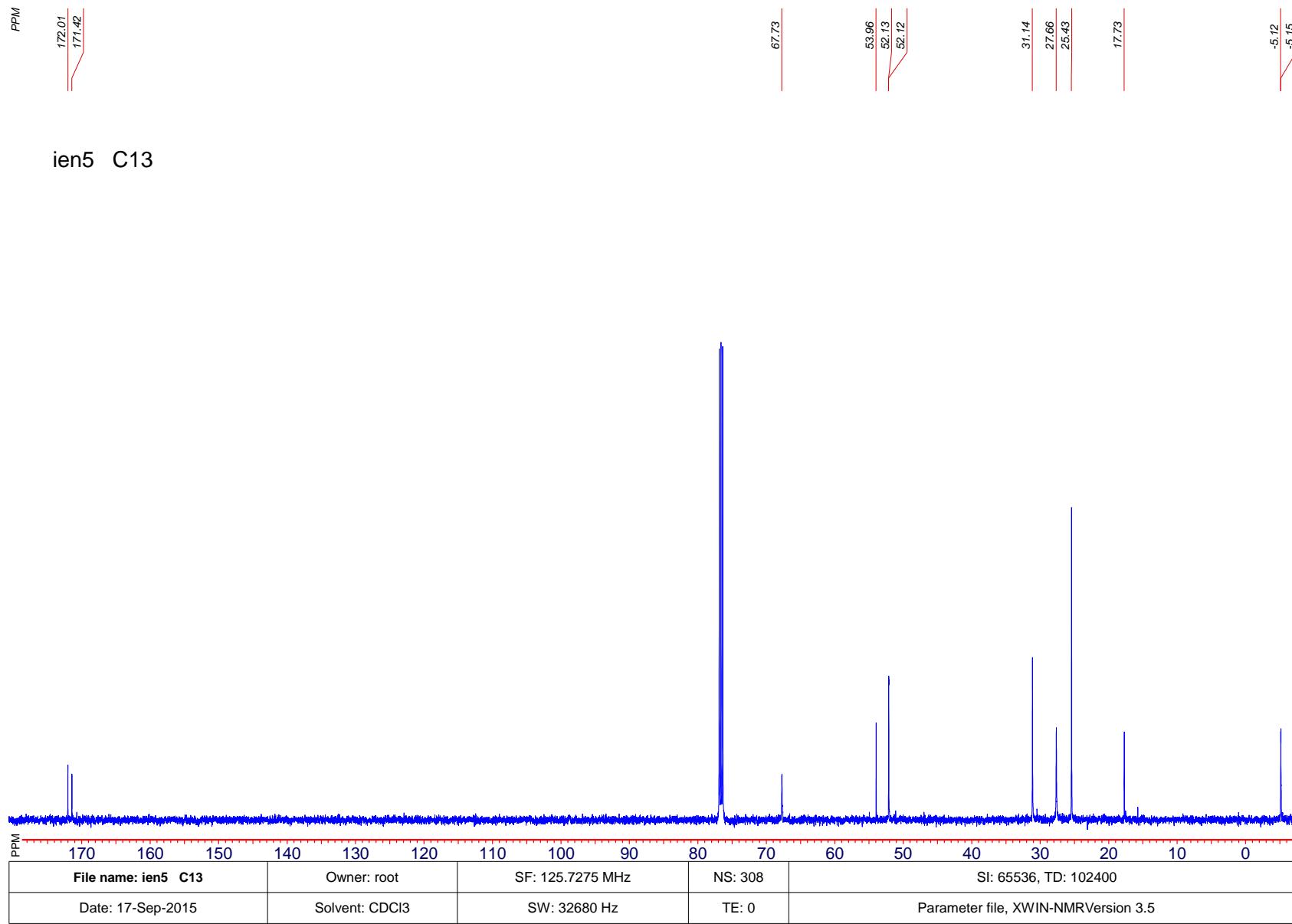


PPM

^1H NMR spectrum of the compound **24**.



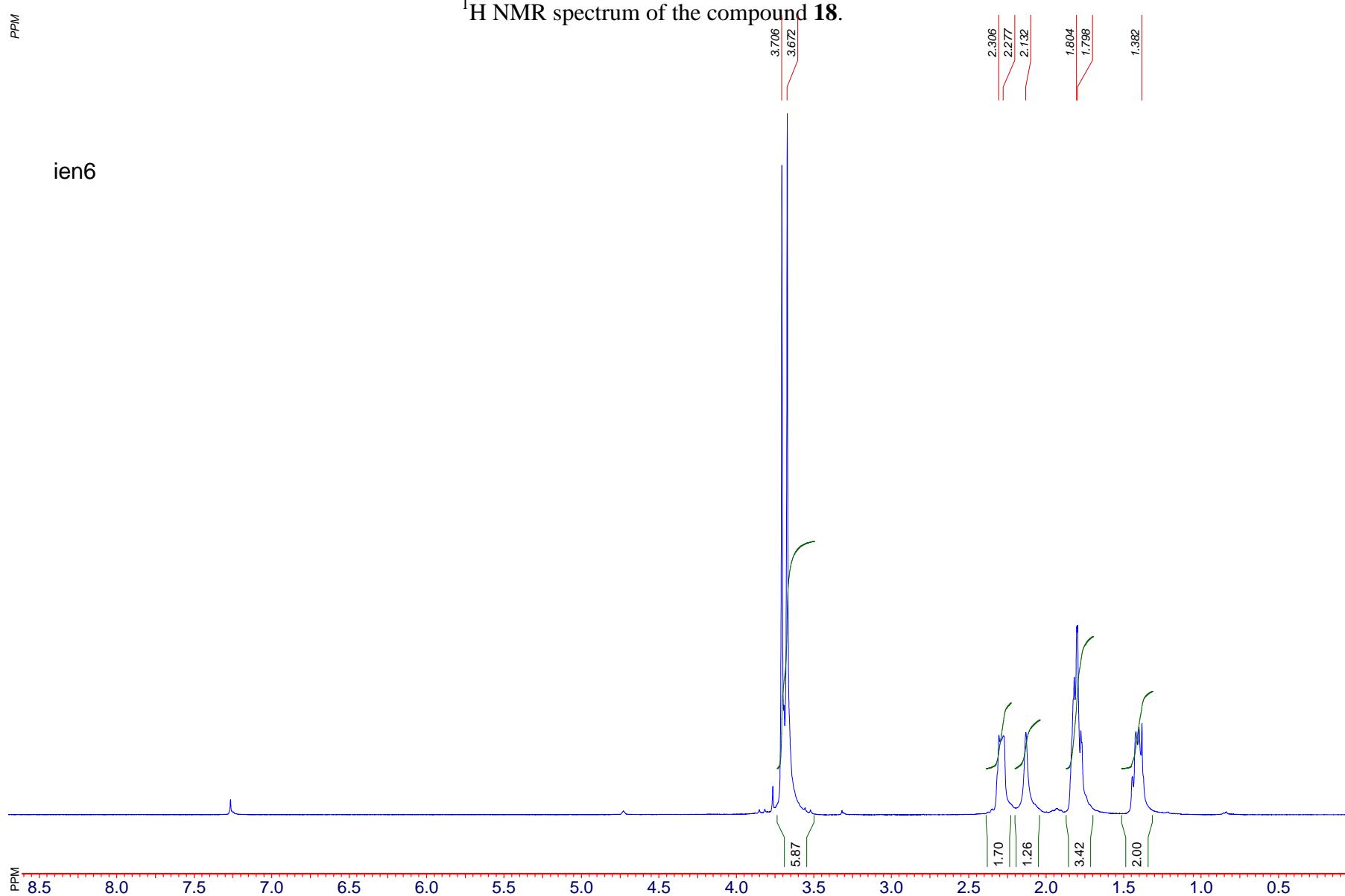
¹³C NMR spectrum of the compound **24**.



PPM

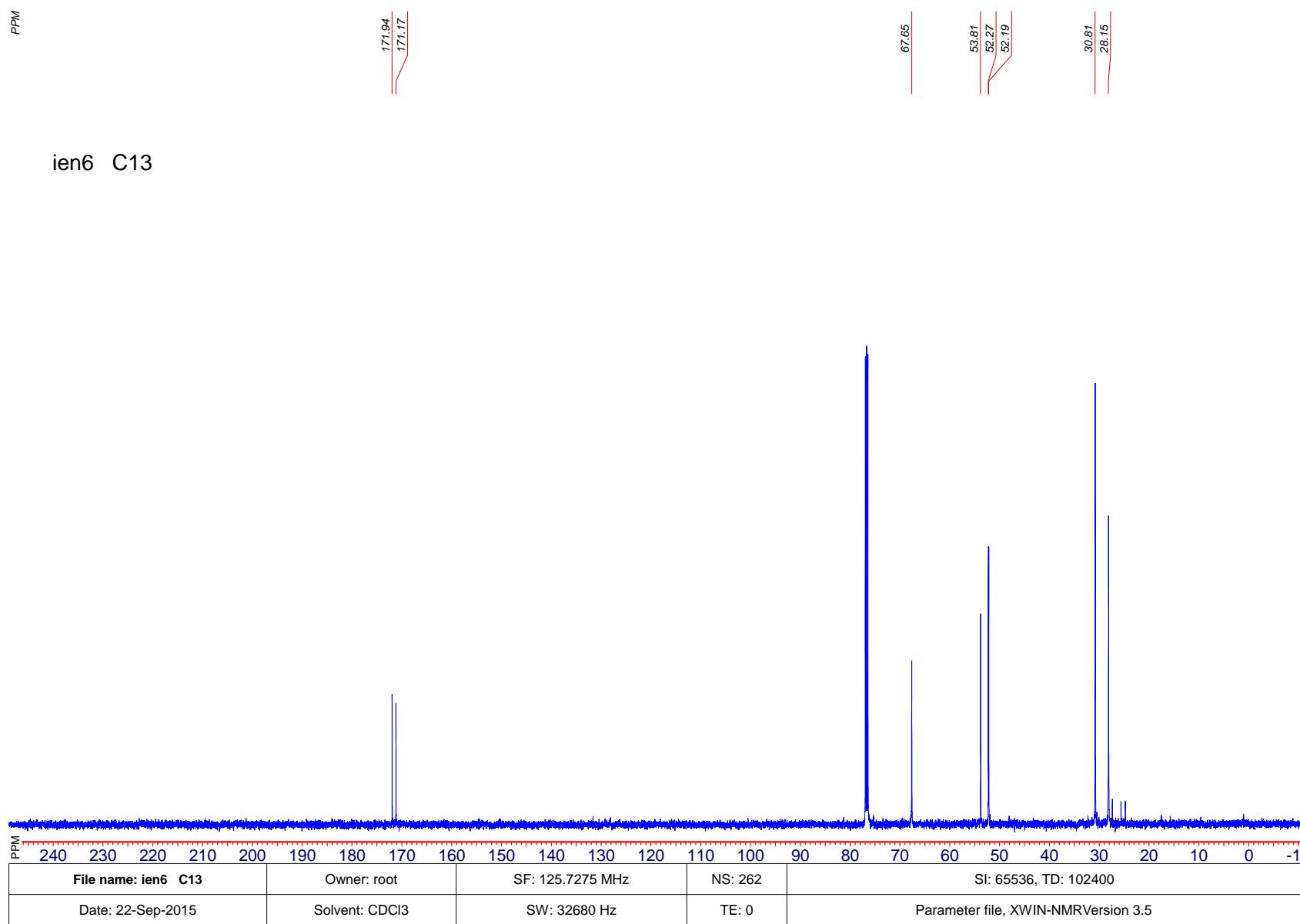
^1H NMR spectrum of the compound **18**.

ien6



S25

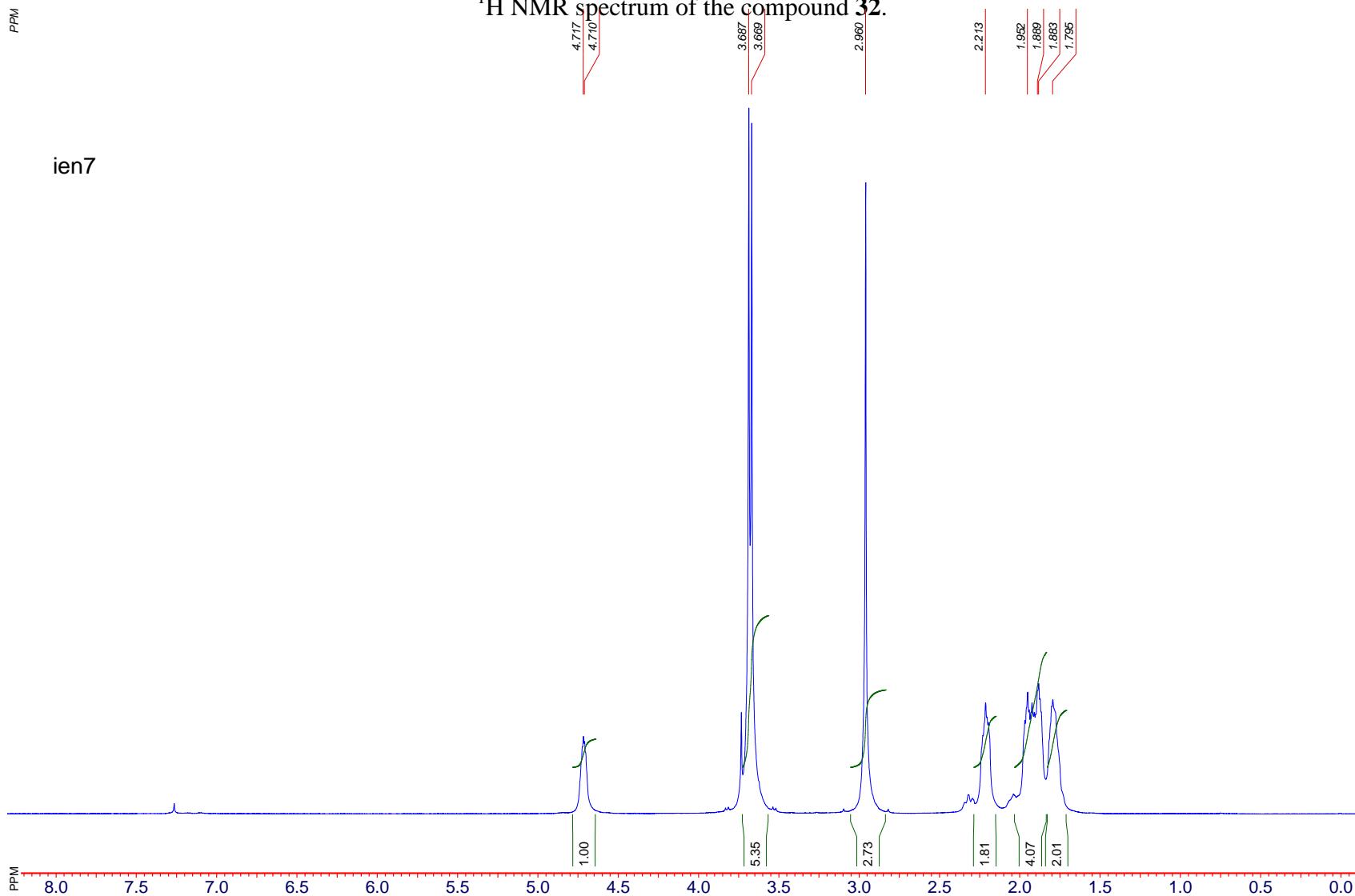
¹³C NMR spectrum of the compound **18**.



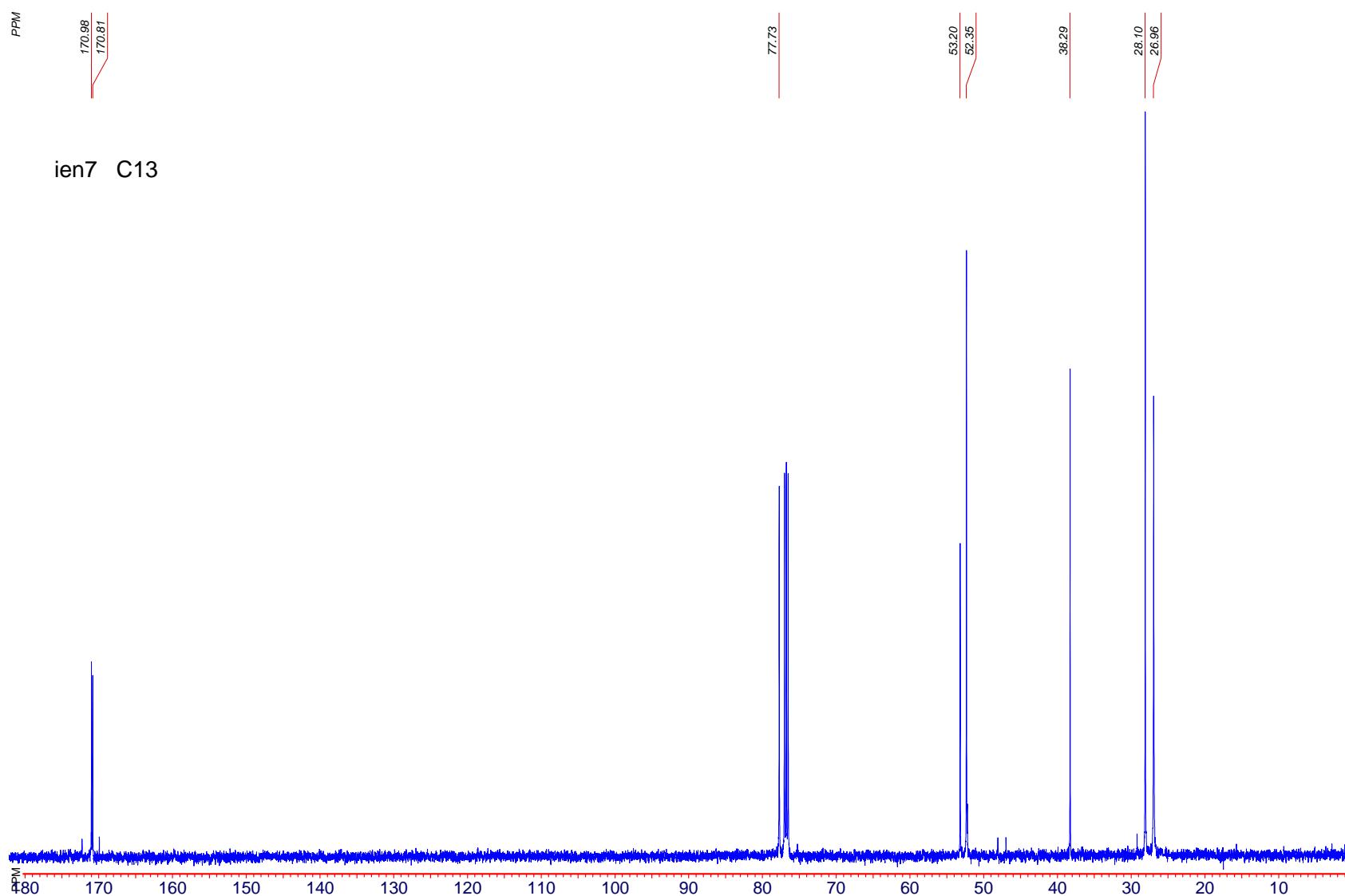
PPM

¹H NMR spectrum of the compound 32.

ien7



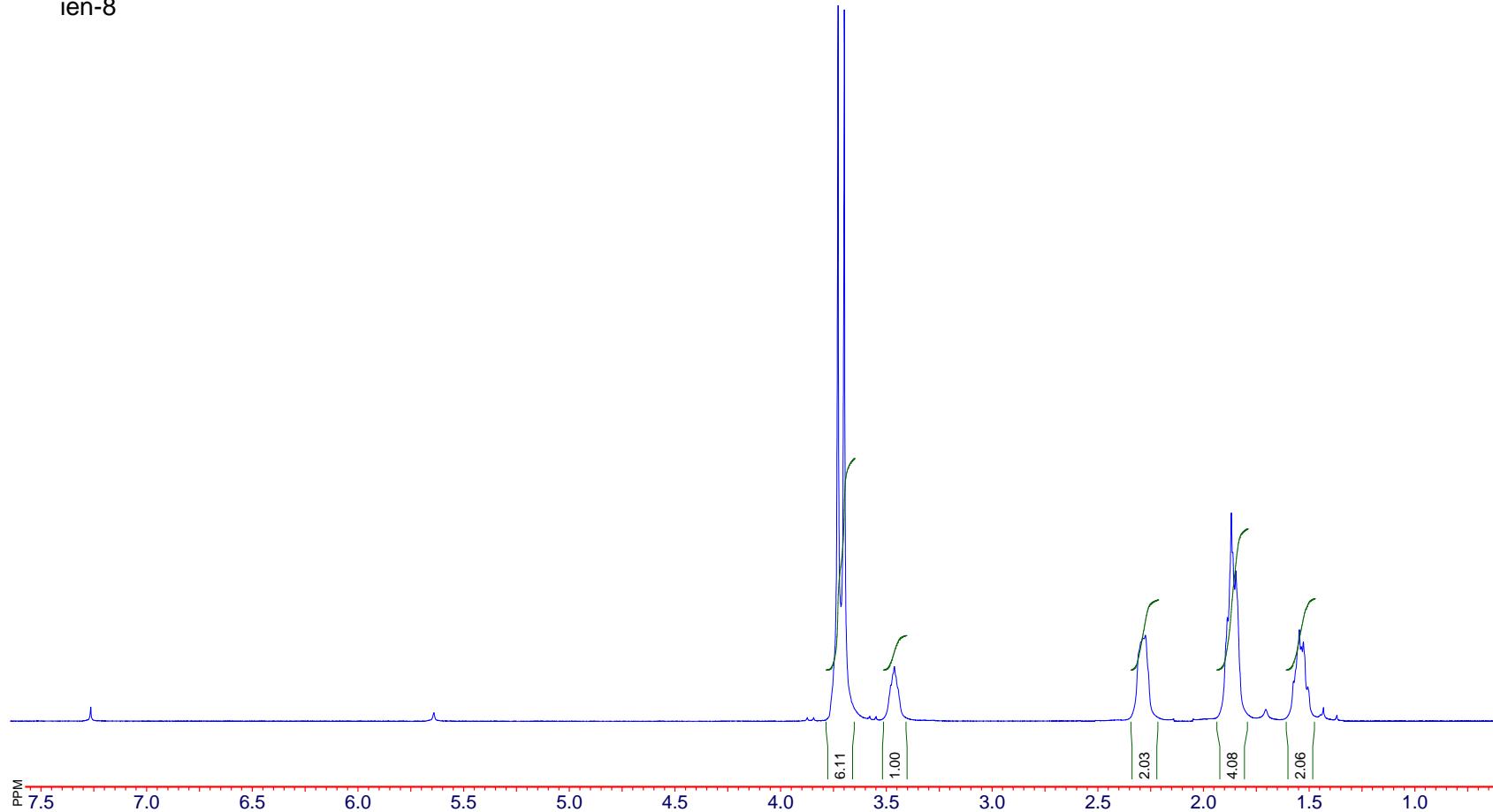
^{13}C NMR spectrum of the compound **32**.



