

Synthesis and Evaluation of New Pyrazoline Derivatives as Potential Anticancer Agents in HepG-2 Cell Line

Wei jie Xu ^{1,†}, Ying Pan ^{1,†}, Hong Wang ^{1,†}, Haiyan Li ², Qing Peng ³, Duncan Wei ¹, Cheng Chen ¹ and Jinhong Zheng ^{1,*}

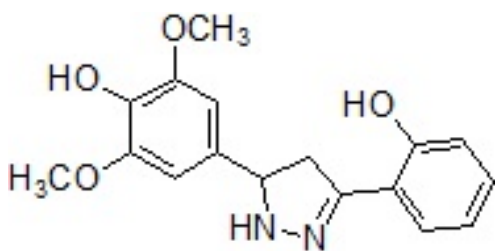
¹ Department of Chemistry, Shantou University Medical College, Shantou 515041, Guangdong, China; 13536920534@163.com (W.X.); ypan@stu.edu.cn (Y.P.); 13592851552@163.com (H.W.); weiduncan2012@163.com (D.W.); 13411963192@163.com (C.C.); jhzheng@stu.edu.cn (J.Z)

² Department of Pharmacology, Shantou University Medical College, Shantou 515041, Guangdong, China; 13066360950@163.com (H.L.)

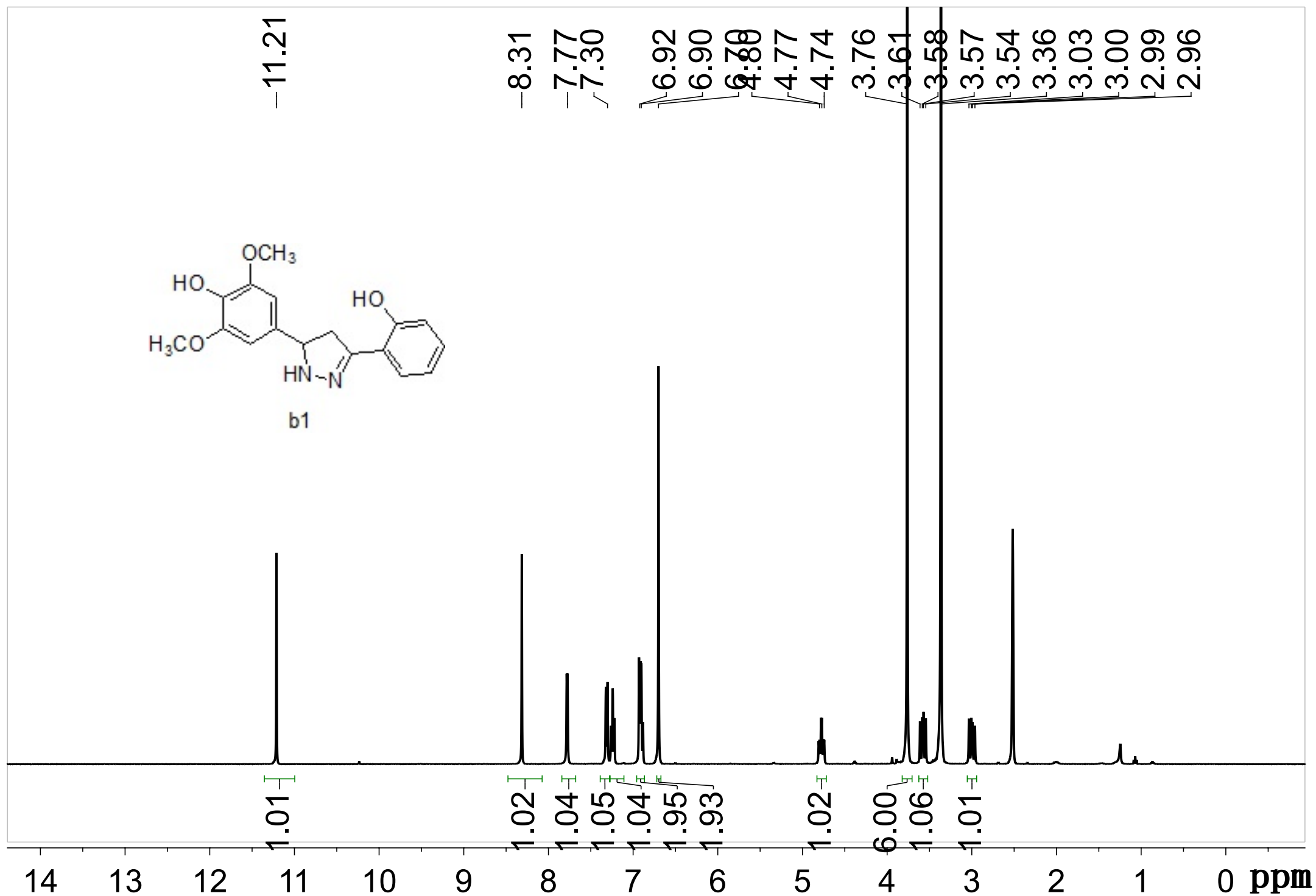
³ Department of Hepatobiliary Surgery II, Zhujiang Hospital of Southern Medical University, Guangzhou 510280, Guangdong, China; zhujiangzhuanhua@163.com (Q.P.)

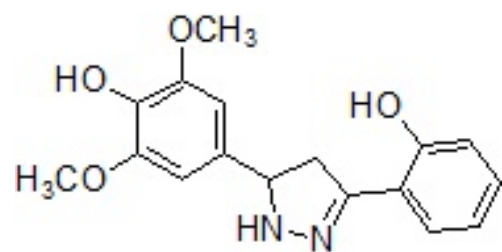
* Correspondence: jhzheng@stu.edu.cn; Tel.: +86-754-8890-0499; Fax: +86-754-8855-7562

† These authors contributed equally to this work.