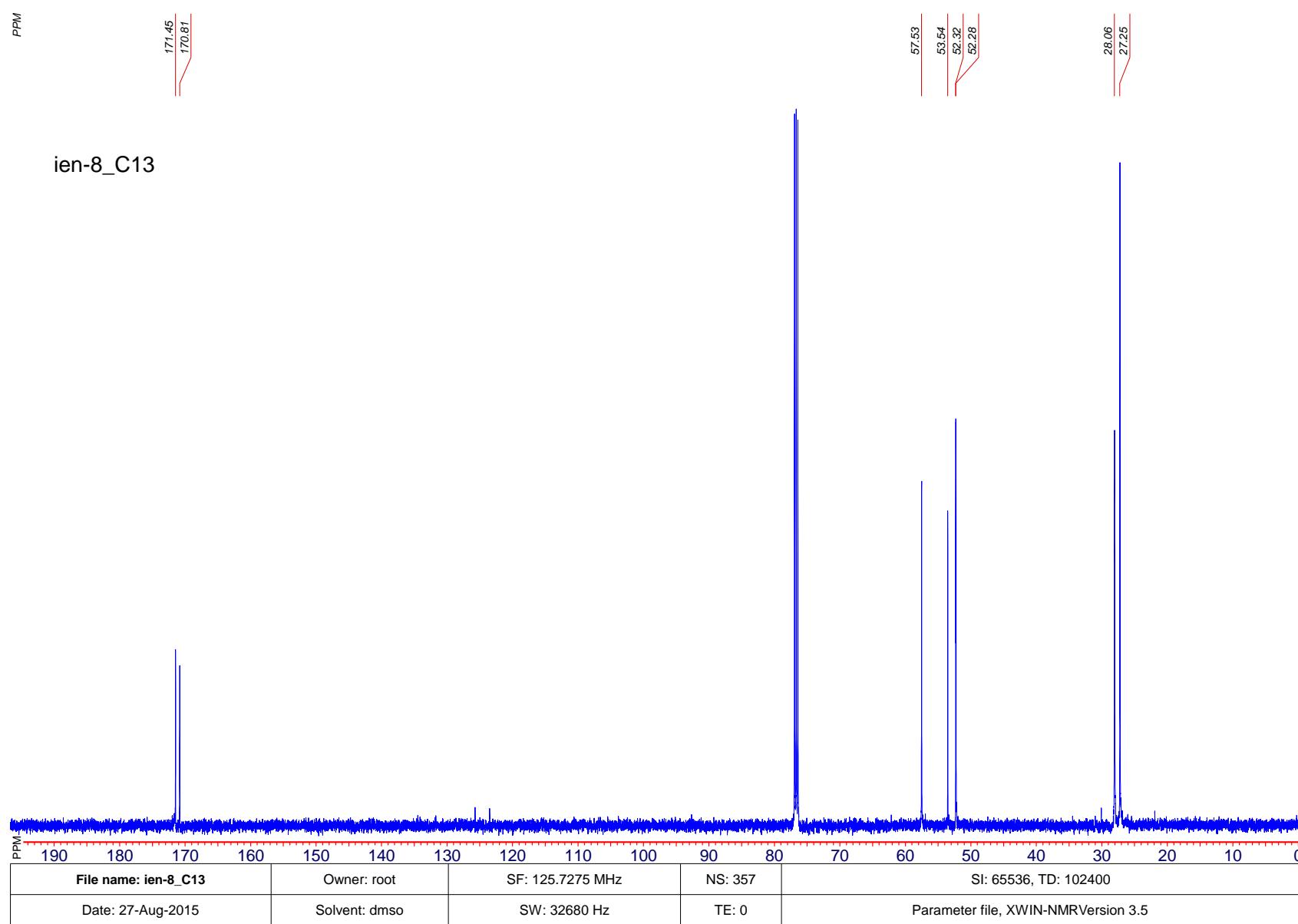
PPM

¹H NMR spectrum of the compound 34.

ien-8



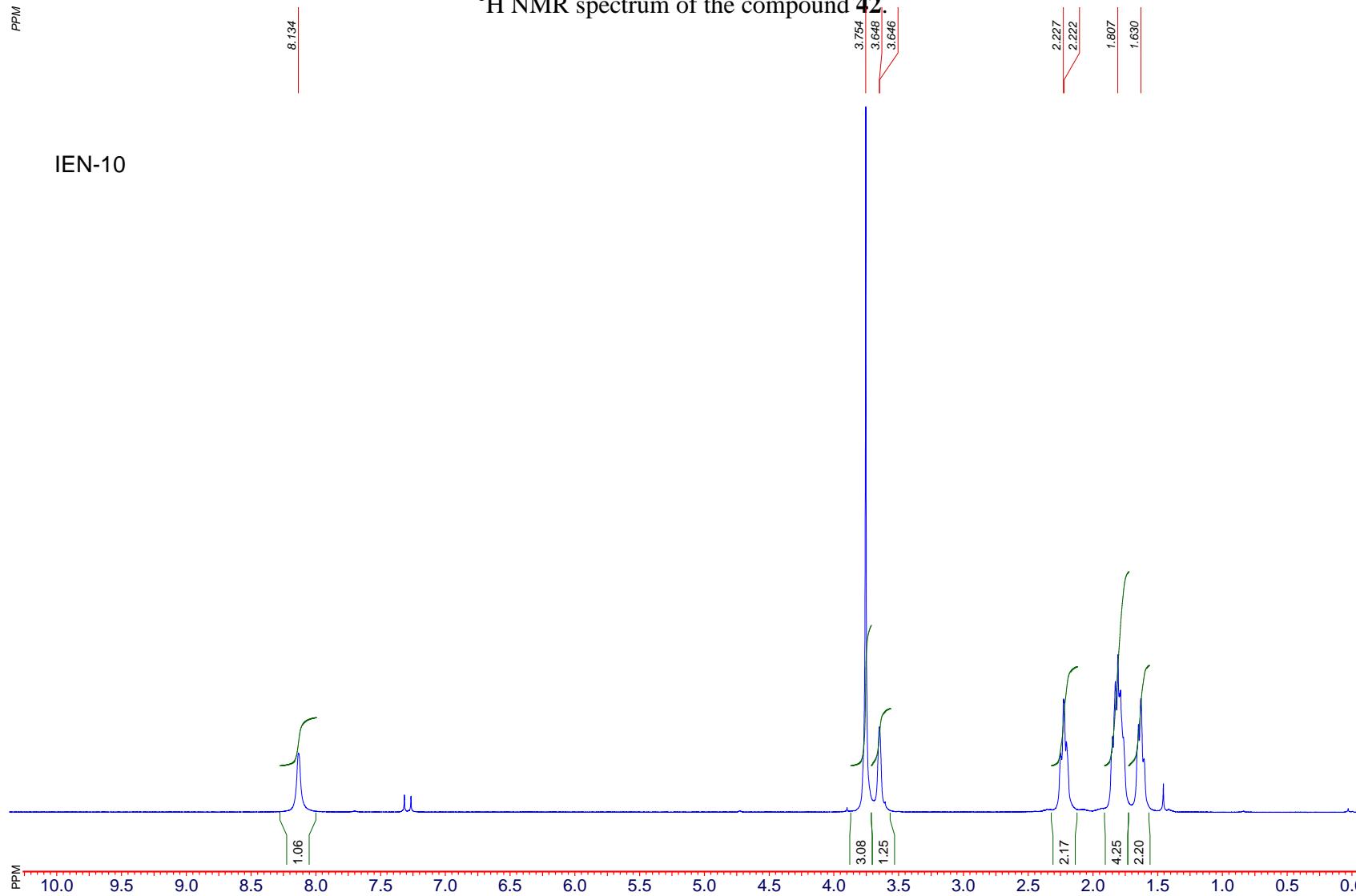
¹³C NMR spectrum of the compound **34**.



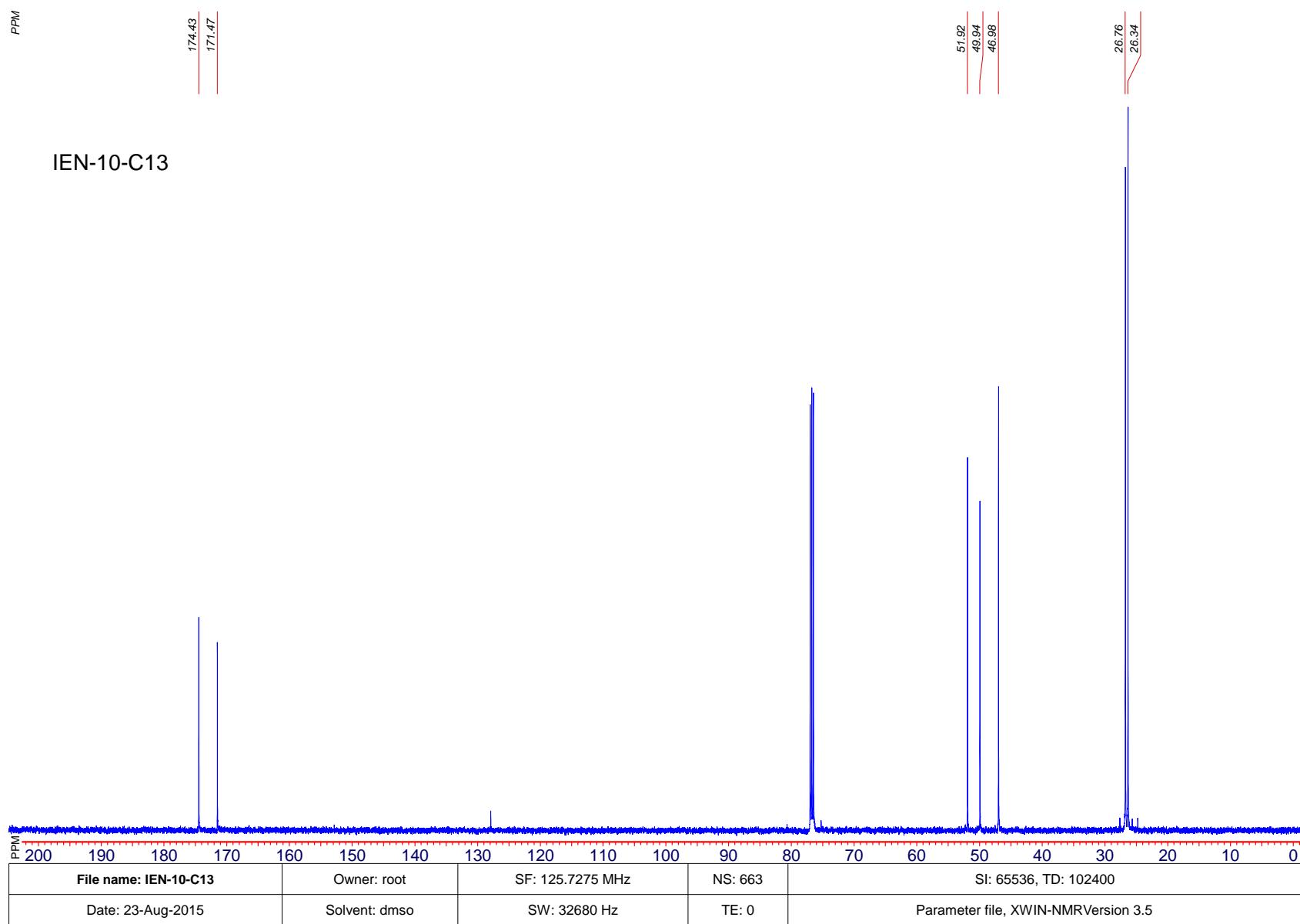
PPM

^1H NMR spectrum of the compound **42**.

IEN-10



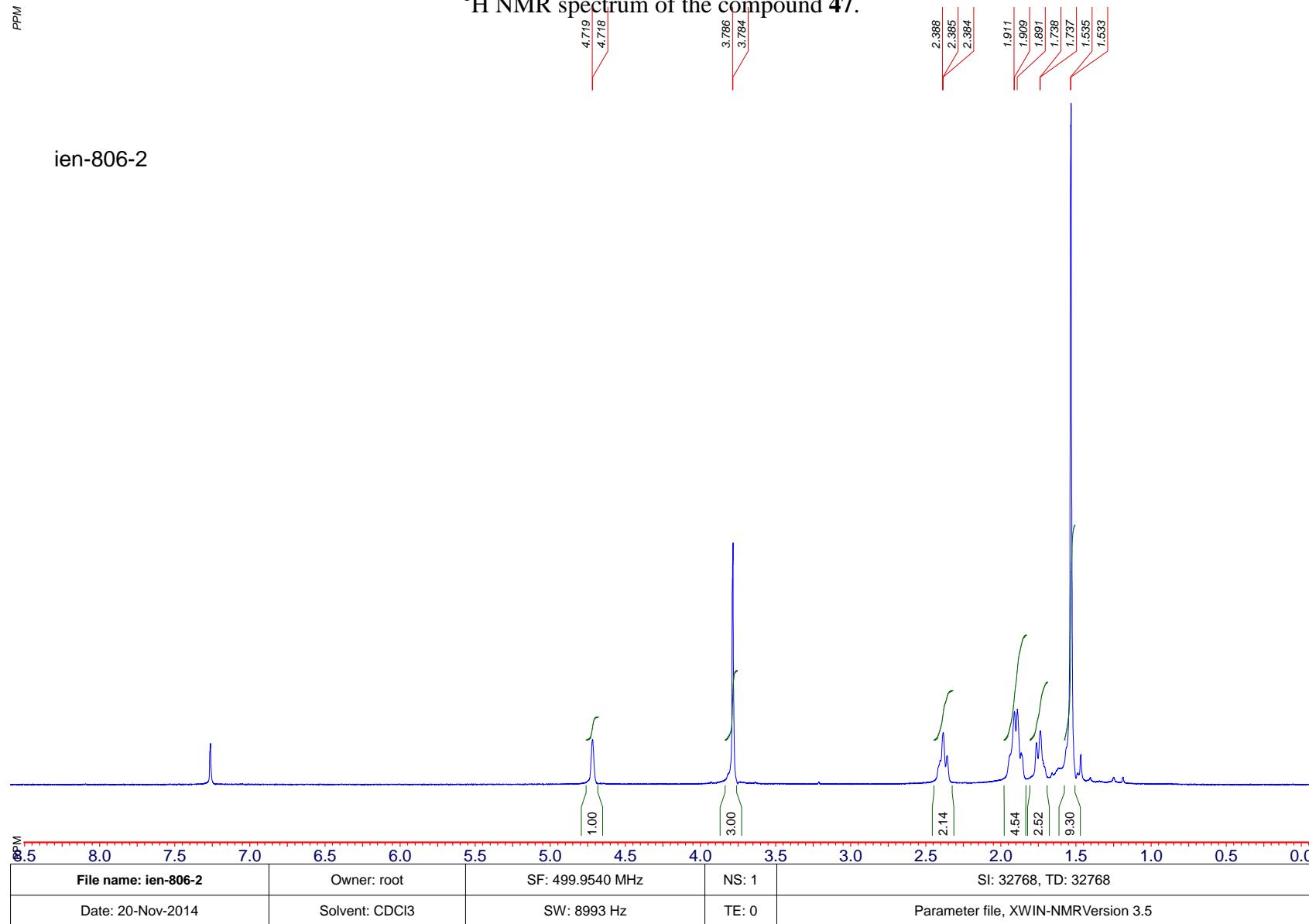
^{13}C NMR spectrum of the compound **42**.



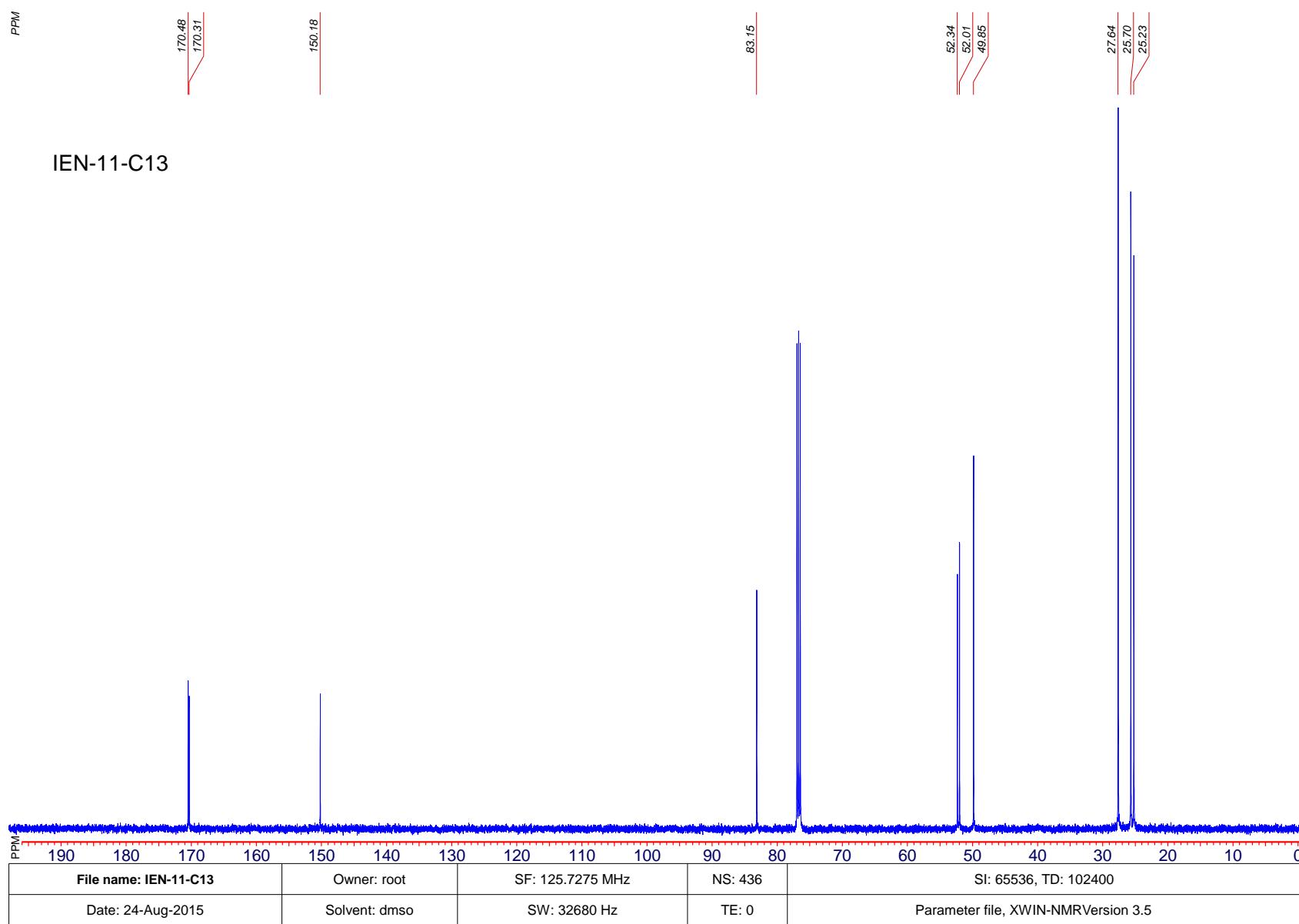
PPM

¹H NMR spectrum of the compound 47.

ien-806-2



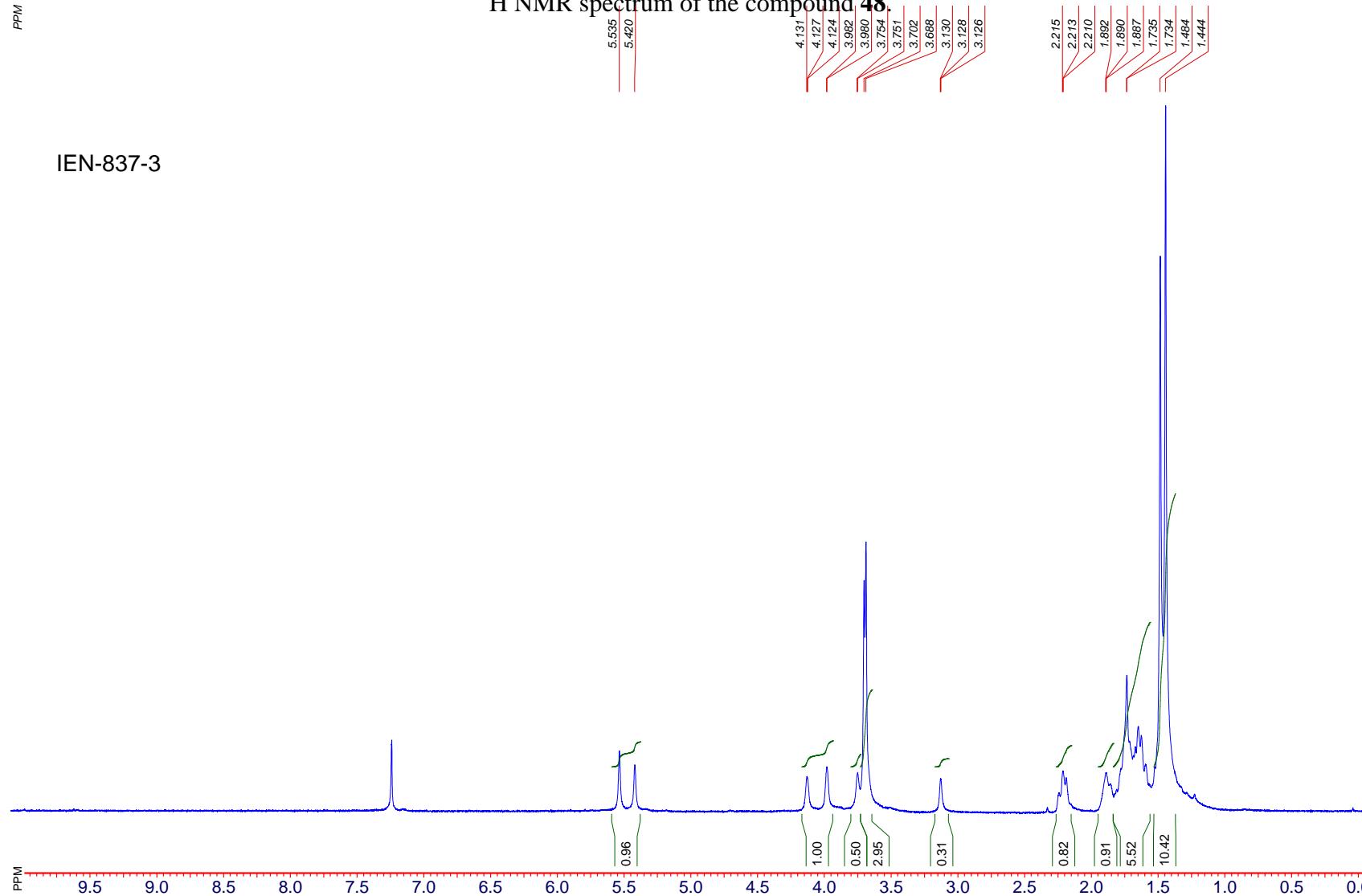
¹³C NMR spectrum of the compound **47**.



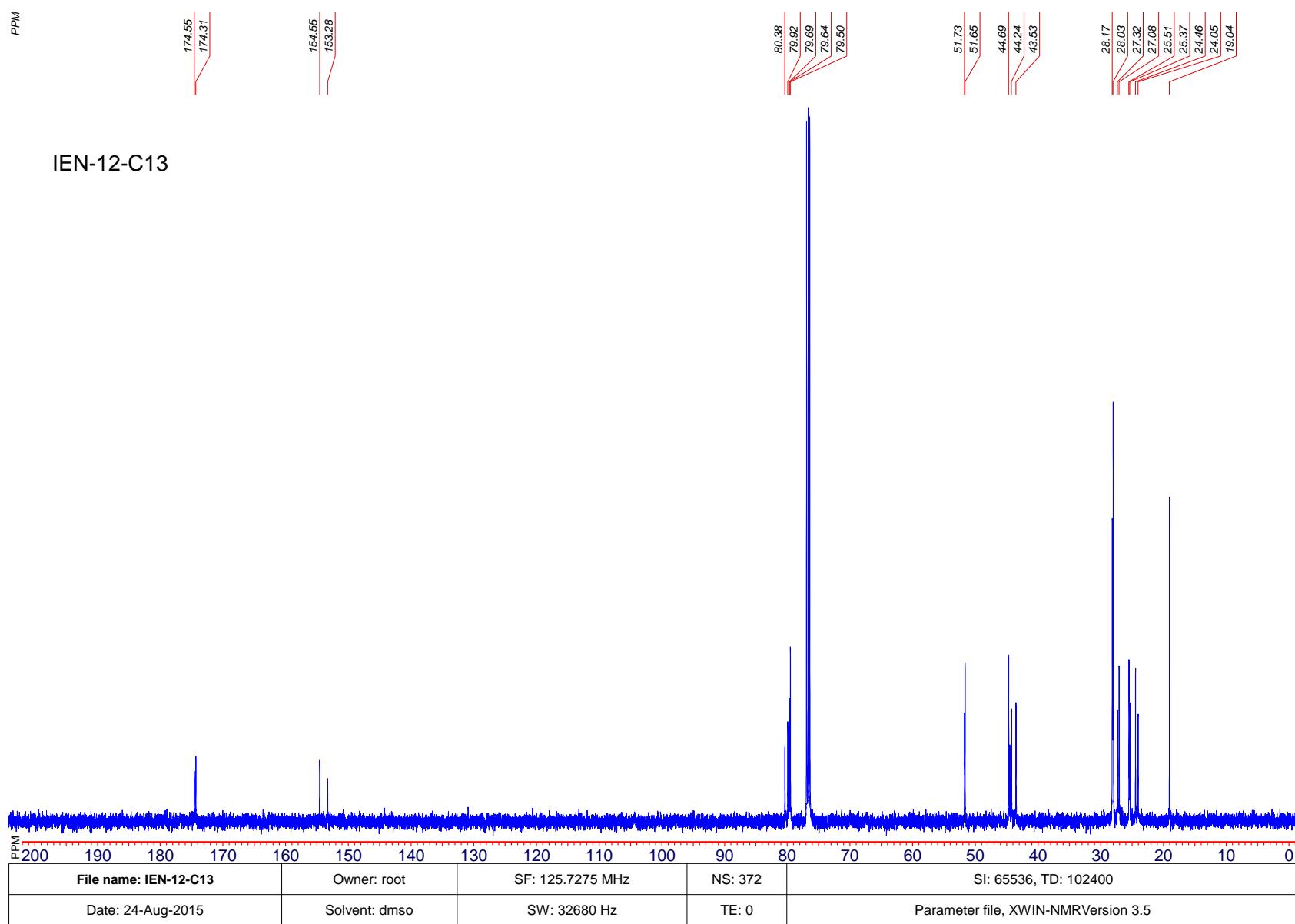
PPM

¹H NMR spectrum of the compound **48**.

IEN-837-3



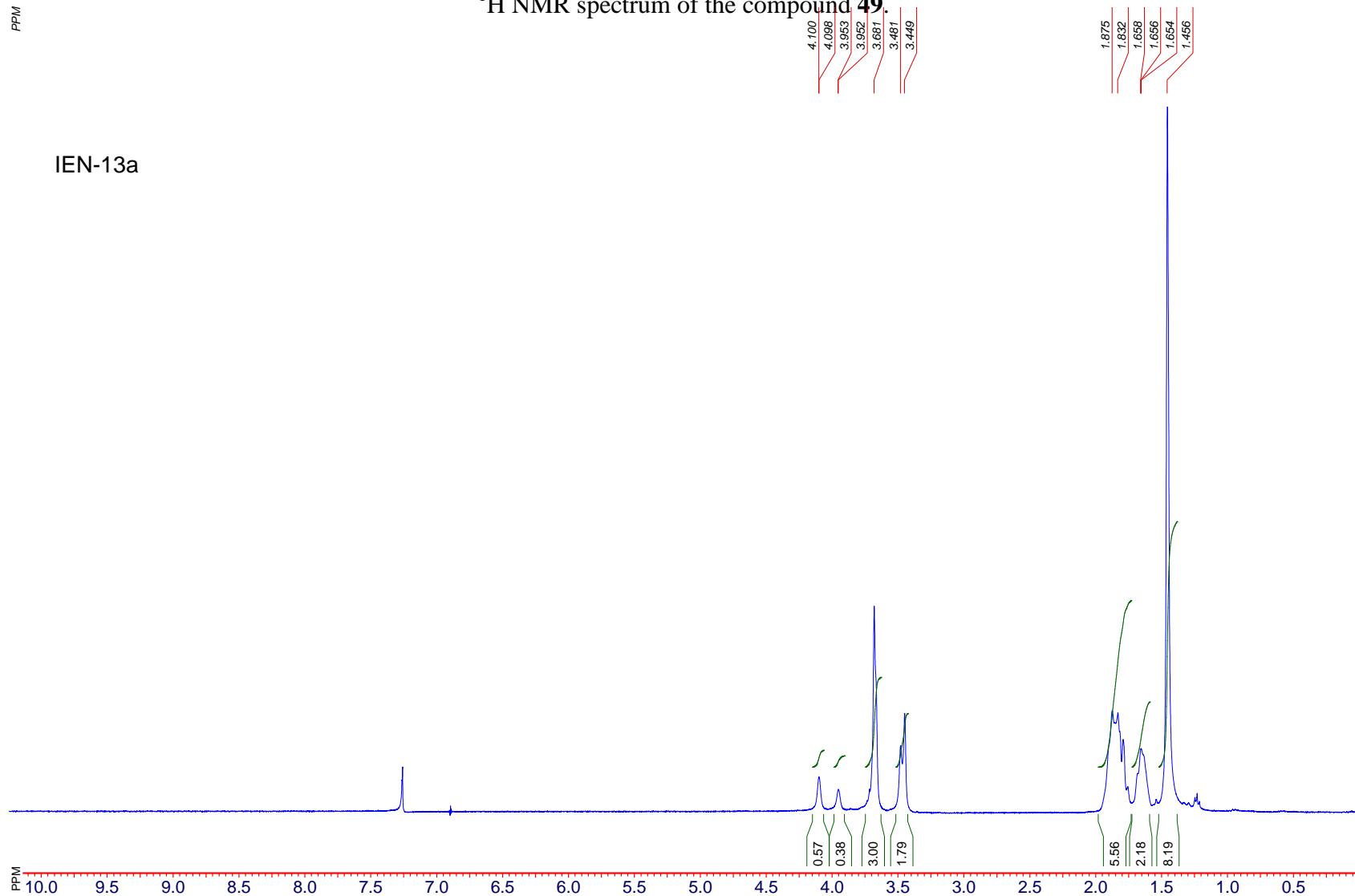
^{13}C NMR spectrum of the compound **48**.



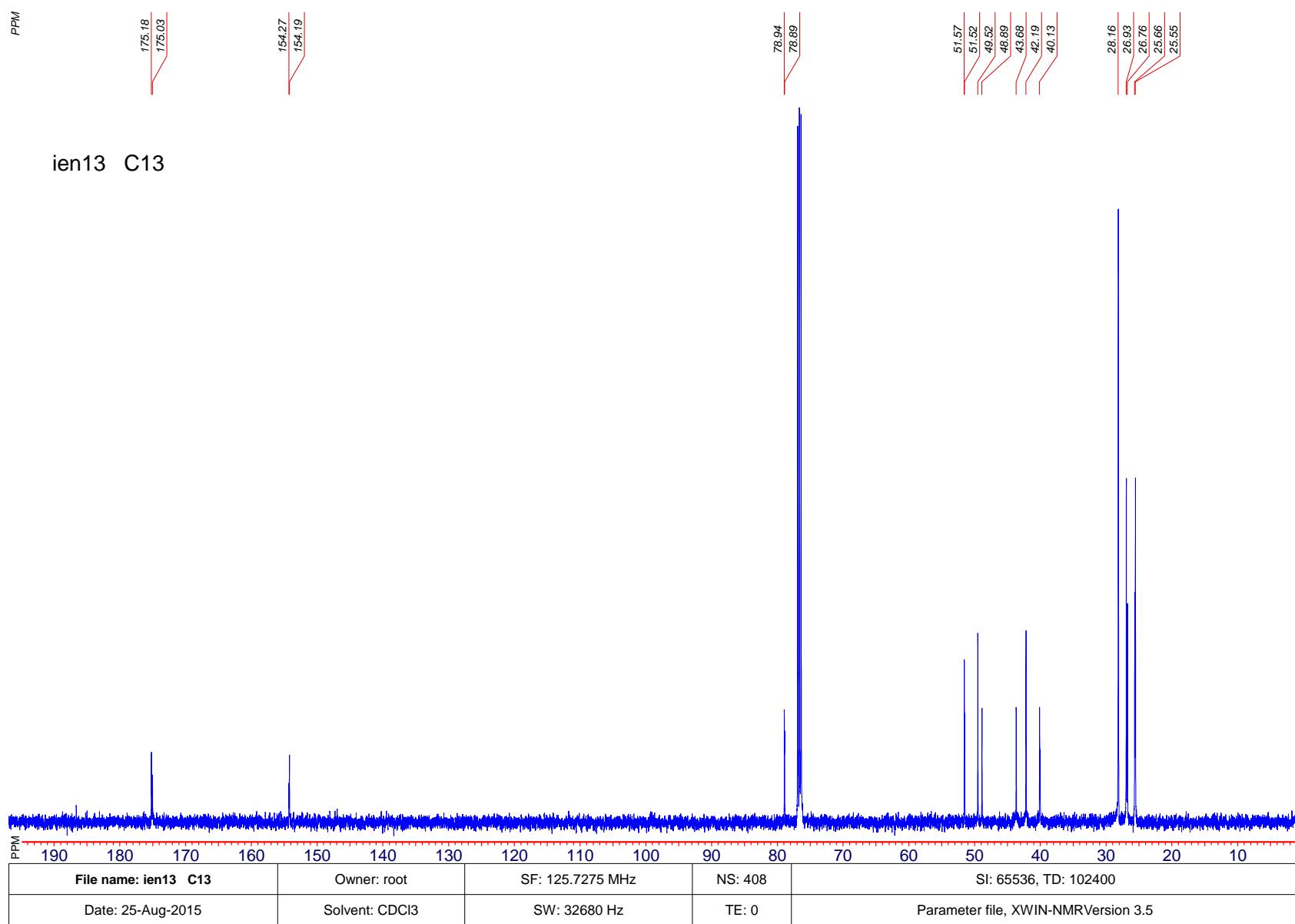
PPM

¹H NMR spectrum of the compound **49**.

IEN-13a



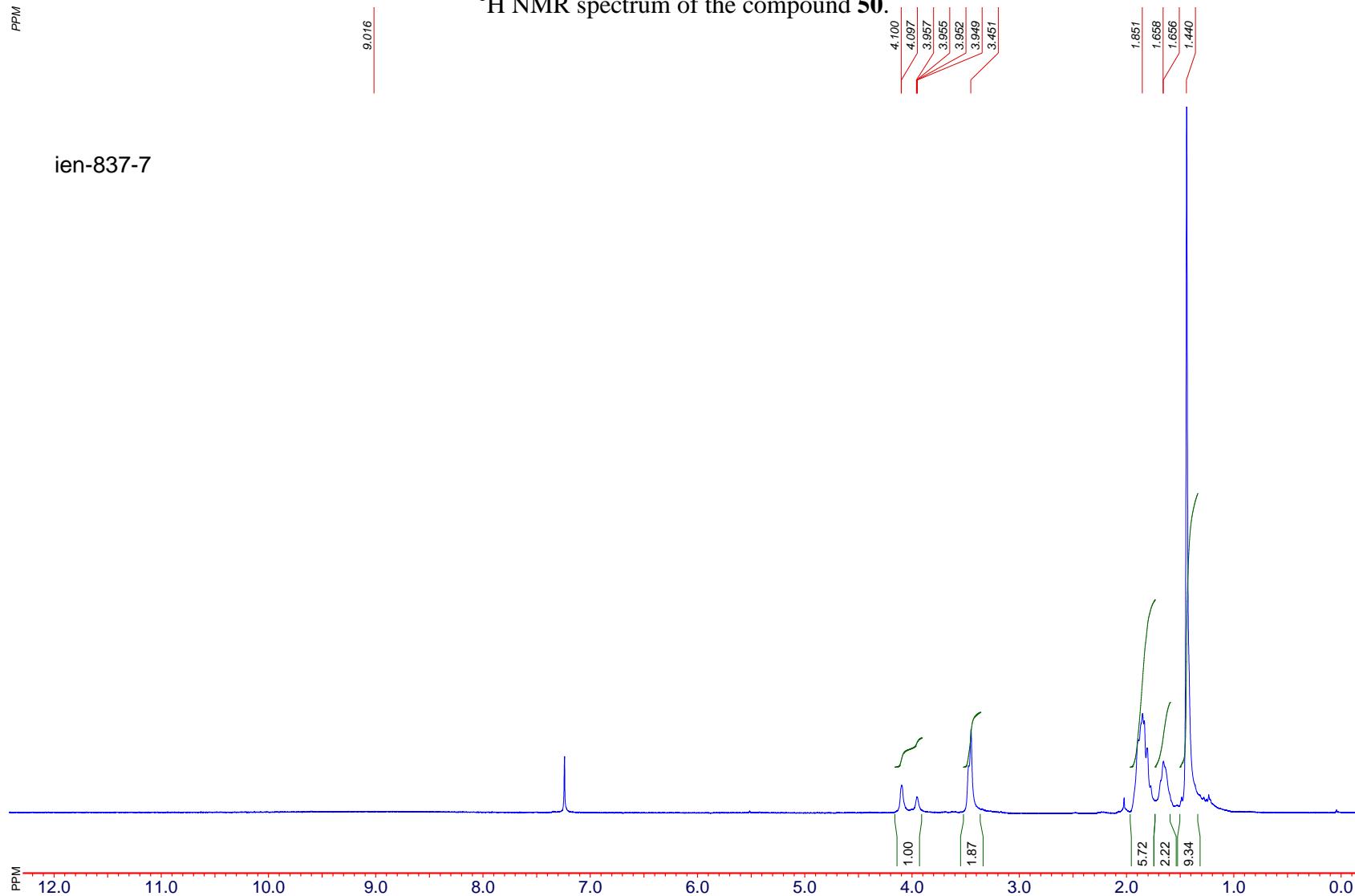
^{13}C NMR spectrum of the compound **49**.