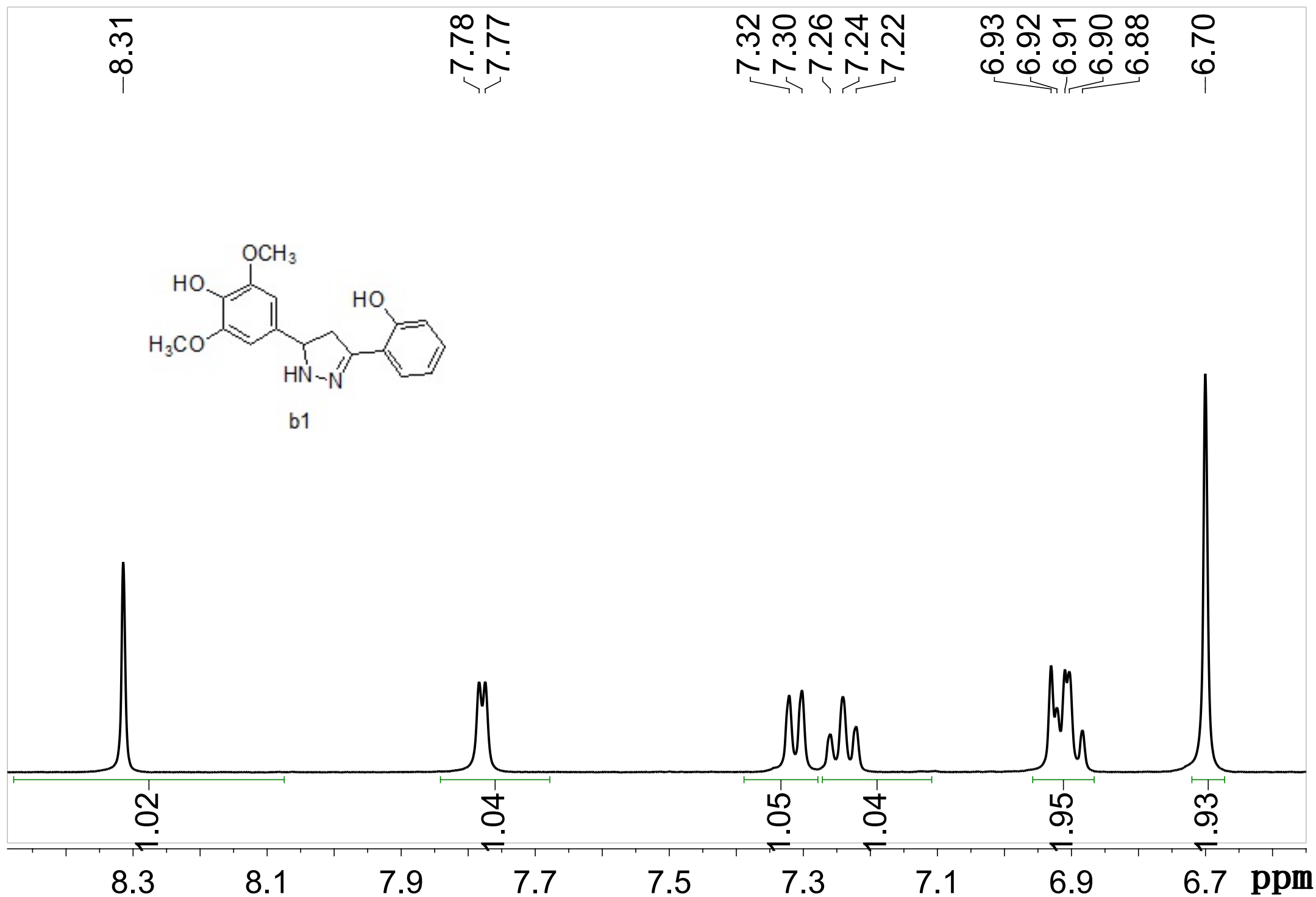


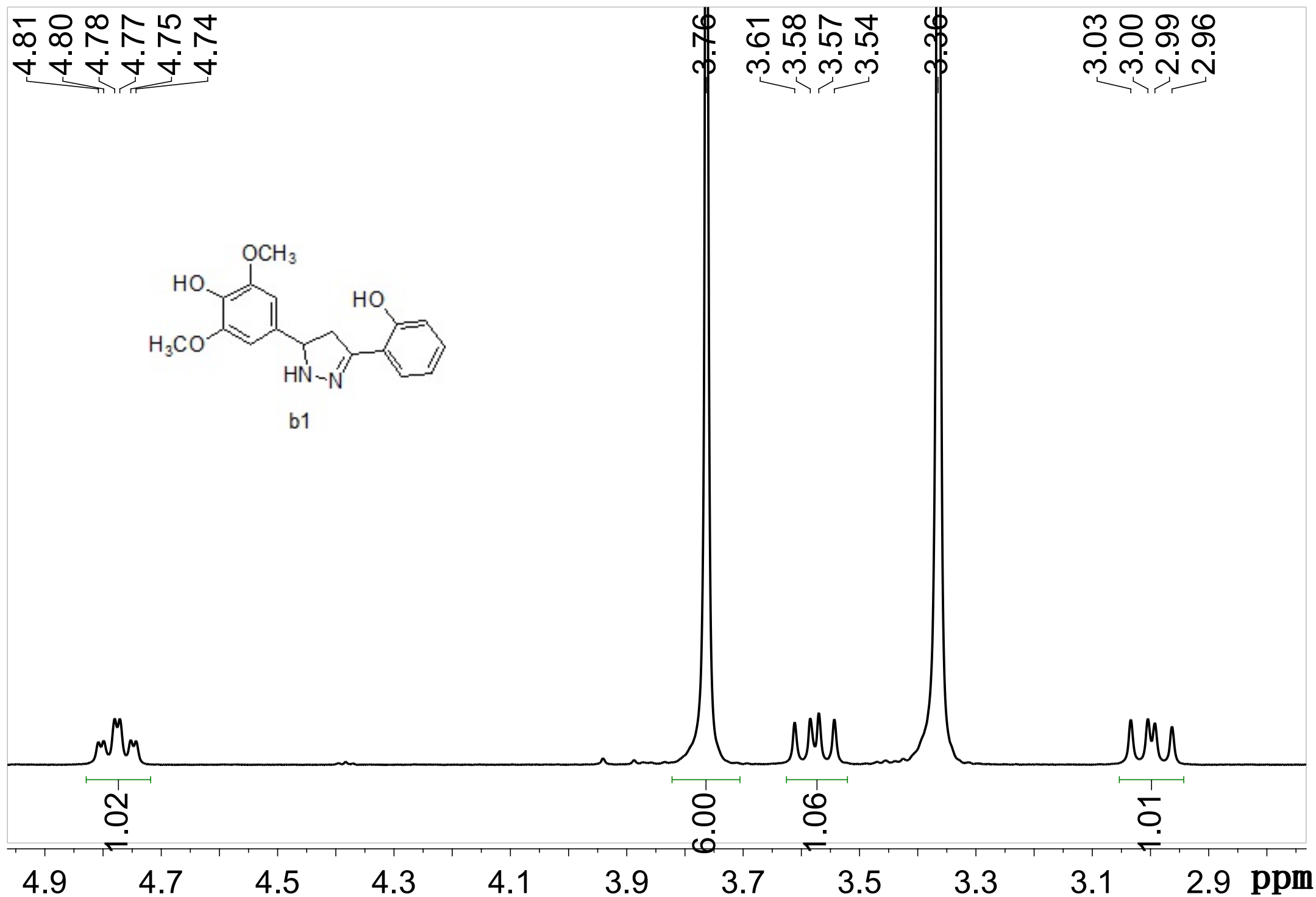
b1

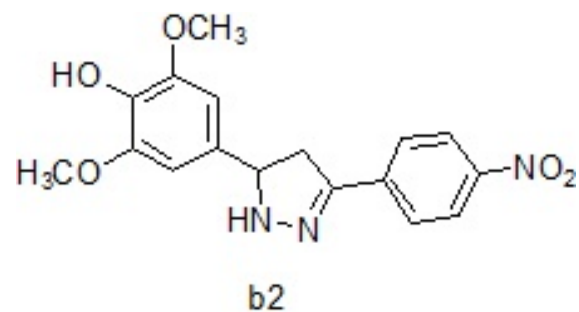




b1







Chemical shift values (ppm) for the aromatic region:

- 8.32
- 8.24
- 8.23
- 8.22
- 8.21
- 8.21
- 8.19
- 8.19
- 7.83
- 7.83
- 7.81
- 7.81
- 7.80
- 6.65

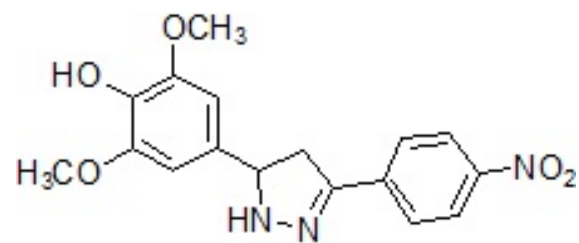
Chemical shift values (ppm) for the aliphatic region:

- 4.91
- 4.88
- 4.85
- 3.74
- 3.50
- 3.47
- 3.45
- 3.38
- 3.38
- 2.96
- 2.93
- 2.92
- 2.89

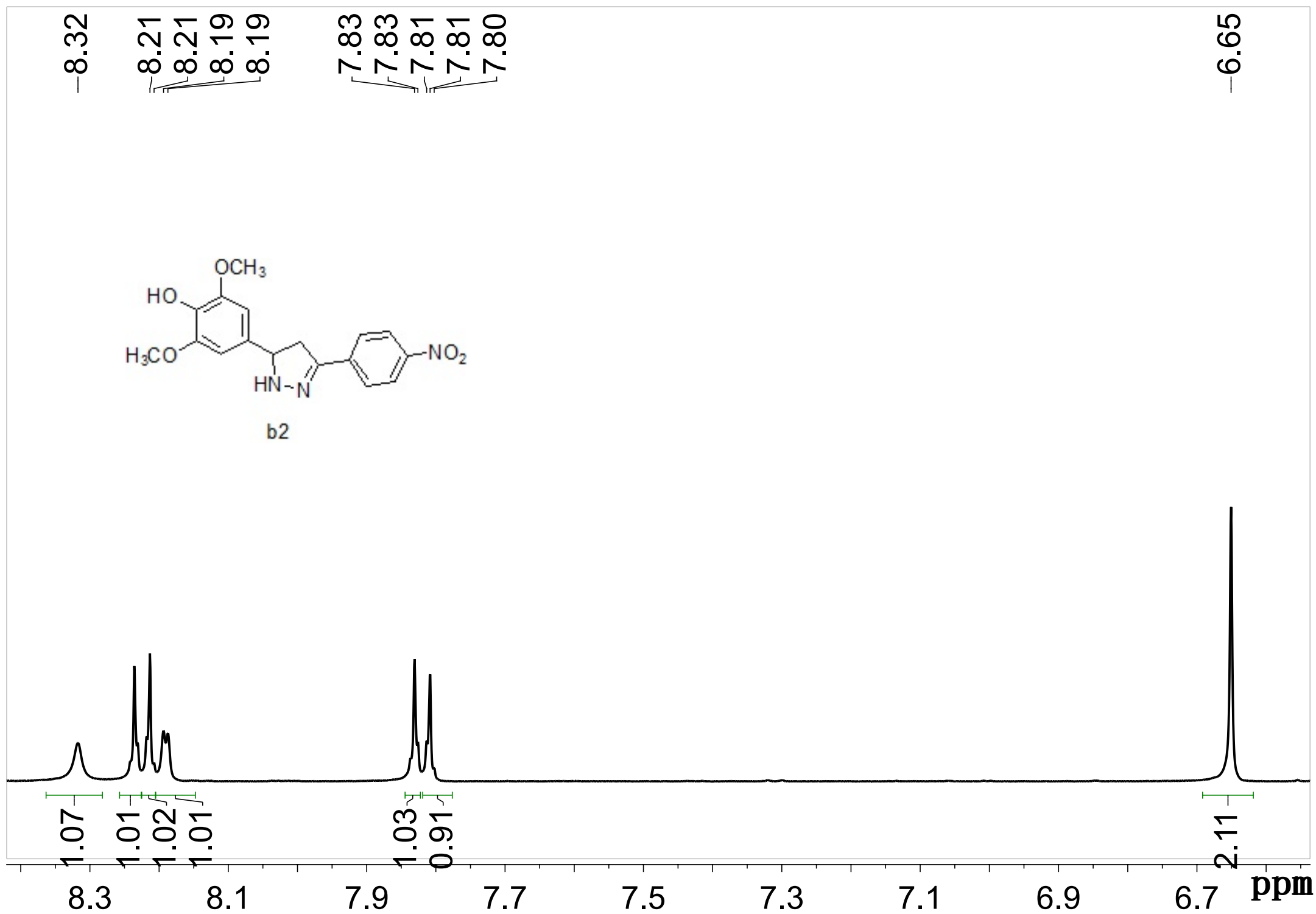
Integration values for the peaks:

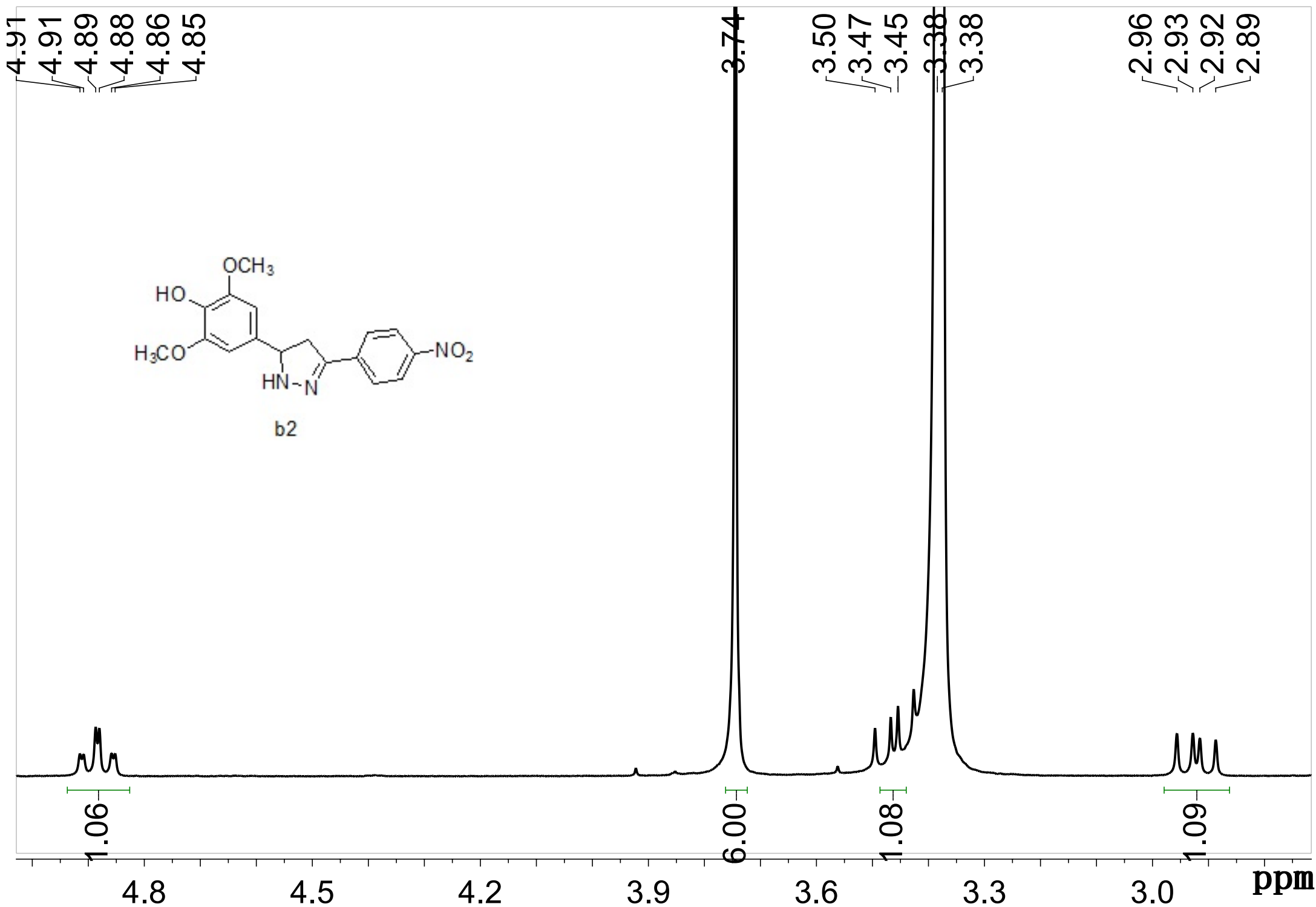
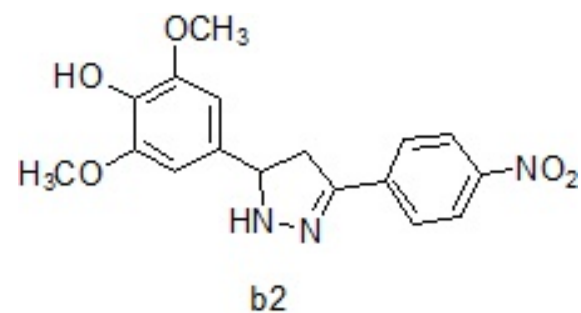
- 1.07
- 1.01
- 1.02
- 1.01
- 1.03
- 0.91
- 2.11
- 1.06
- 6.00
- 1.08
- 1.09

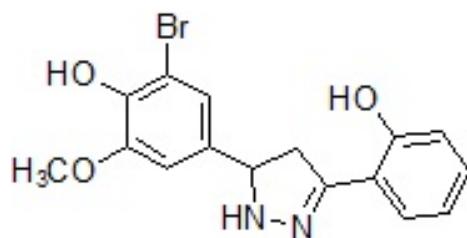
Chemical shift scale (ppm): 14, 13, 12, 11, 10, 9, 8, 7, 6, 5, 4, 3, 2, 1, 0 ppm



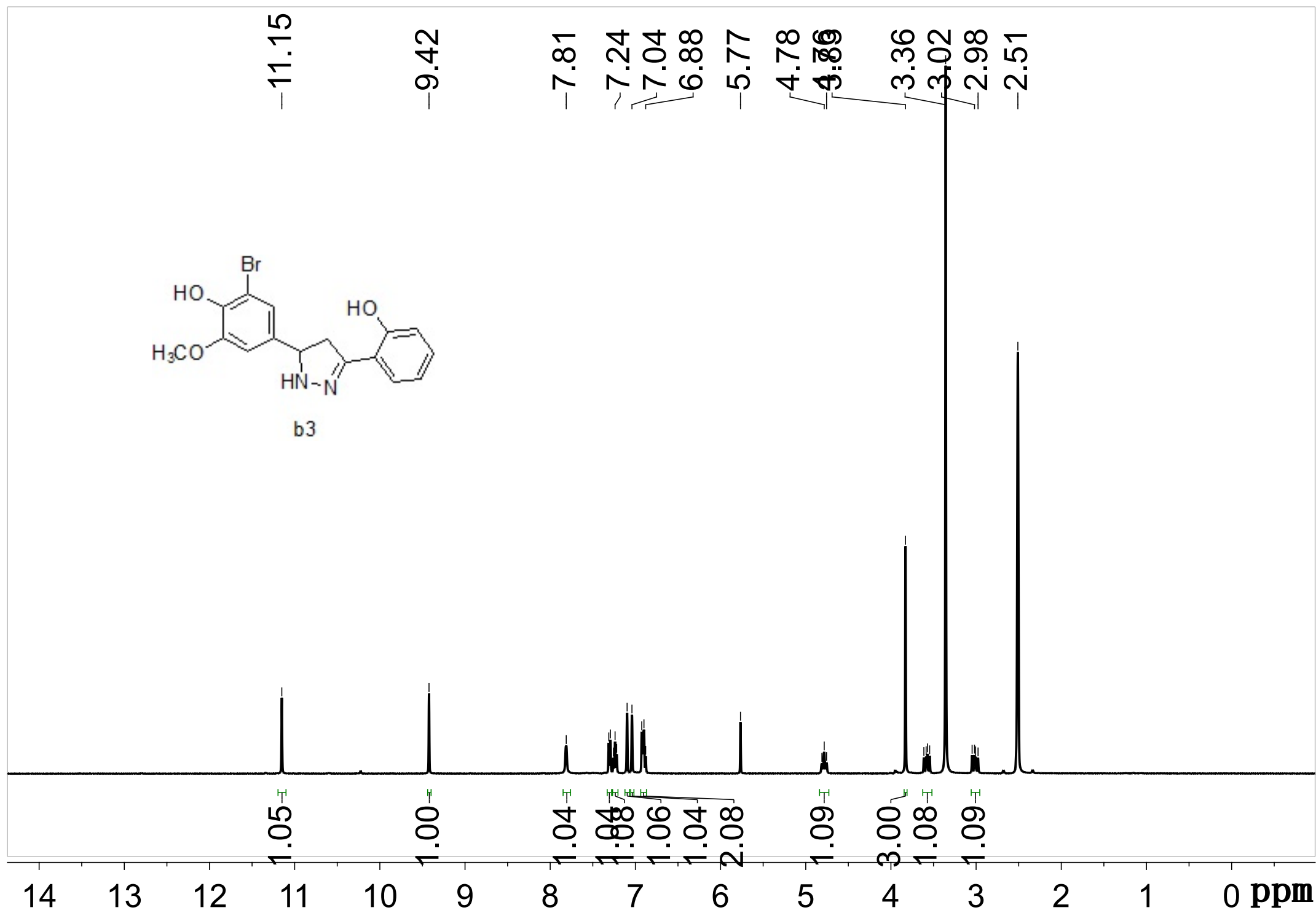
b2

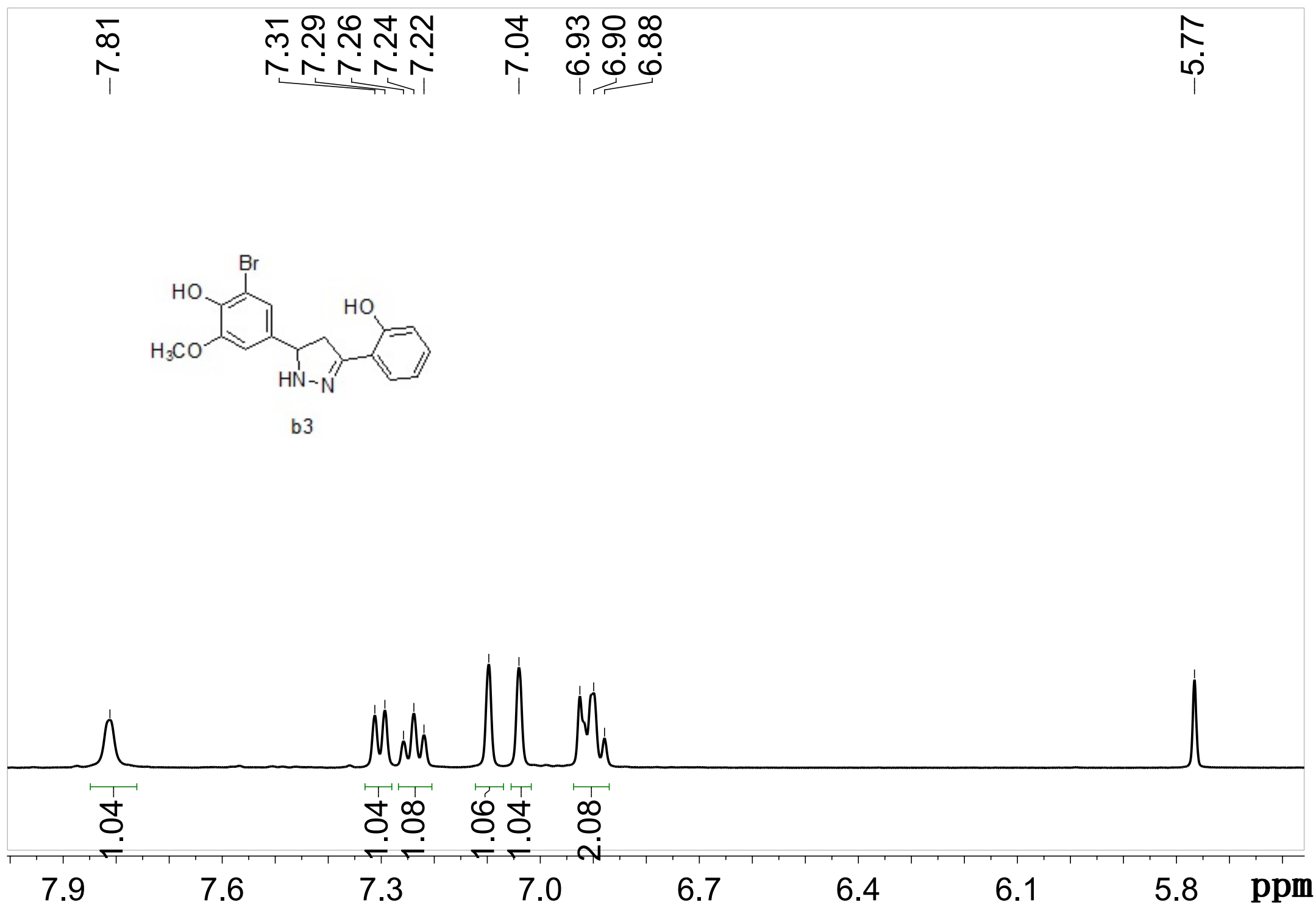


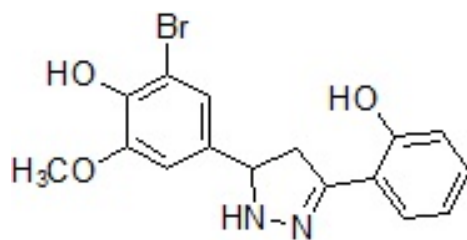




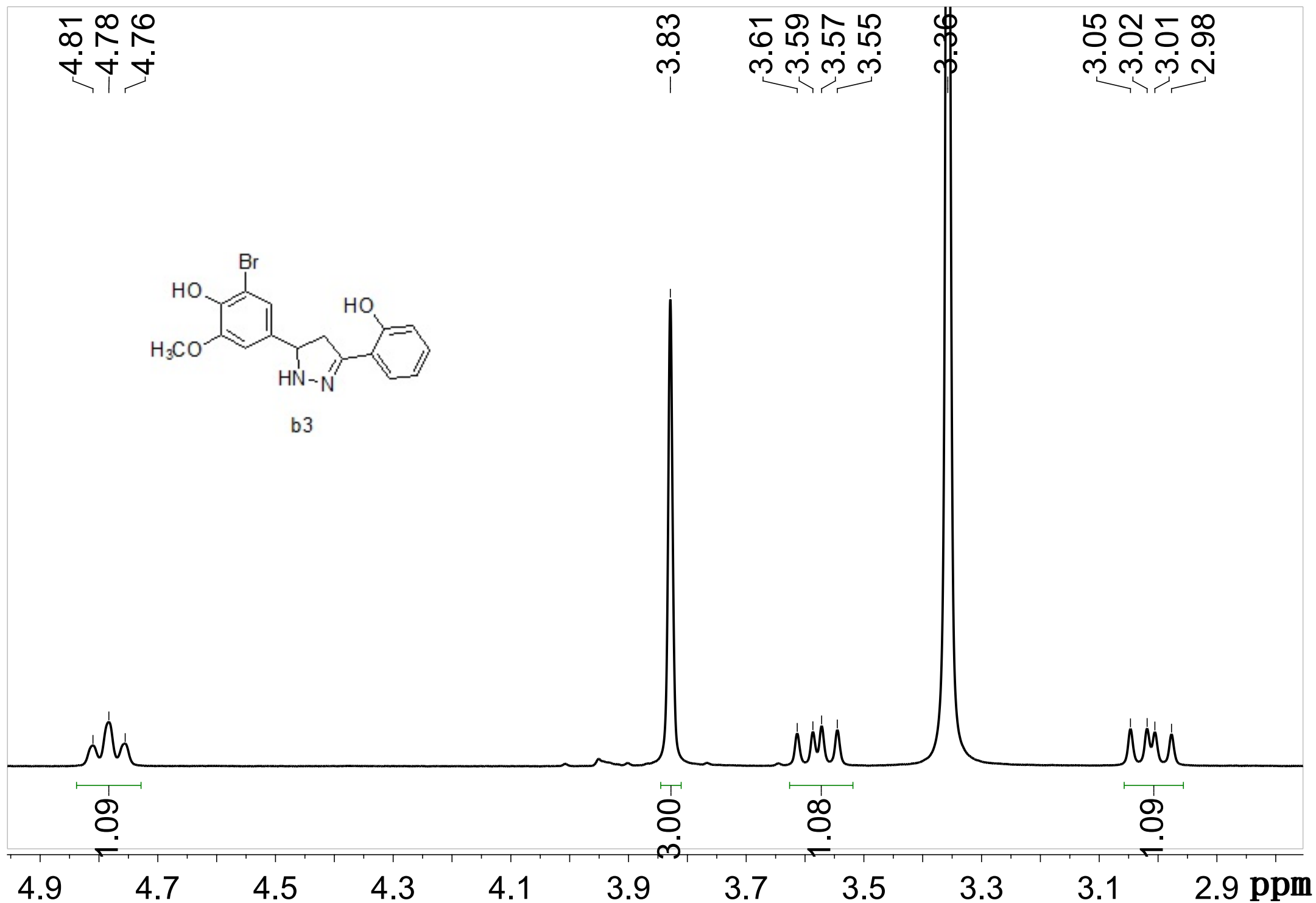