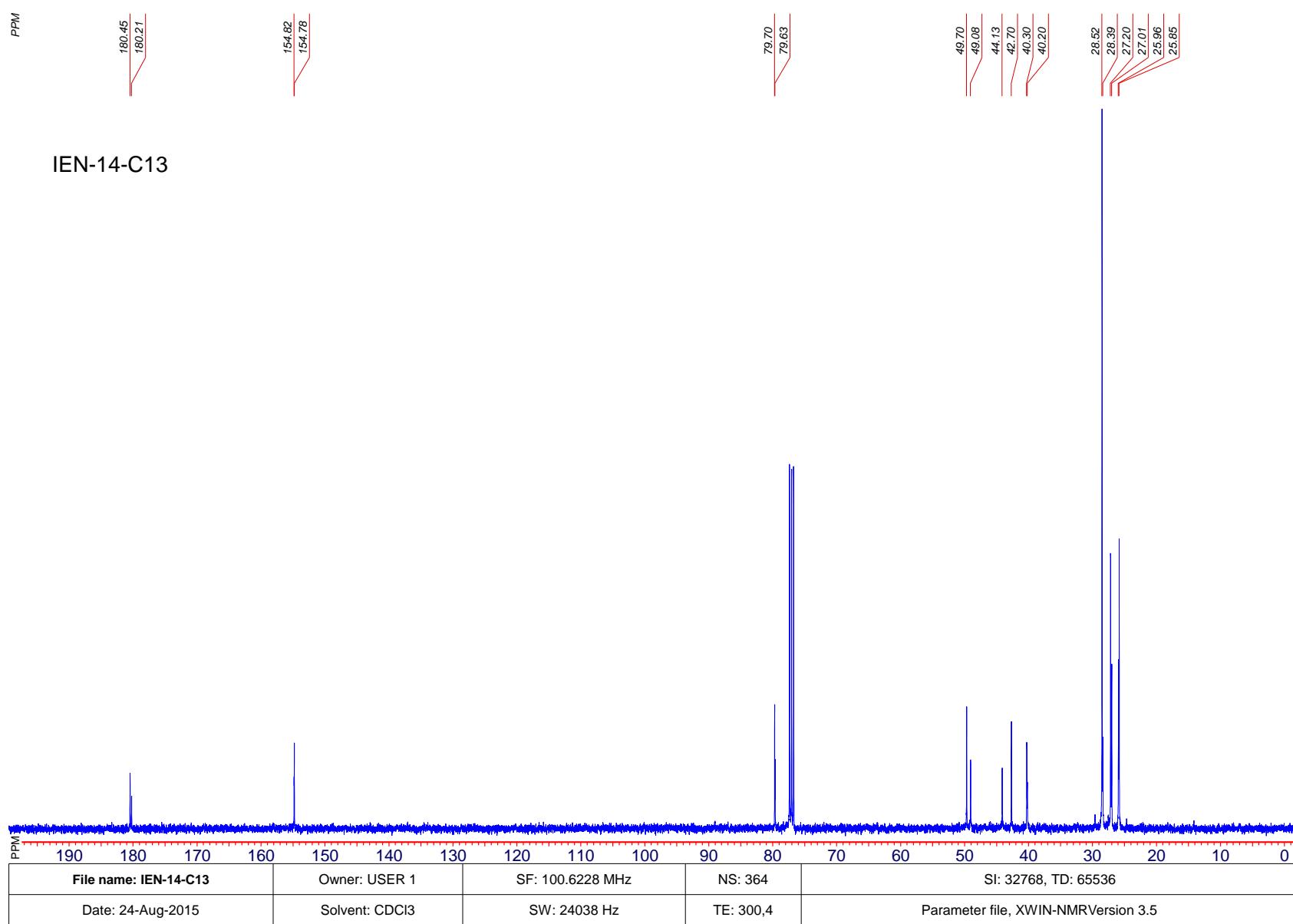
PPM

^1H NMR spectrum of the compound **50**.

ien-837-7



¹³C NMR spectrum of the compound **50**.



PPM

¹H NMR spectrum of the compound **11**·HCl.

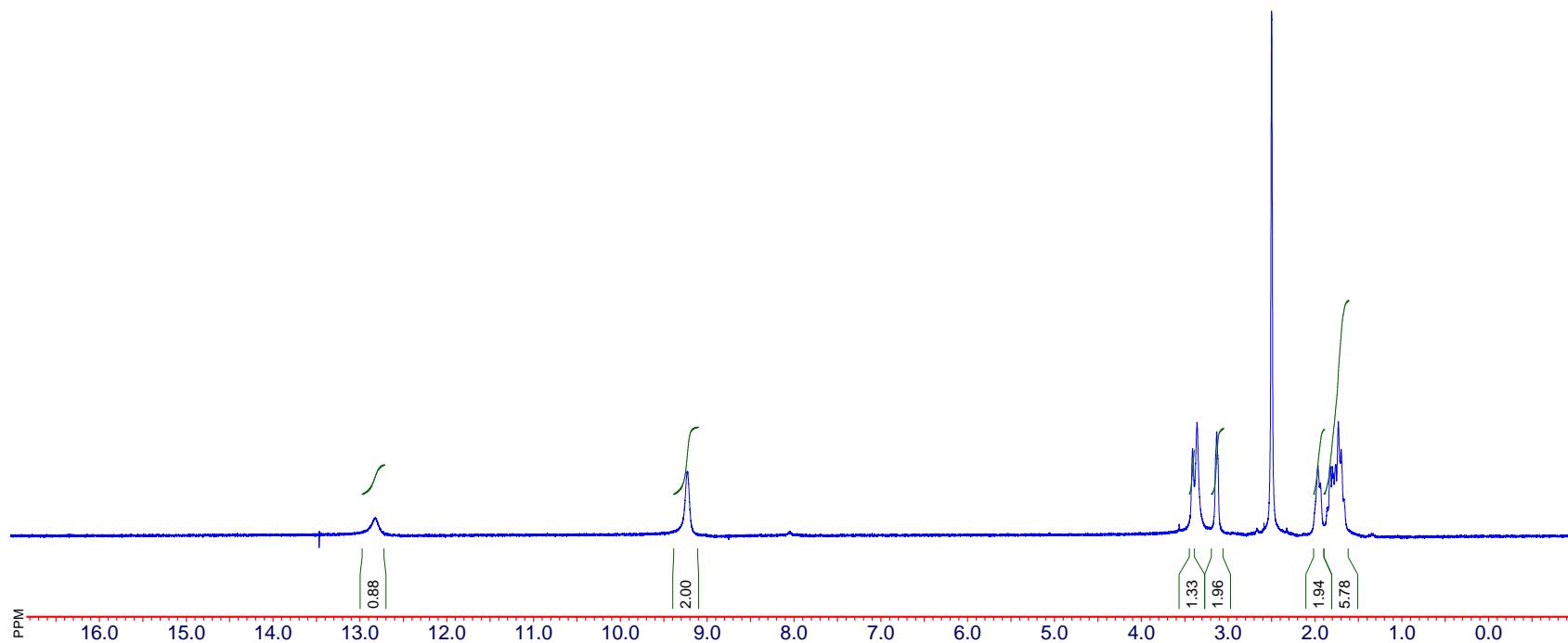
12.818

9.231
9.228
9.224

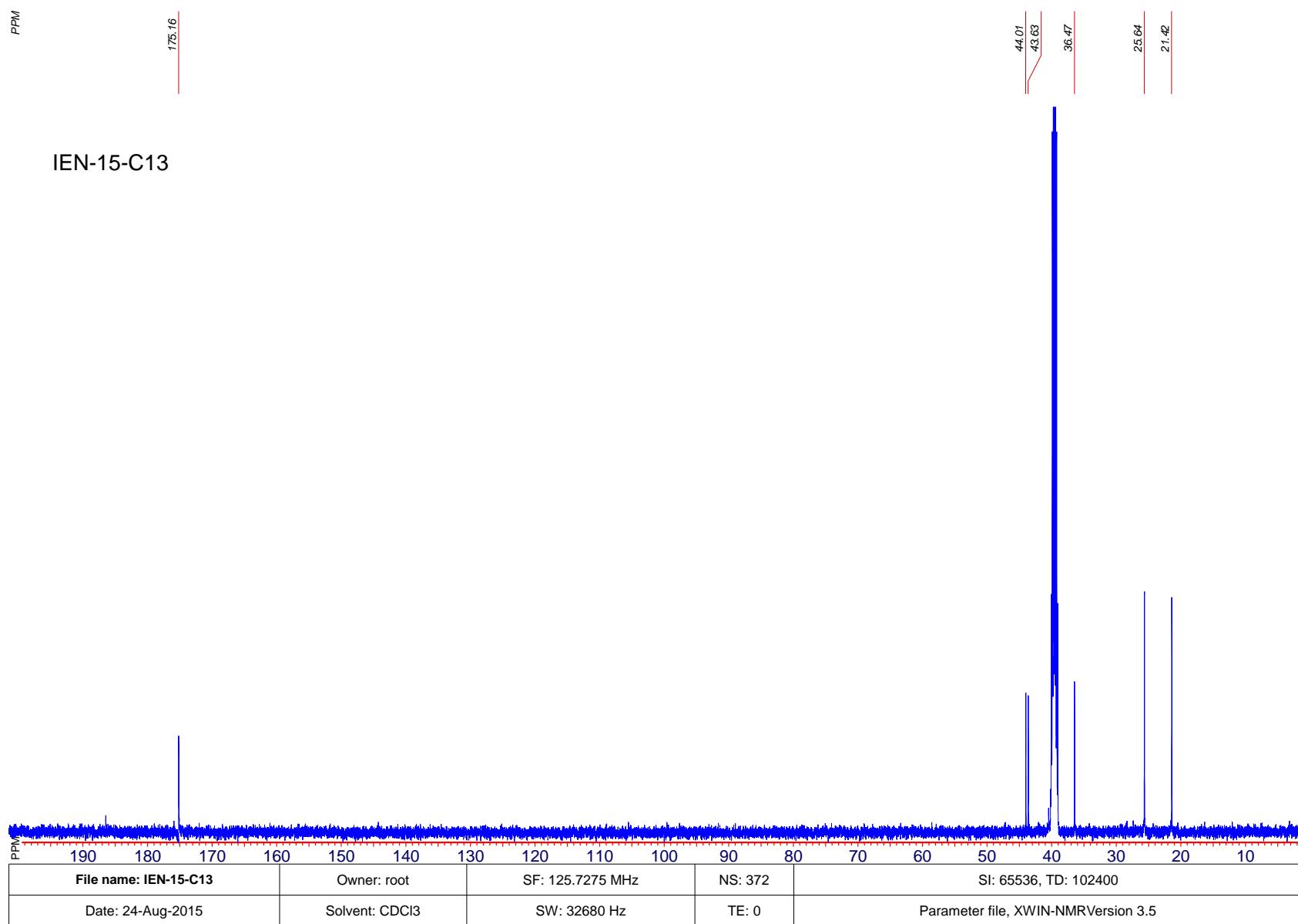
3.360
3.359
3.133
3.132
3.130

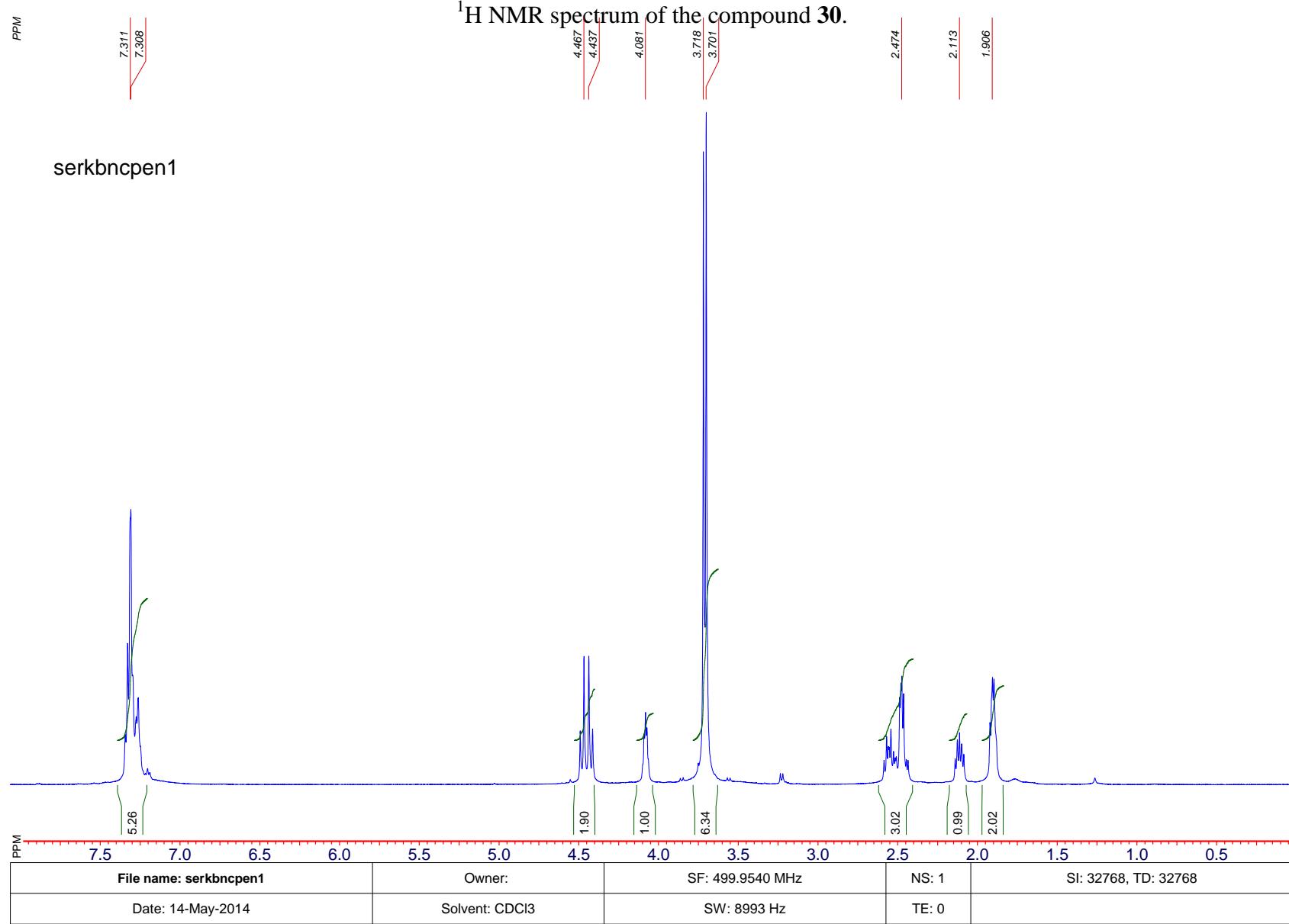
1.967
1.965
1.734
1.732

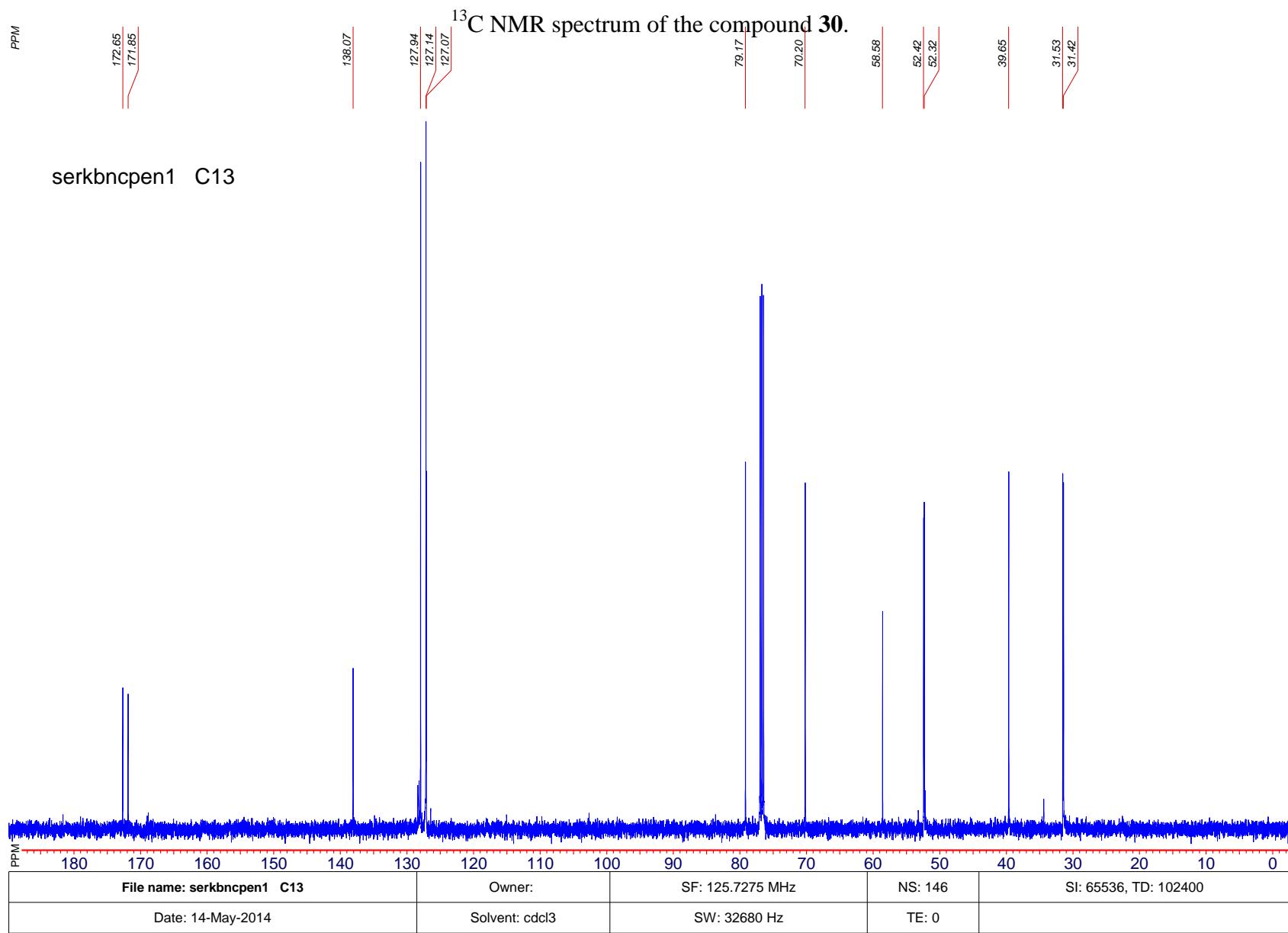
ien-844-1



¹³C NMR spectrum of the compound **11·HCl**.