b3

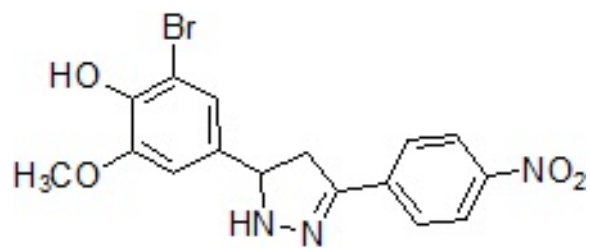




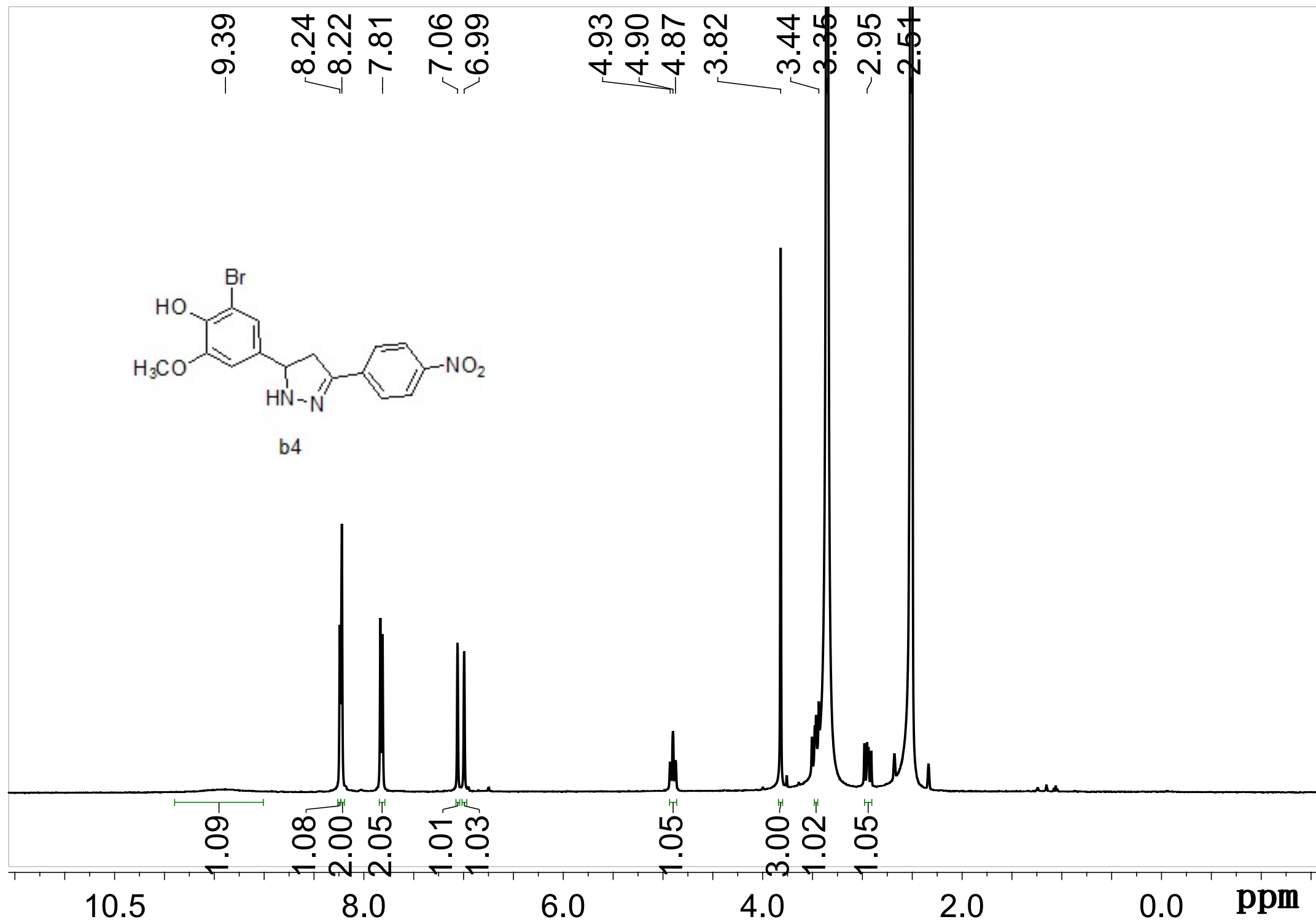


b3





b4

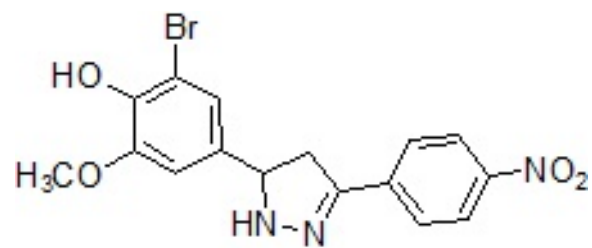


8.24
8.22

7.83
7.81

7.06
6.99

4.93
4.90
4.87



b4

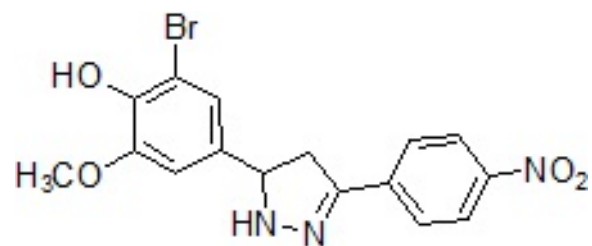
1.08
2.00

2.05

1.01
1.03

1.05

ppm



b4

3.82

3.48

3.46

3.44

3.35

2.98

2.95

2.94

2.91

3.00

1.02

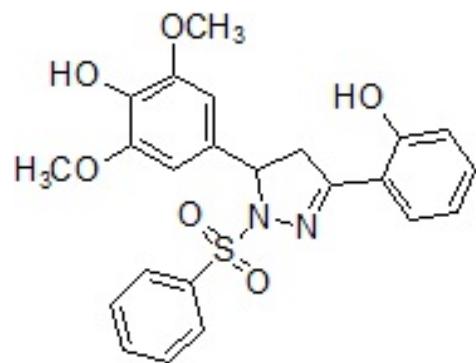
1.05

4.00

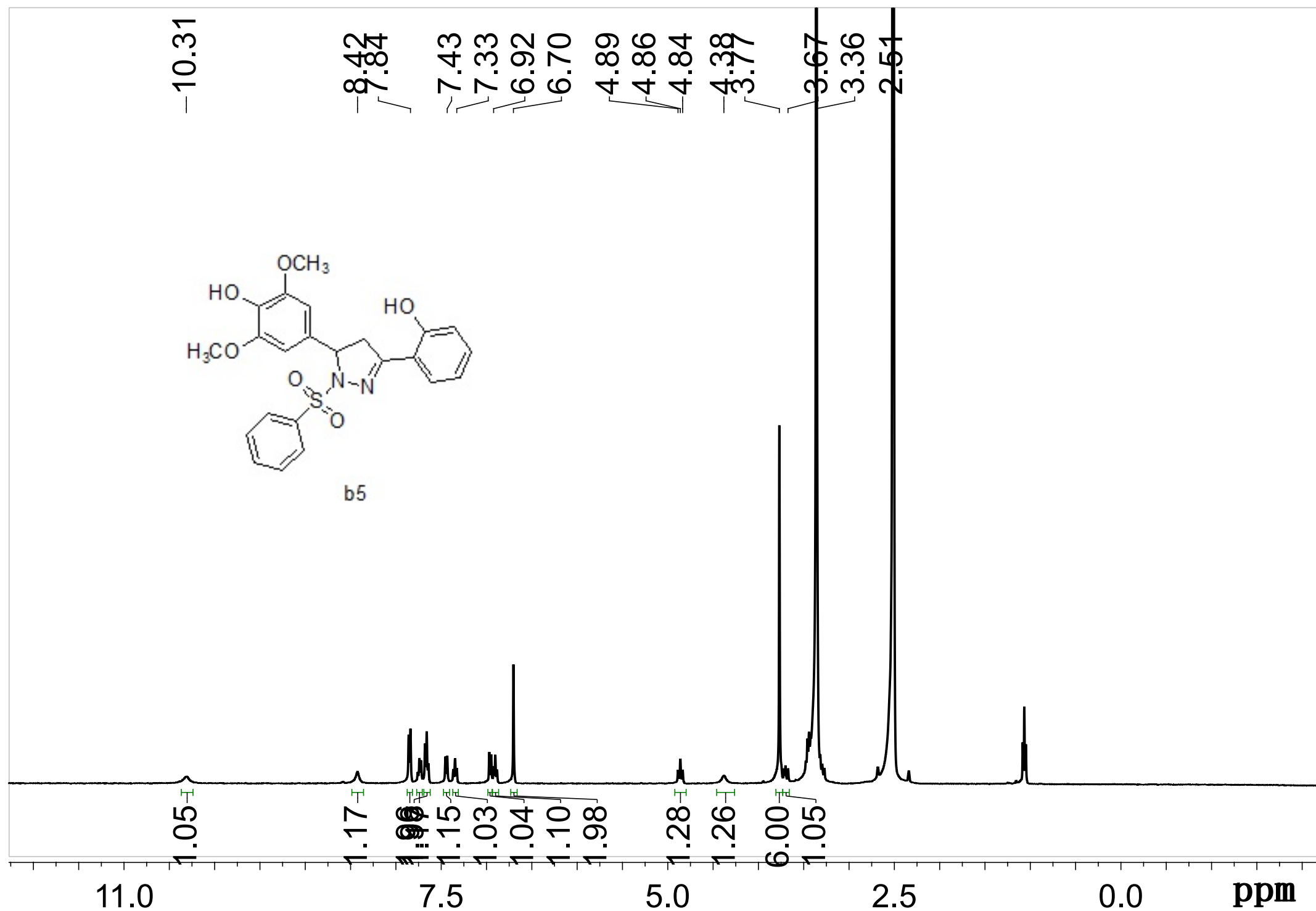
3.55

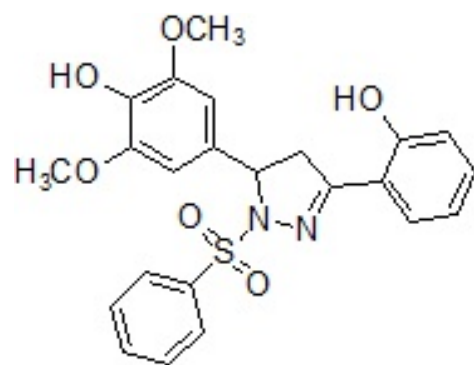
3.10

ppm



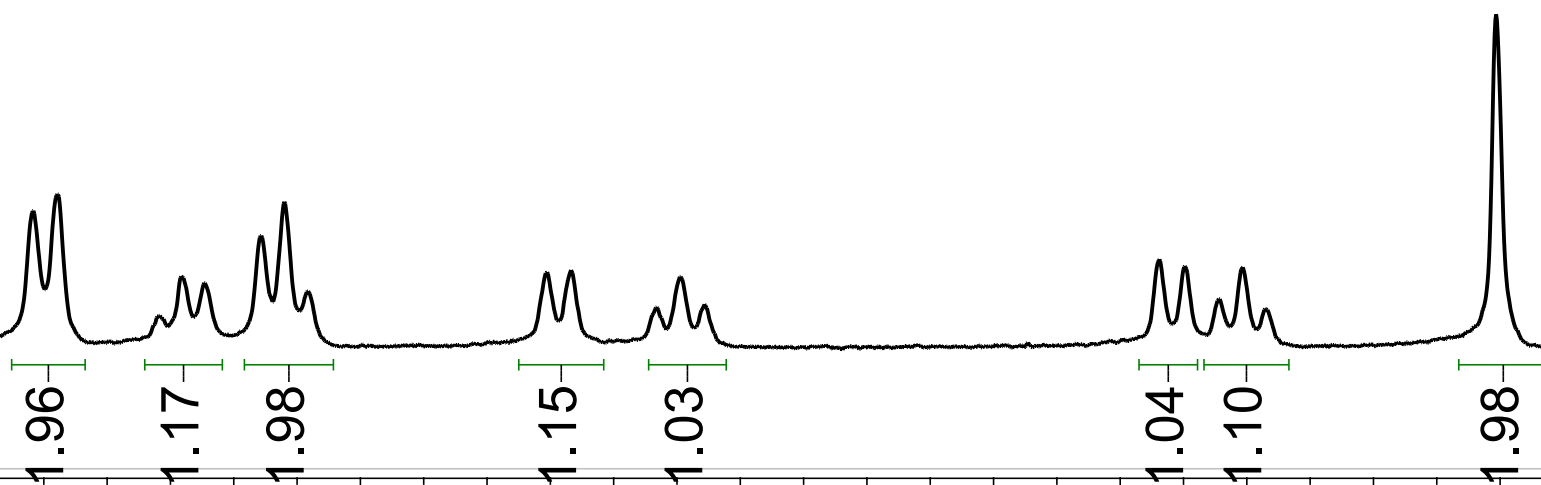