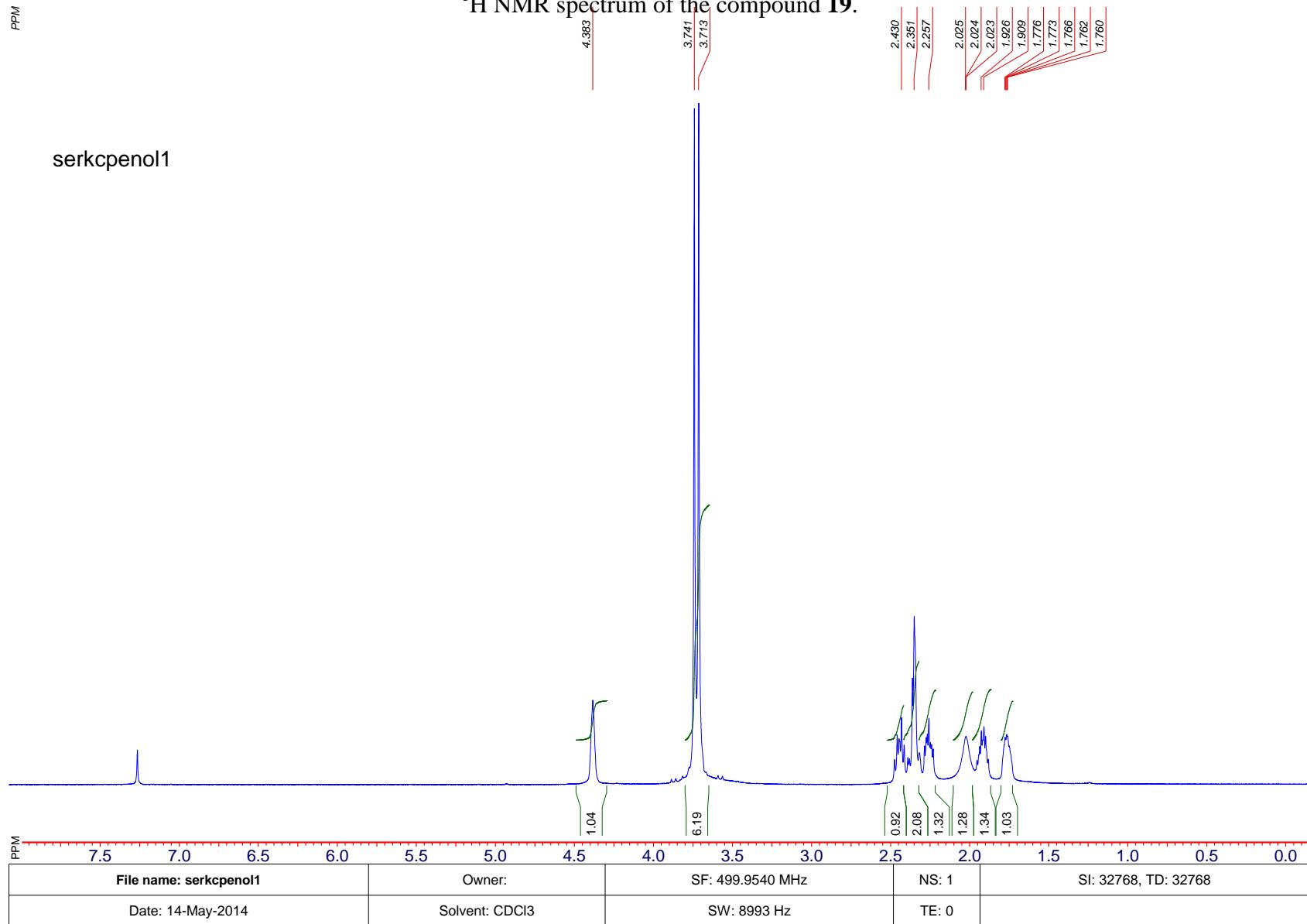




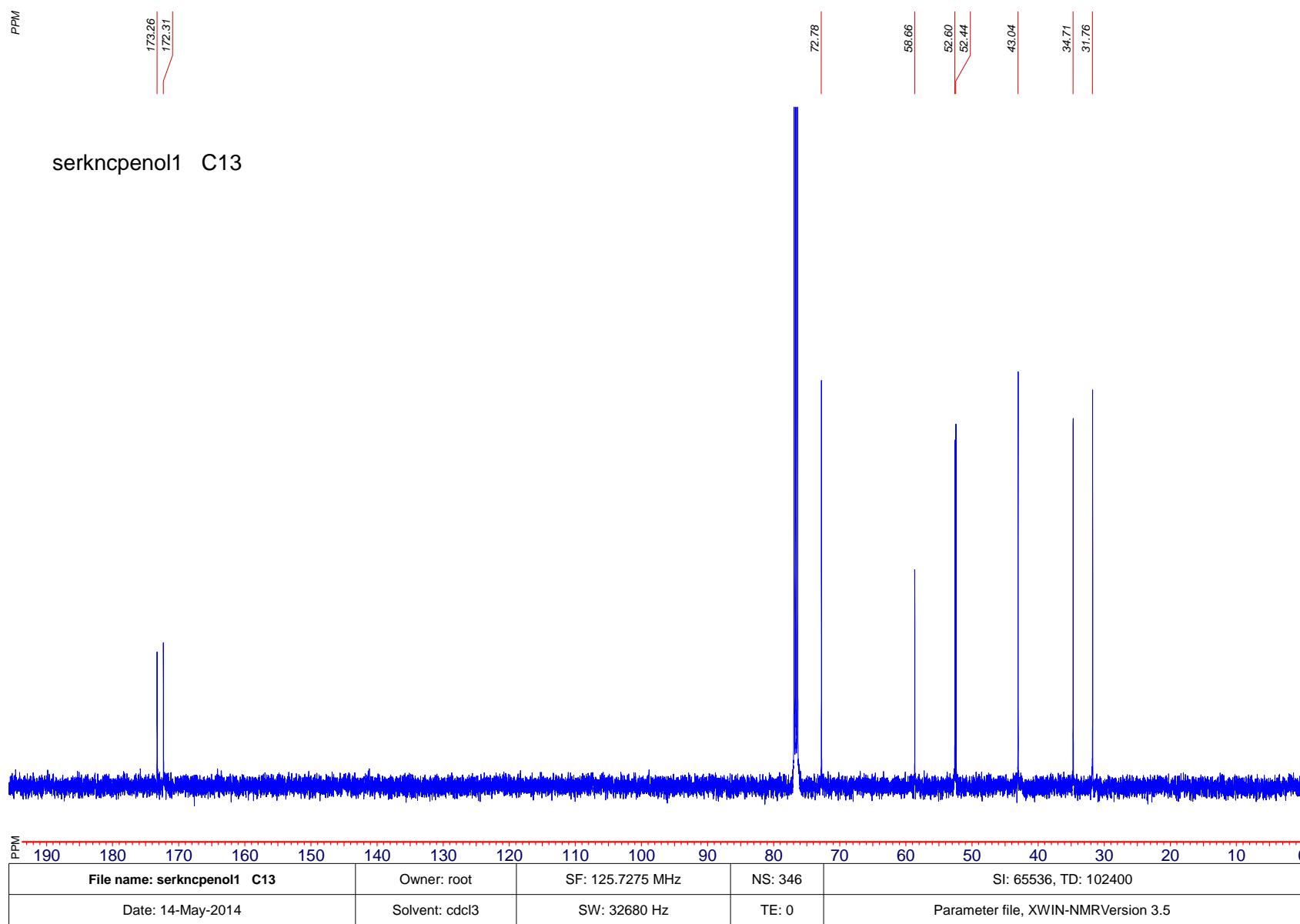
PPM

 ^1H NMR spectrum of the compound **19**.

serkcpenol1



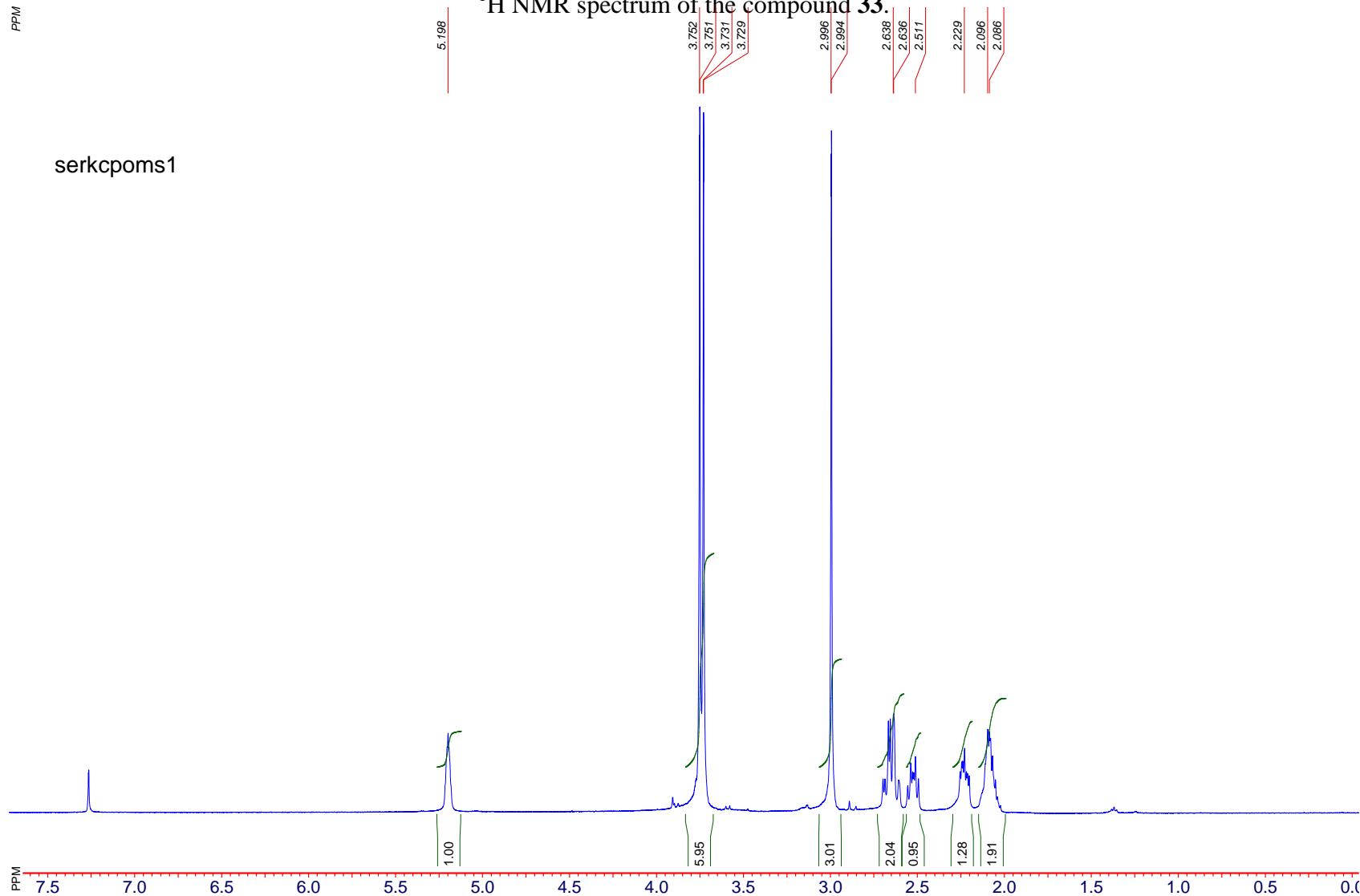
¹³C NMR spectrum of the compound **19**.



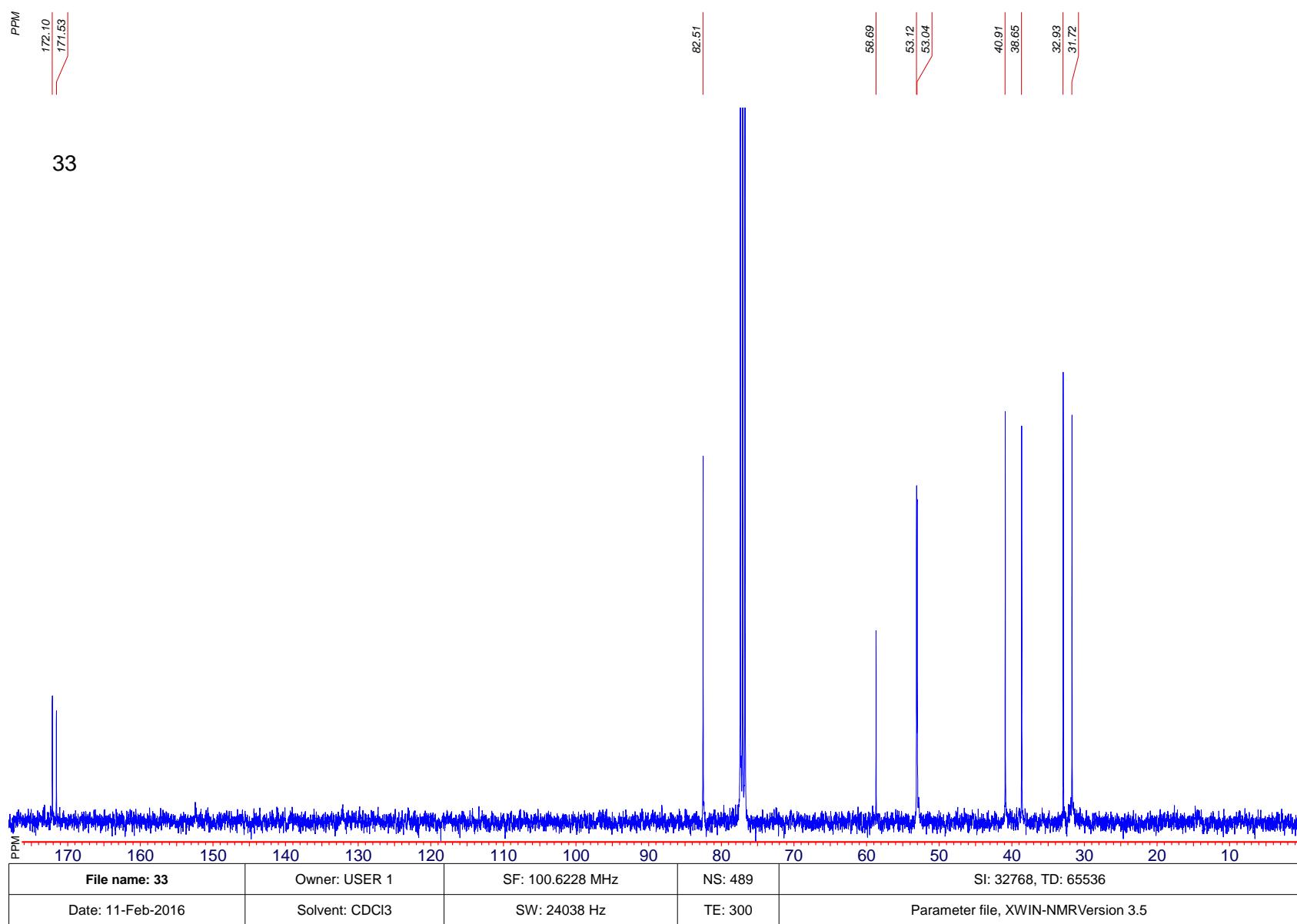
PPM

^1H NMR spectrum of the compound 33.

serkcpoms1



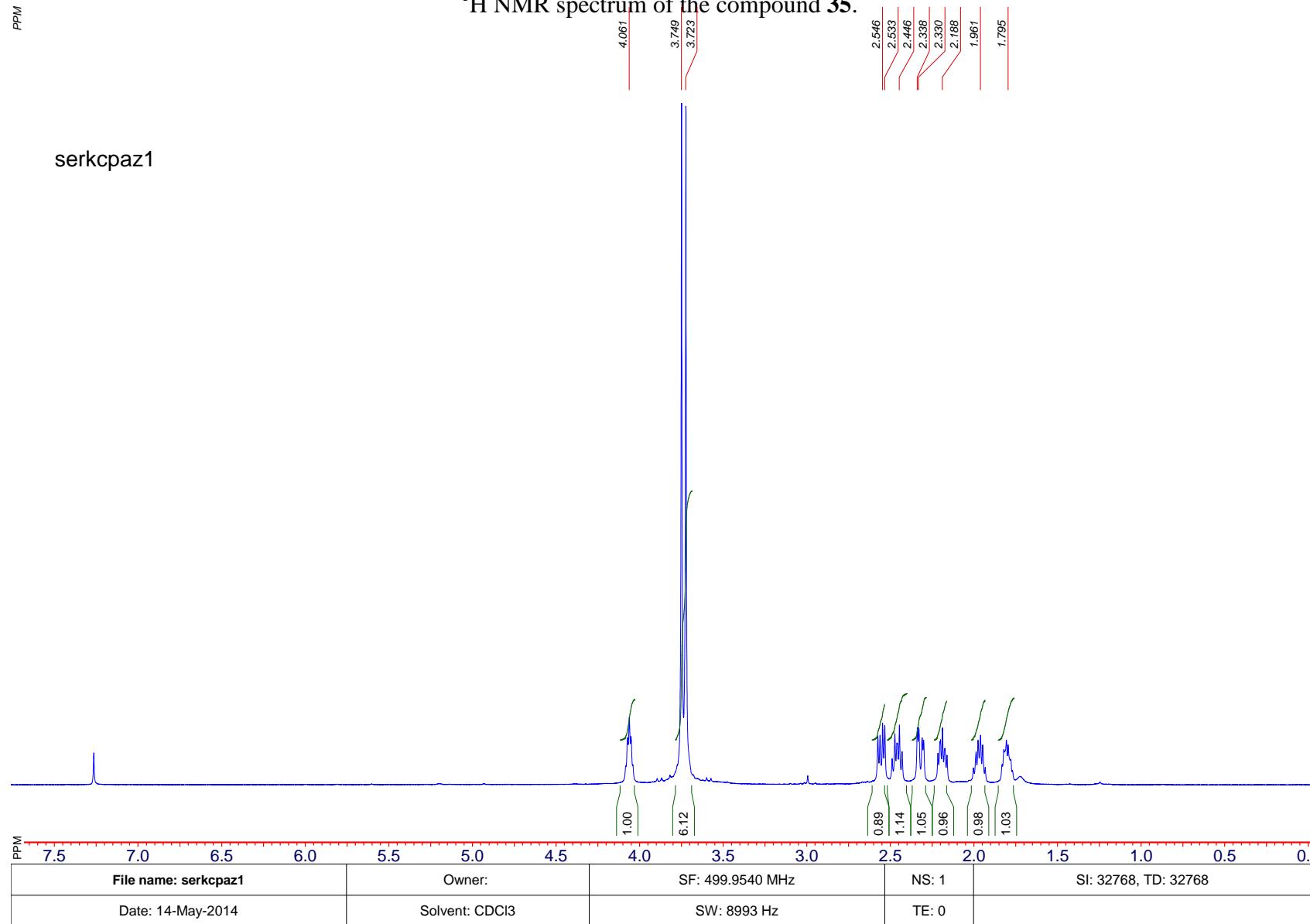
¹³C NMR spectrum of the compound 33.