b5



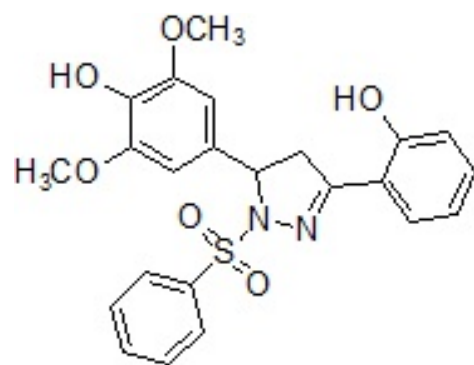


b5

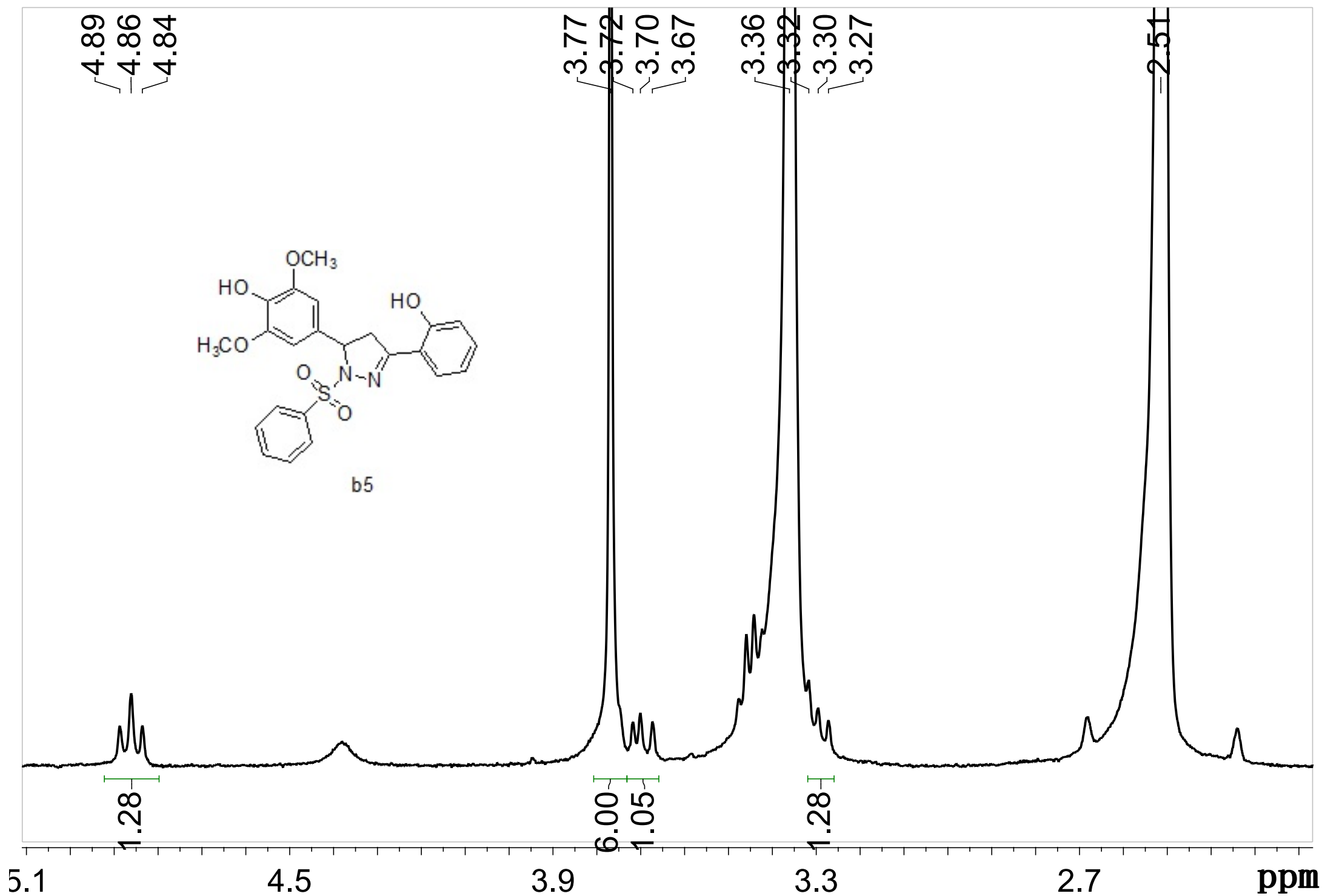
7.86
7.84
7.74
7.72
7.68
7.66
7.64
7.45
7.43
7.37
7.35
7.33
6.97
6.95
6.92
6.90
6.89
6.70

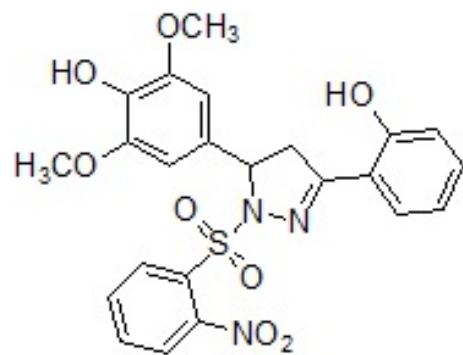


8.0 7.8 7.6 7.4 7.2 7.0 6.8 6.6 ppm