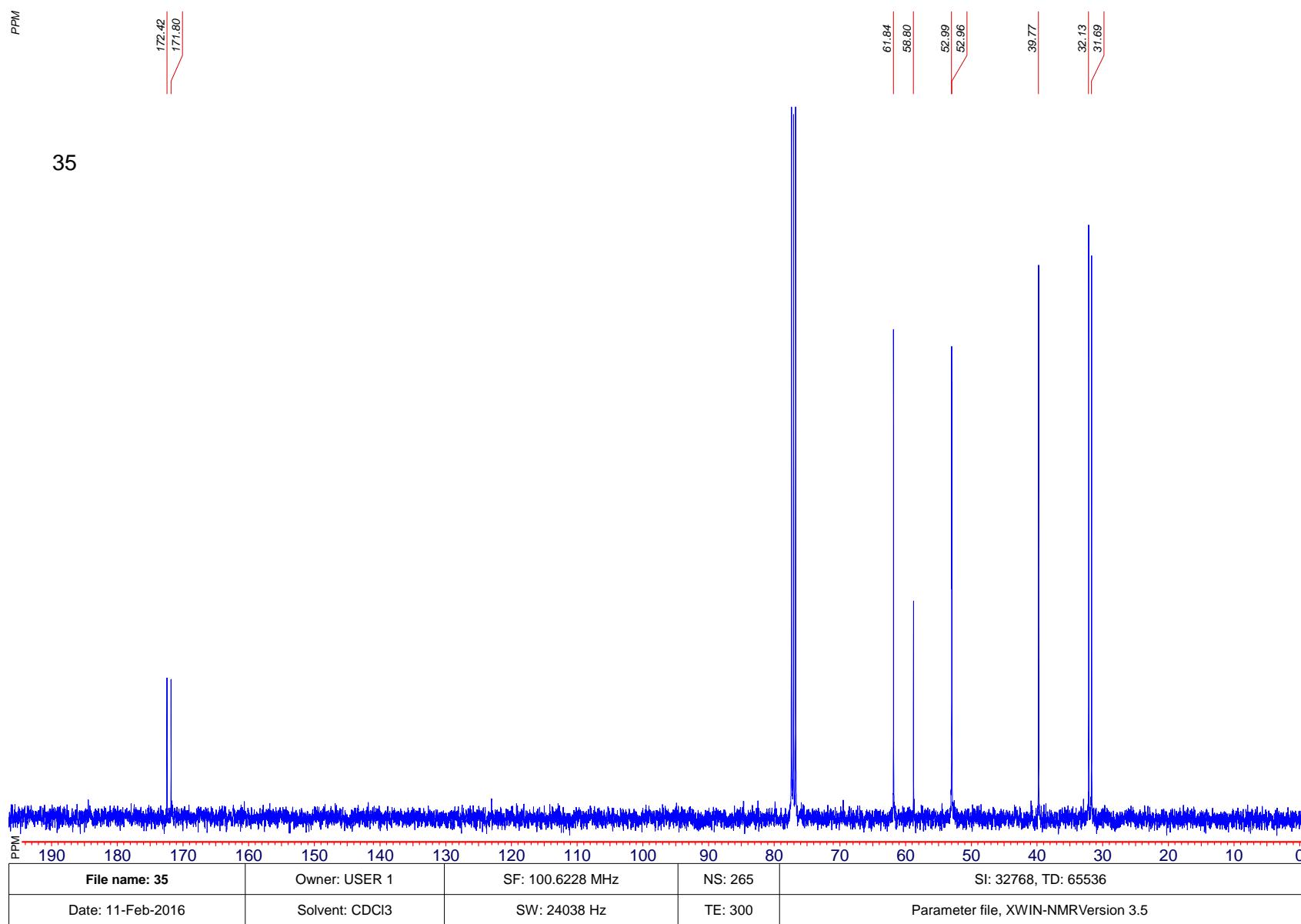
ppm

¹H NMR spectrum of the compound 35.

serkcpaz1



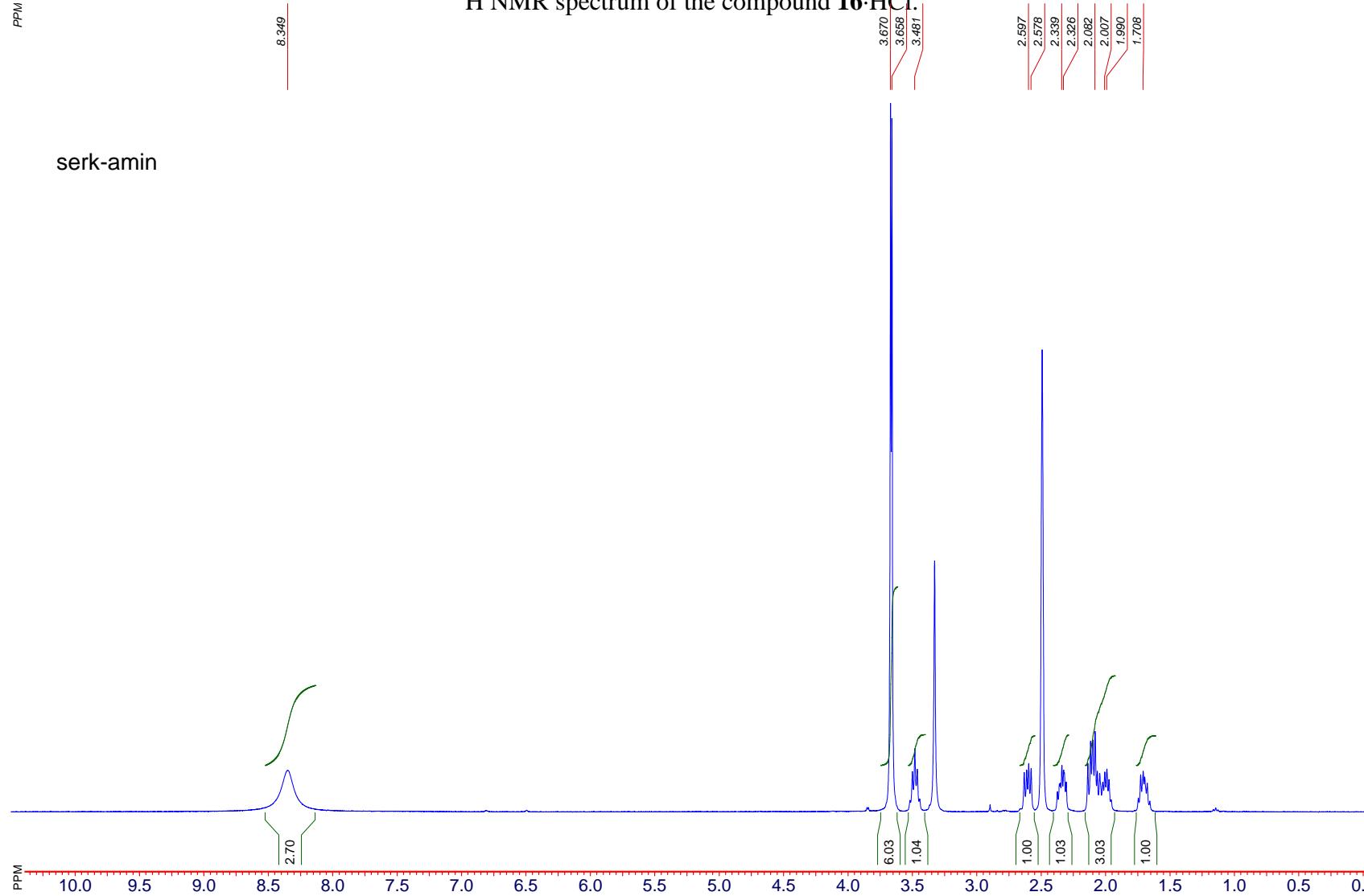
¹³C NMR spectrum of the compound **35**.



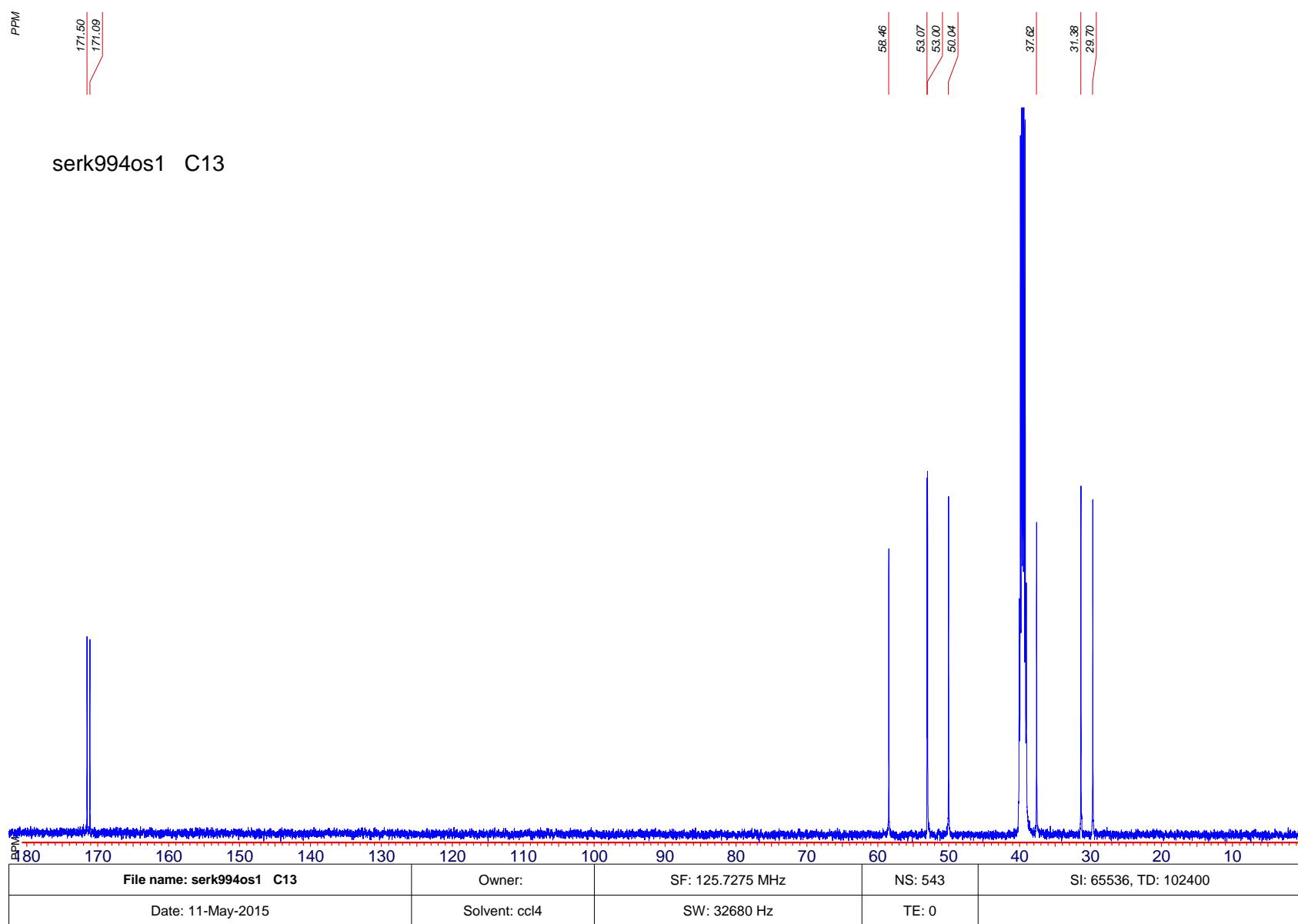
PPM

^1H NMR spectrum of the compound **16·HCl**

serk-amin



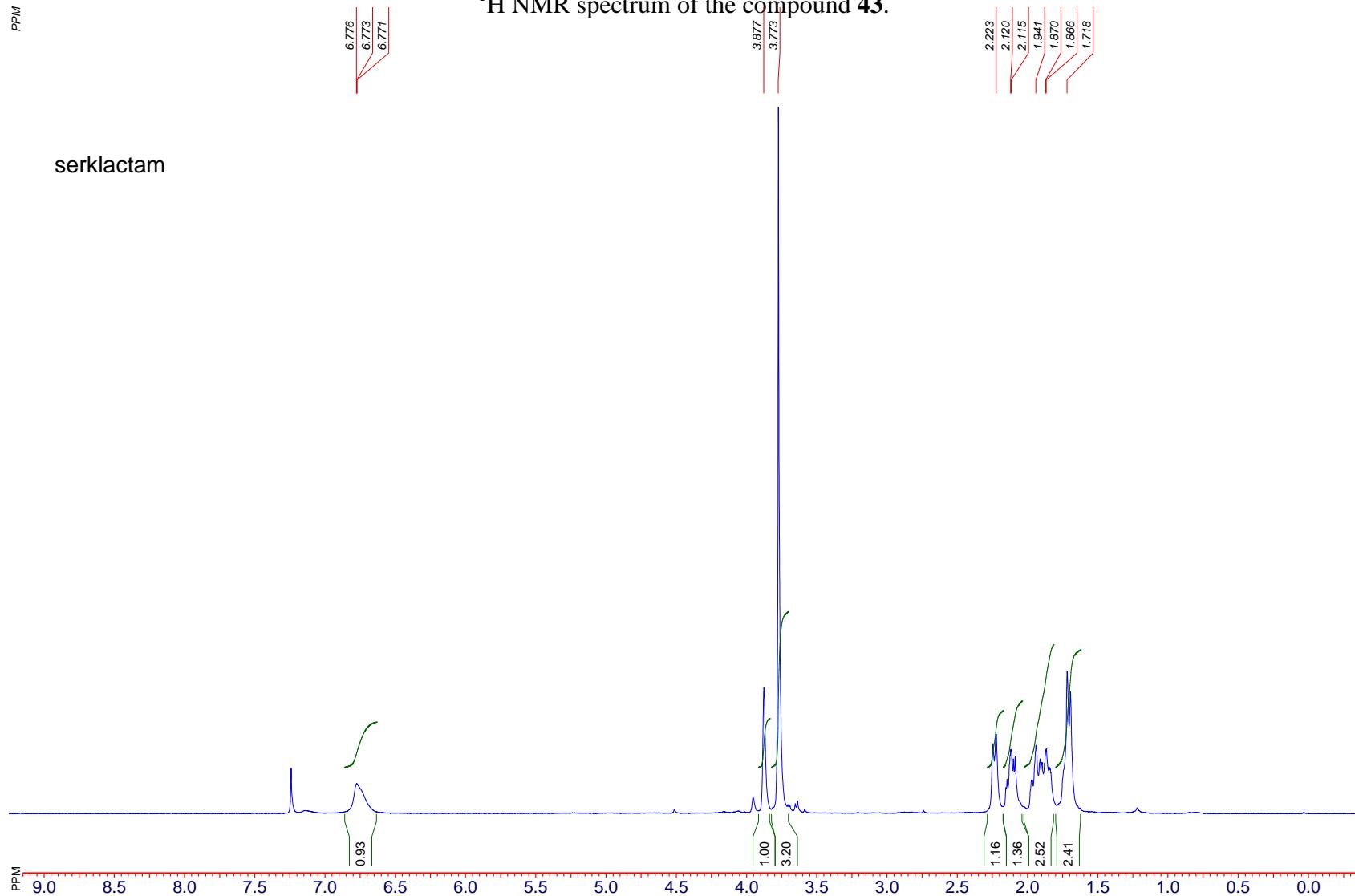
¹³C NMR spectrum of the compound **16·HCl**.



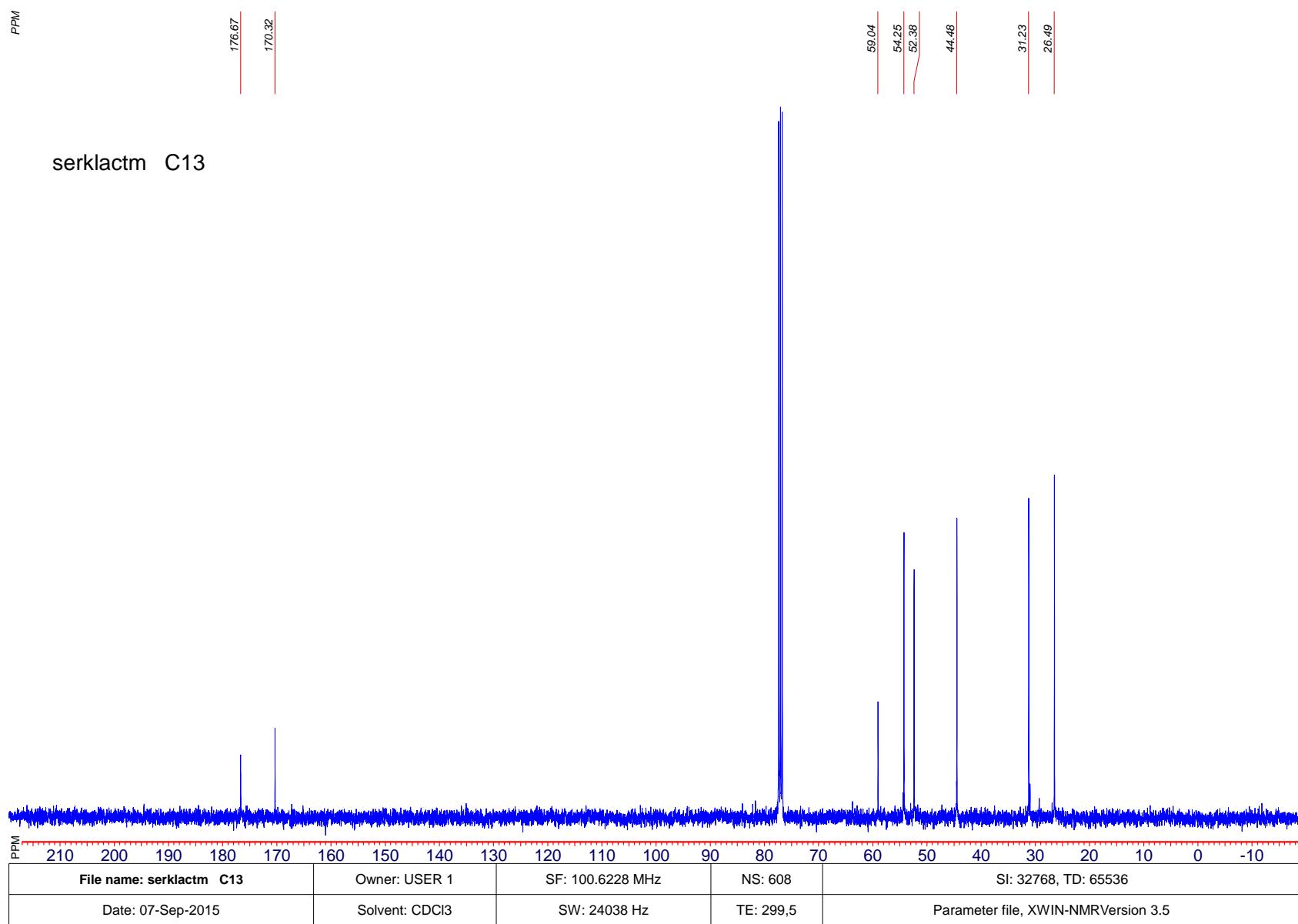
PPM

¹H NMR spectrum of the compound 43.

serklactam



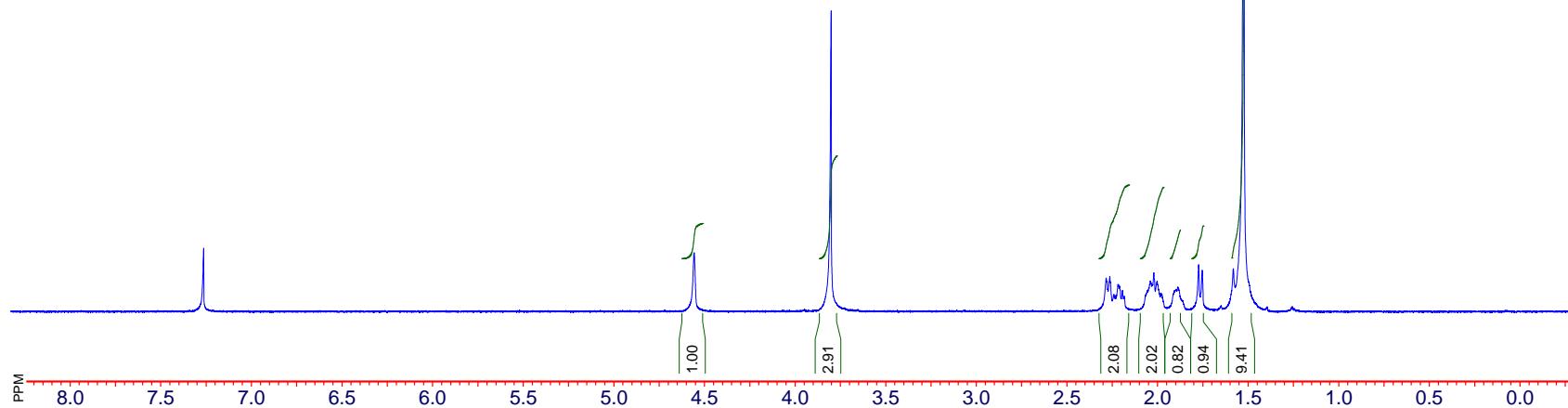
¹³C NMR spectrum of the compound **43**.



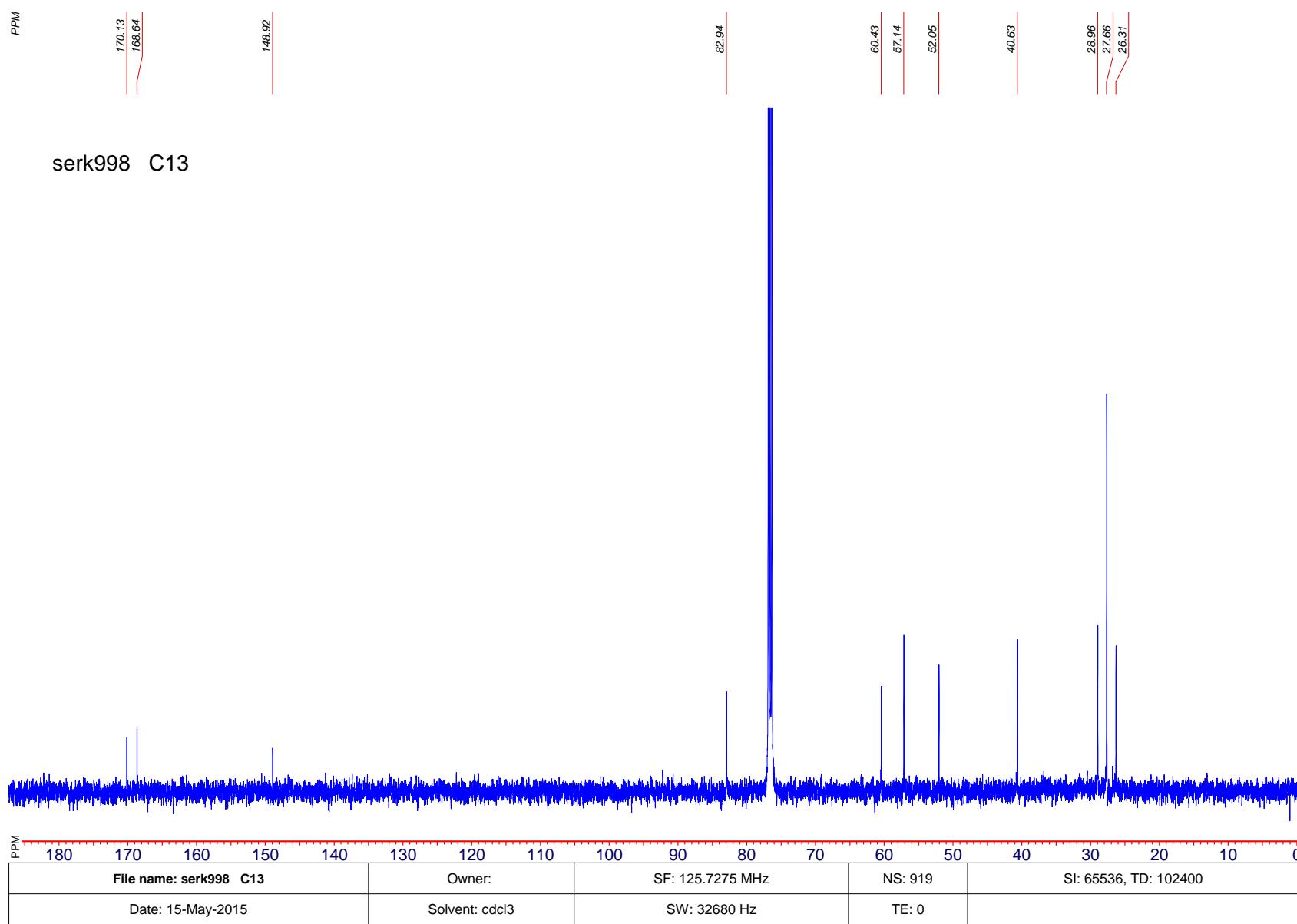
PPM

^1H NMR spectrum of the compound **51**.

serk998



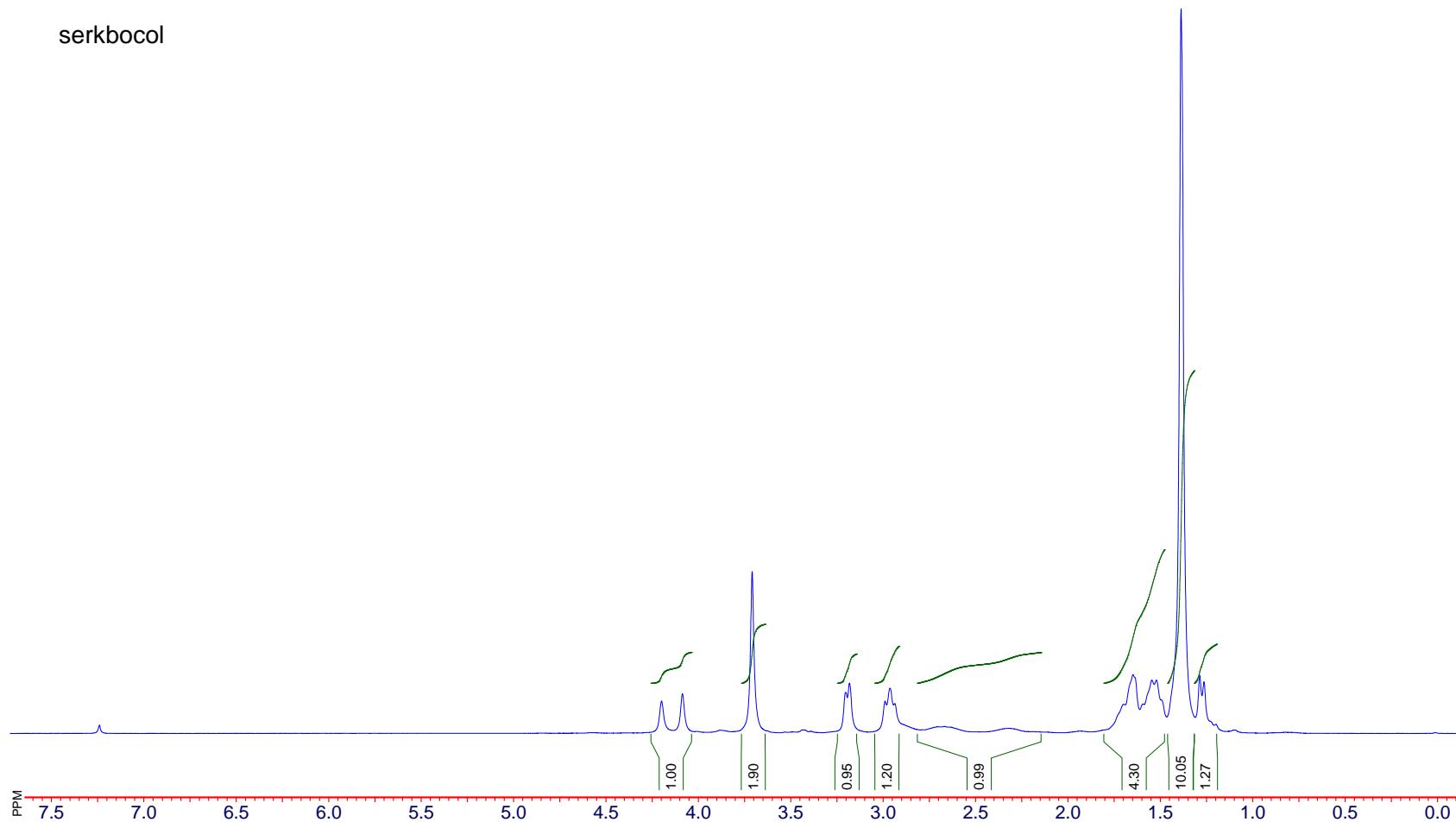
^{13}C NMR spectrum of the compound **51**.



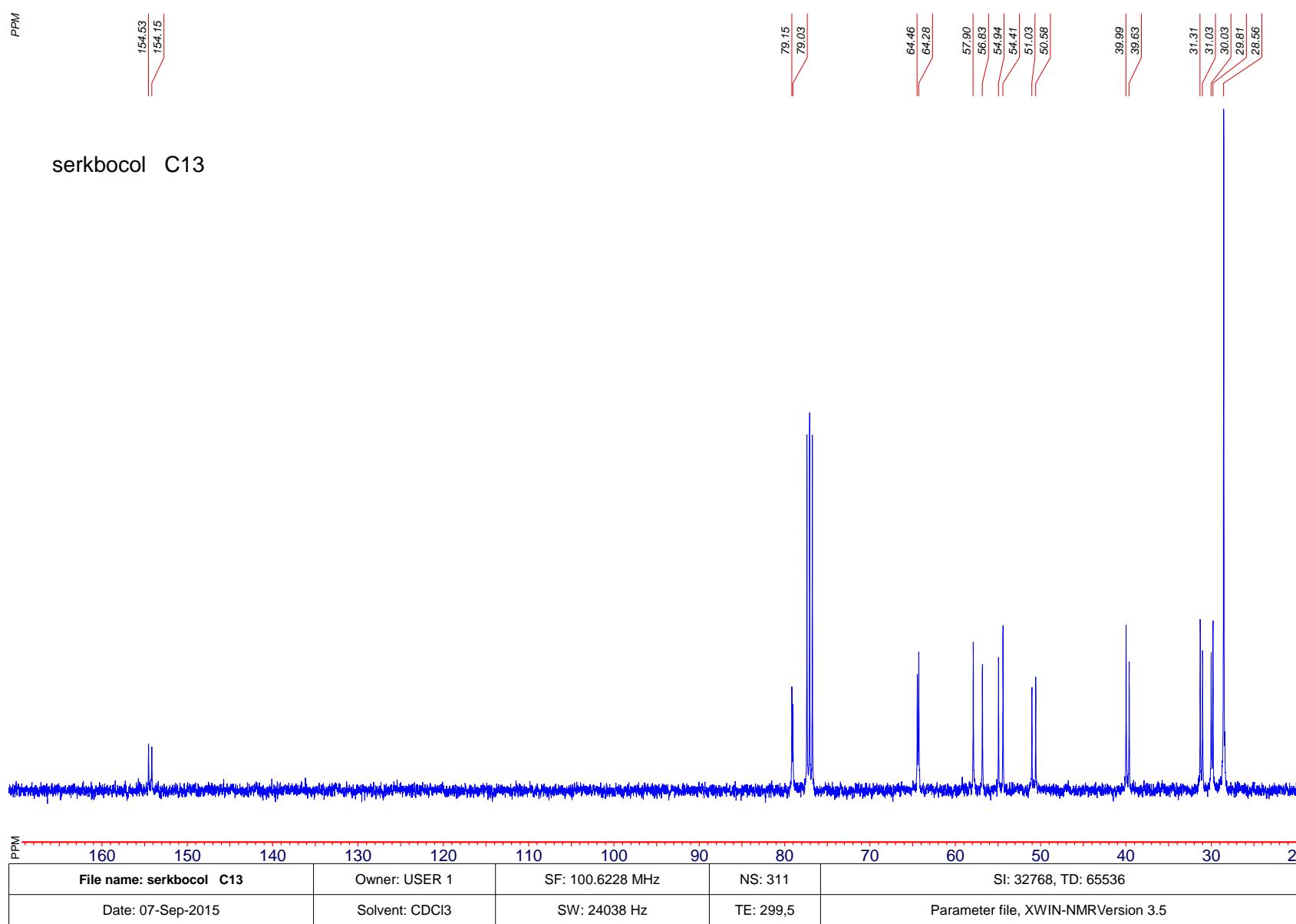
PPM

^1H NMR spectrum of the compound **52**.

serkbocol



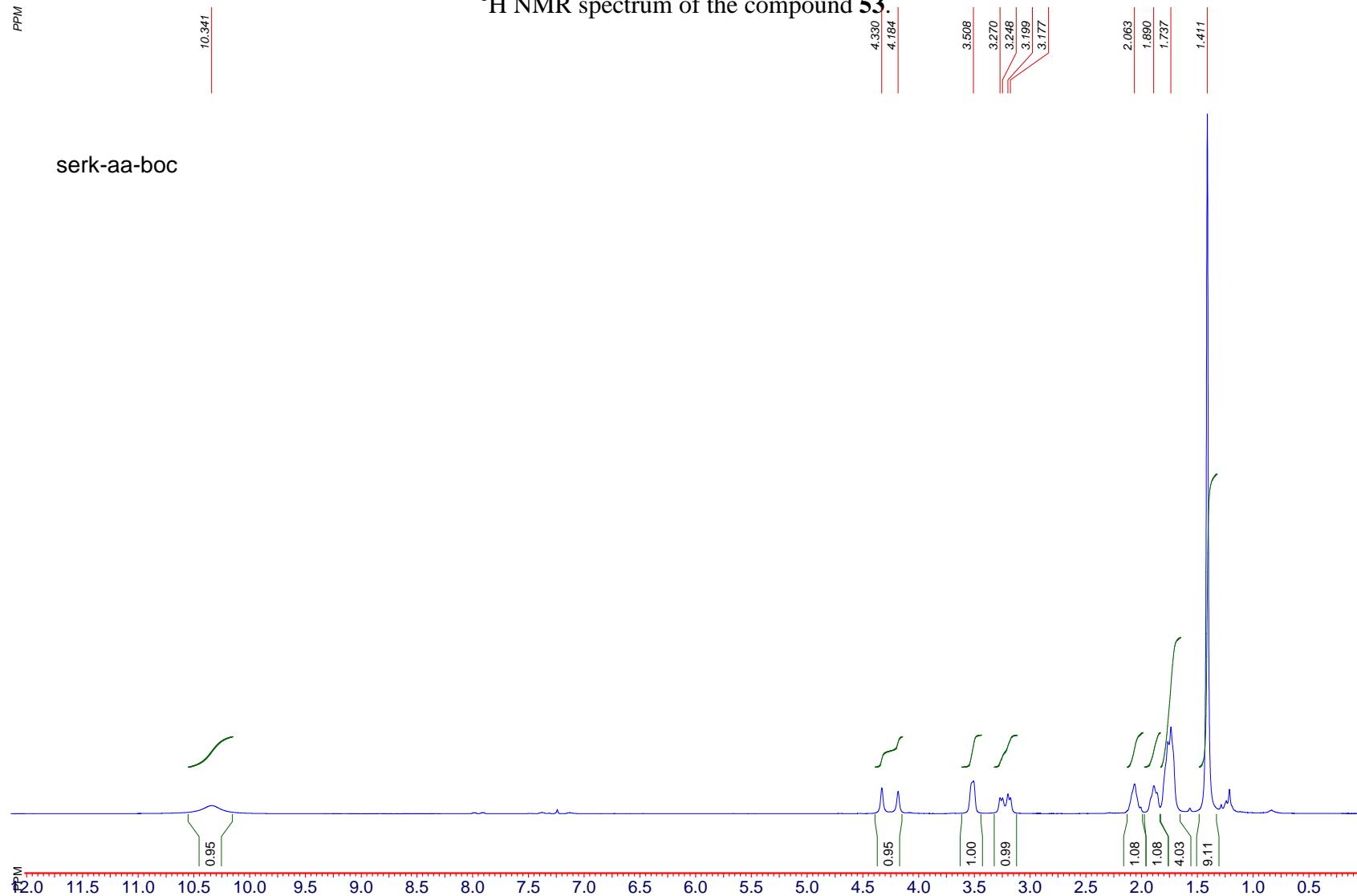
¹³C NMR spectrum of the compound **52**.



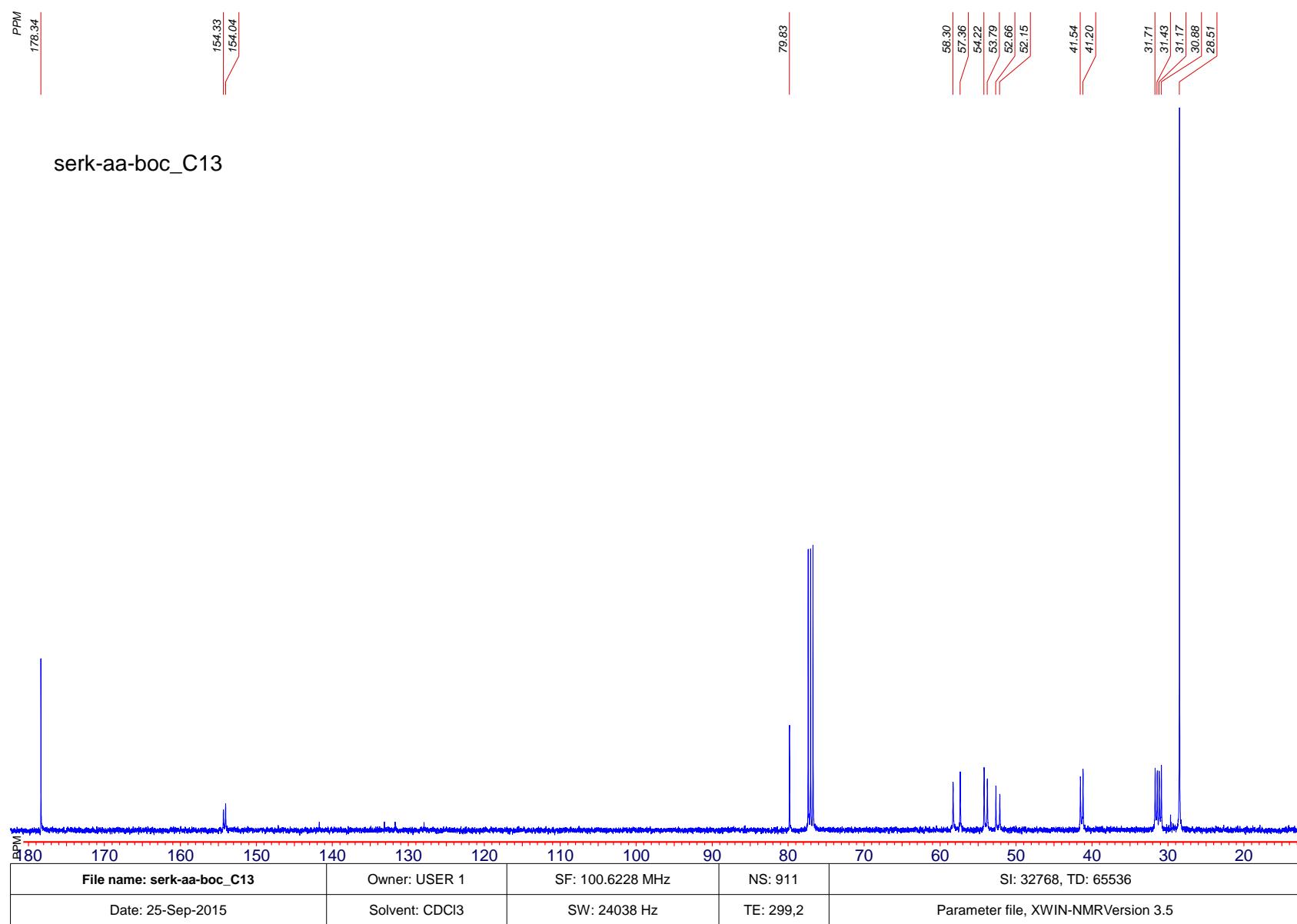
PPM

^1H NMR spectrum of the compound **53**.

serk-aa-boc



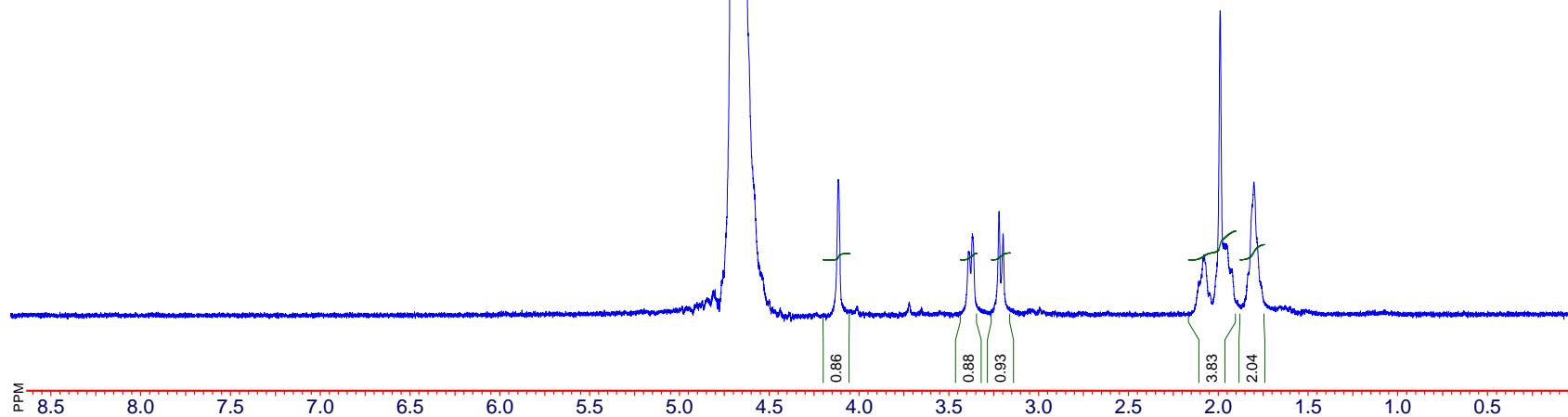
^{13}C NMR spectrum of the compound **53**.



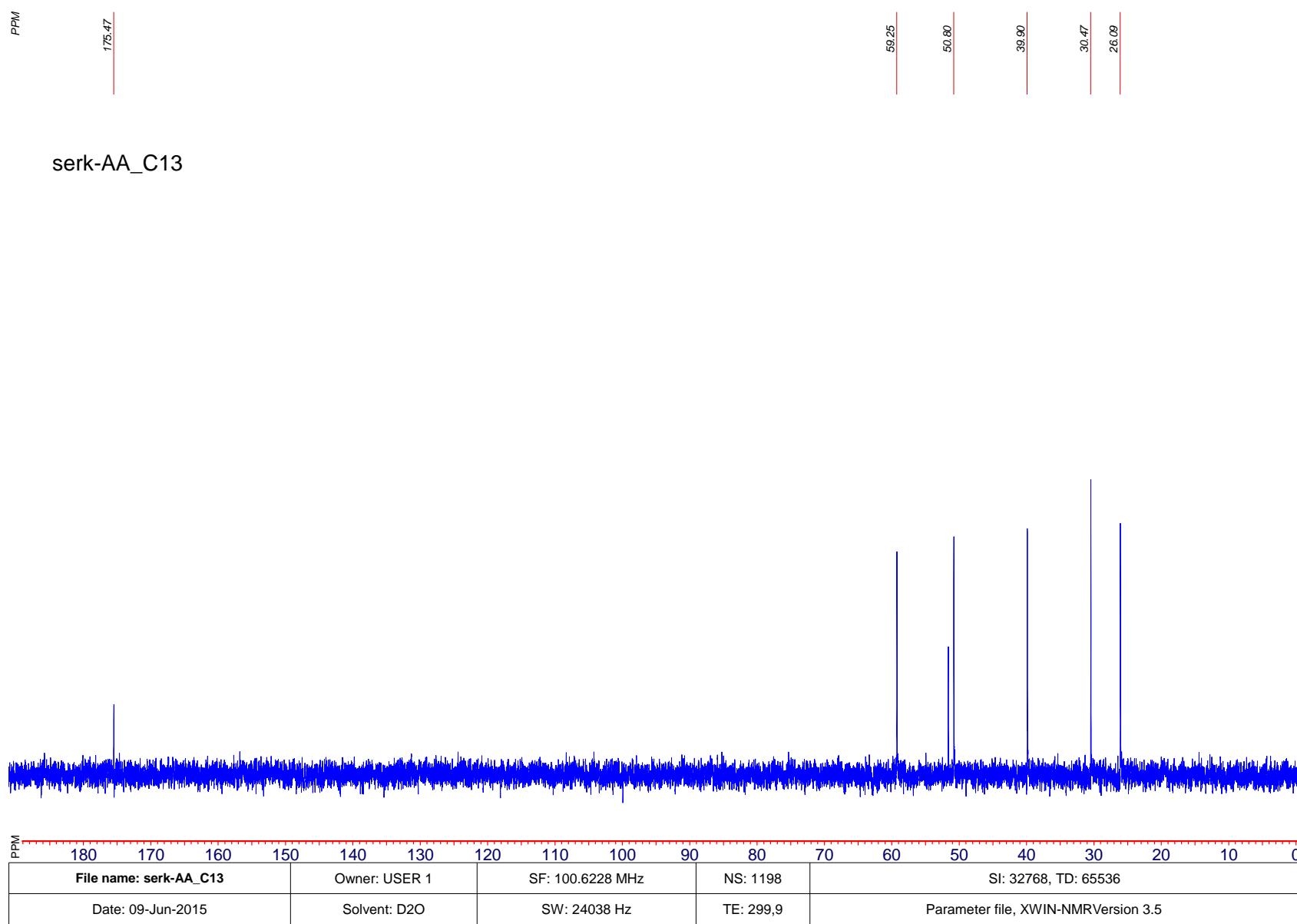
PPM

^1H NMR spectrum of the compound **12·HCl**.

serkaa1



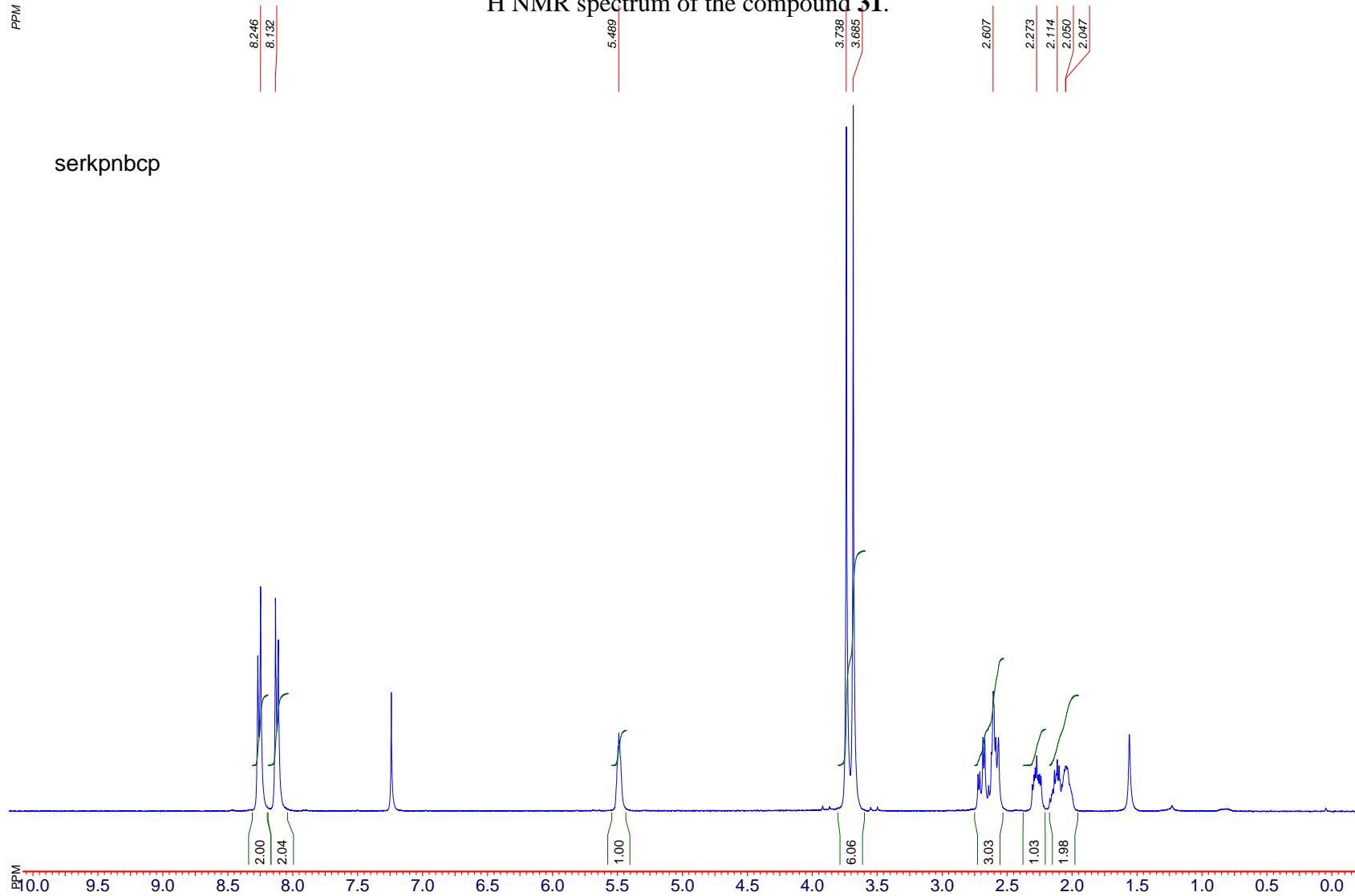
¹³C NMR spectrum of the compound **12·HCl**.



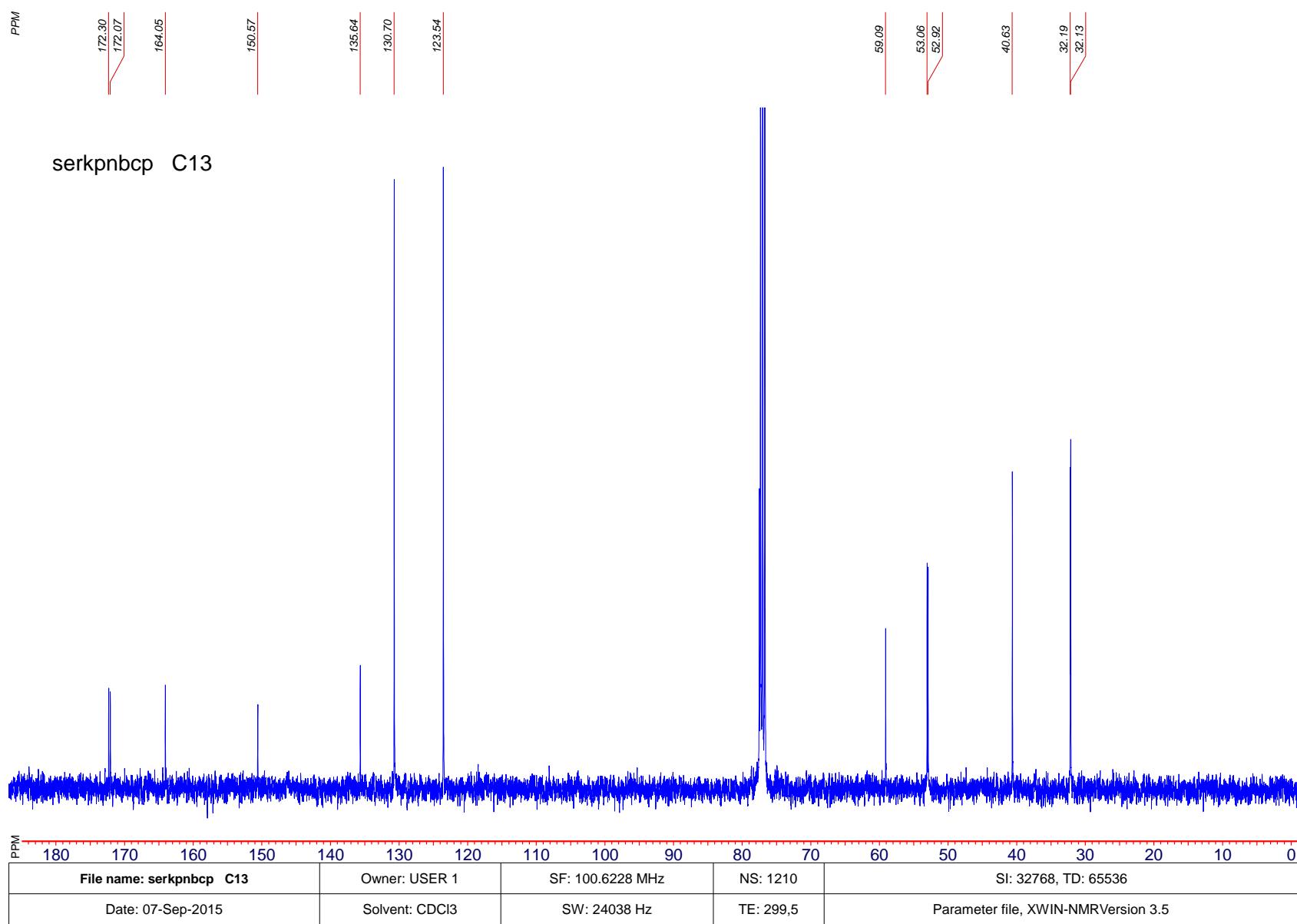
PPM

^1H NMR spectrum of the compound **31**.

serkpnbcp



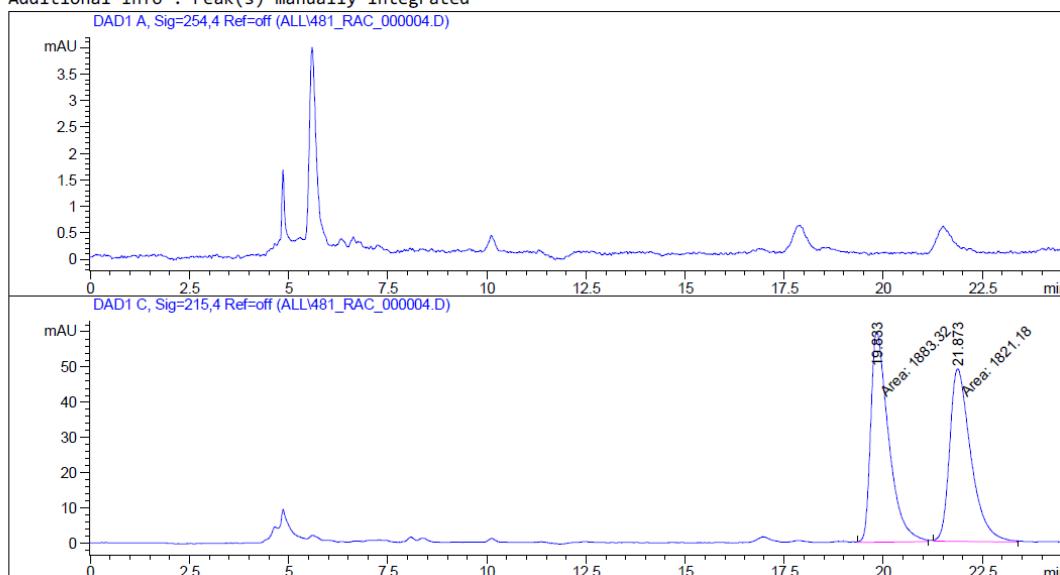
¹³C NMR spectrum of the compound **31**.



Chiral stationary phase HPLC trace of racemic **52**

Acq. Operator :
 Acq. Instrument : Instrument 1 Location : Vial 51
 Injection Date : 8/14/2015 11:19:22 AM Inj Volume : 2.0 μ l
 Acq. Method : C:\CHEM32\1\METHODS\HEXANE-IPA 90-10.M
 Last changed : 8/14/2015 11:44:02 AM
 (modified after loading)
 Analysis Method : C:\CHEM32\1\METHODS\HEXANE-IPA 90-10.M
 Last changed : 8/14/2015 1:00:41 PM
 (modified after loading)
 Sample Info : IA, Hexane-IPA 97-3 0.7 ml/min

Additional Info : Peak(s) manually integrated



=====

Area Percent Report

=====

Sorted By : Signal
 Multiplier: : 1.0000
 Dilution: : 1.0000
 Use Multiplier & Dilution Factor with ISTDs
 Signal 1: DAD1 A, Sig=254,4 Ref=off

Signal 2: DAD1 C, Sig=215,4 Ref=off

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	19.833	MM T	0.5243	1883.32080	59.87222	50.8387
2	21.873	MM T	0.6189	1821.18469	49.04562	49.1613

Totals : 3704.50549 108.91784

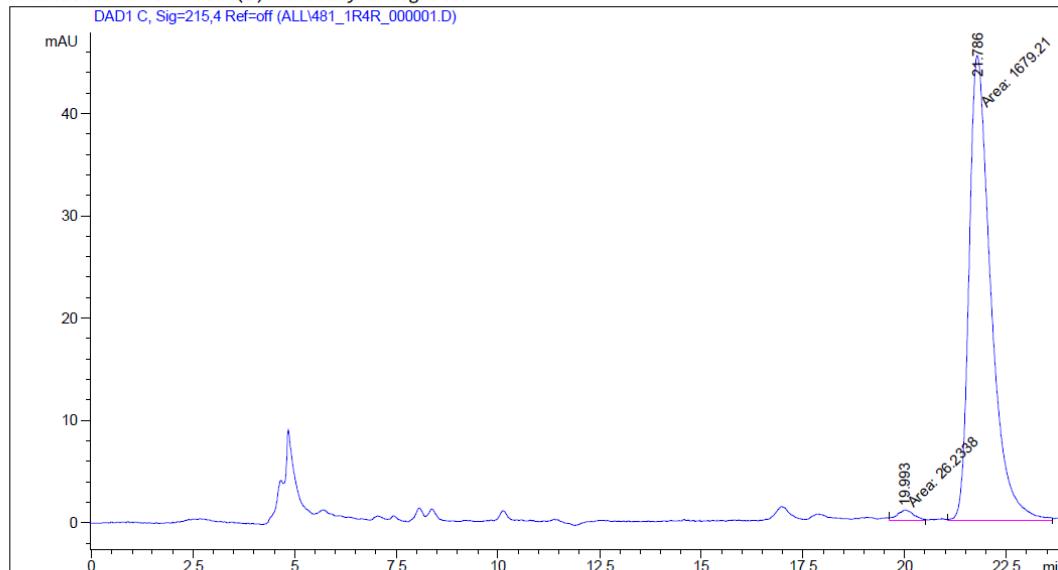
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*** End of Report ***

Chiral stationary phase HPLC trace of (*R,R*)-52

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 52
Injection Date : 8/14/2015 11:49:49 AM Inj Volume : 2.0 μ l
Acq. Method : C:\CHEM32\1\METHODS\HEXANE-IPA 90-10.M
Last changed : 8/14/2015 11:49:36 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\HEXANE-IPA 90-10.M
Last changed : 8/14/2015 1:00:41 PM
(modified after loading)
Sample Info : IA, Hexane-IPA 97-3 0.7 ml/min

Additional Info : Peak(s) manually integrated



===== Area Percent Report =====

Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
Signal 1: DAD1 C, Sig=215,4 Ref=off

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	19.993	MM T	0.4617	26.23381	9.46990e-1	1.5382
2	21.786	MM T	0.7039	1679.20972	45.36599	98.4618

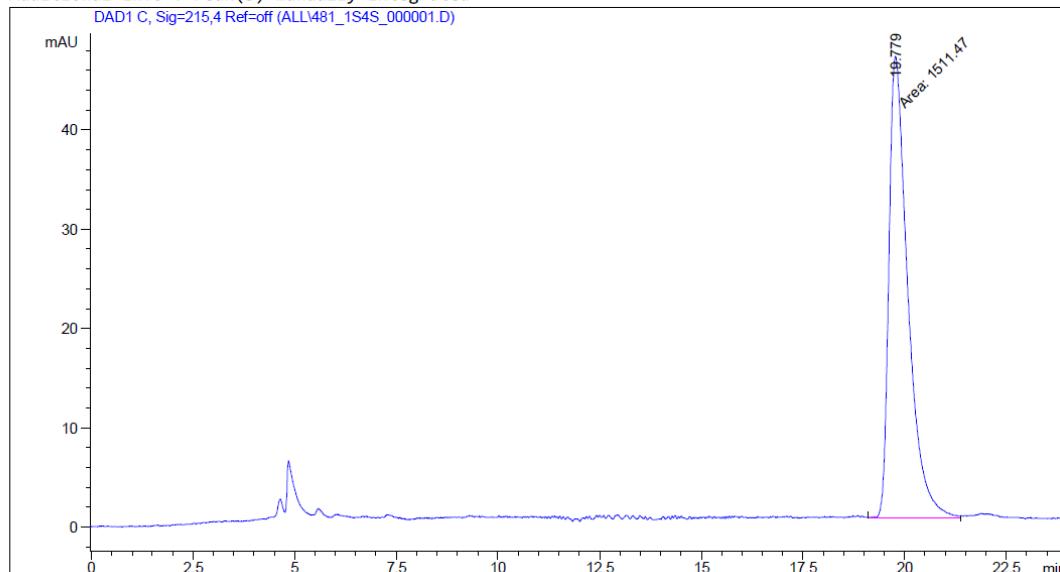
Totals : 1705.44353 46.31298

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*** End of Report ***

Chiral stationary phase HPLC trace of (S,S)-52

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 53
Injection Date : 8/14/2015 12:34:19 PM Inj Volume : 2.0 μ l
Acq. Method : C:\CHEM32\1\METHODS\HEXANE-IPA 90-10.M
Last changed : 8/14/2015 12:14:03 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\HEXANE-IPA 90-10.M
Last changed : 8/14/2015 11:49:17 AM
(modified after loading)
Sample Info : IA, Hexane-IPA 97-3 0.7 ml/min

Additional Info : Peak(s) manually integrated



===== Area Percent Report =====

Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
Signal 1: DAD1 C, Sig=215,4 Ref=off

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	19.779	MM T	0.5413	1511.47327	46.53747	100.0000

Totals : 1511.47327 46.53747

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*** End of Report ***