b5





b6

10.16

8.45

7.95

7.84

7.38

6.96

6.67

5.24

5.22

3.94

3.76

3.44

3.42

3.37

2.51

1.02

2.05

1.02

1.07

1.03

1.04

2.04

2.01

1.01

1.07

6.00

0.90

11.0

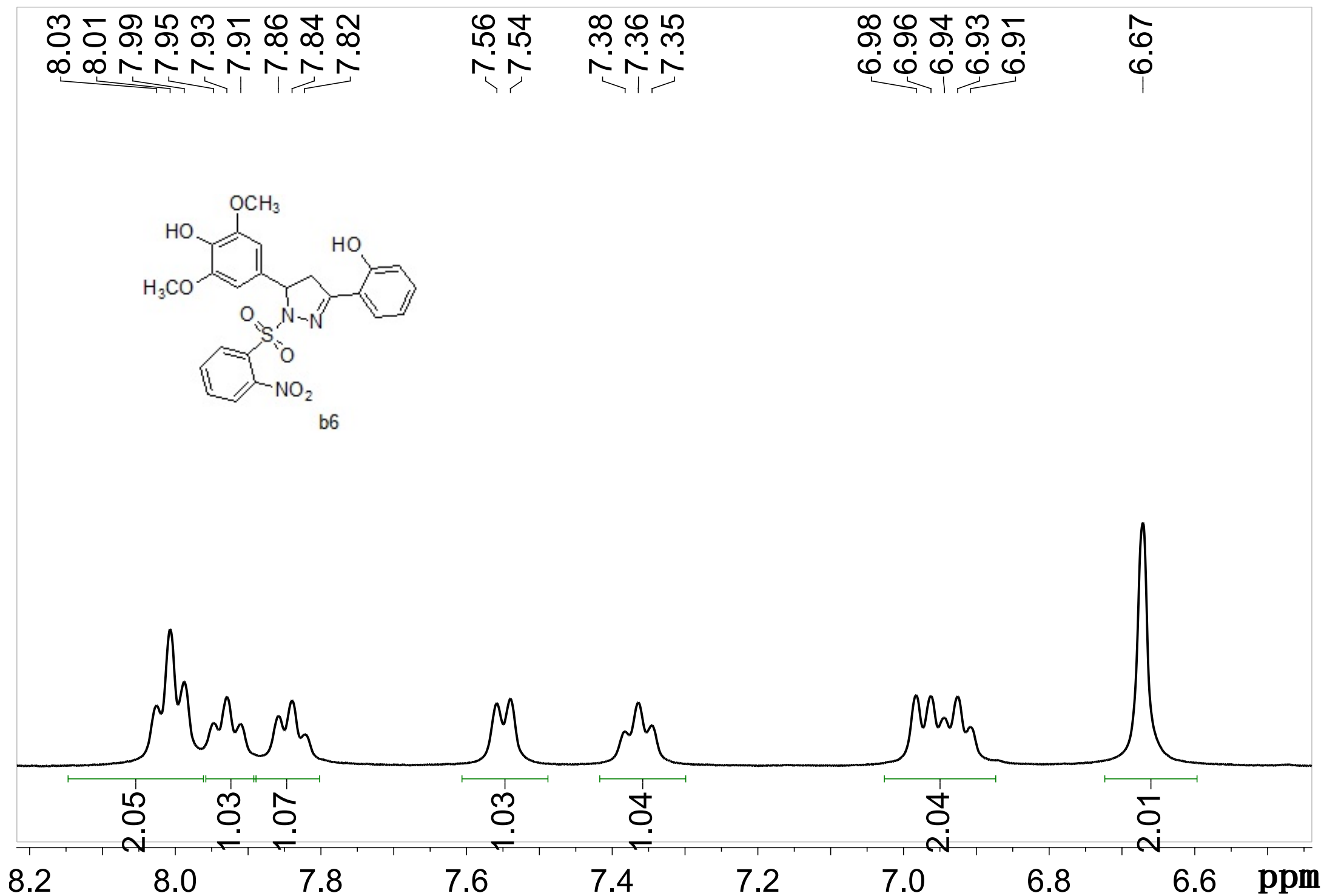
7.5

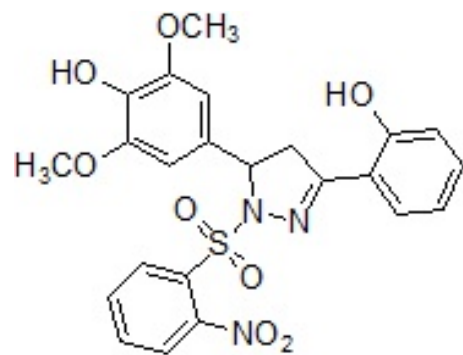
5.0

2.5

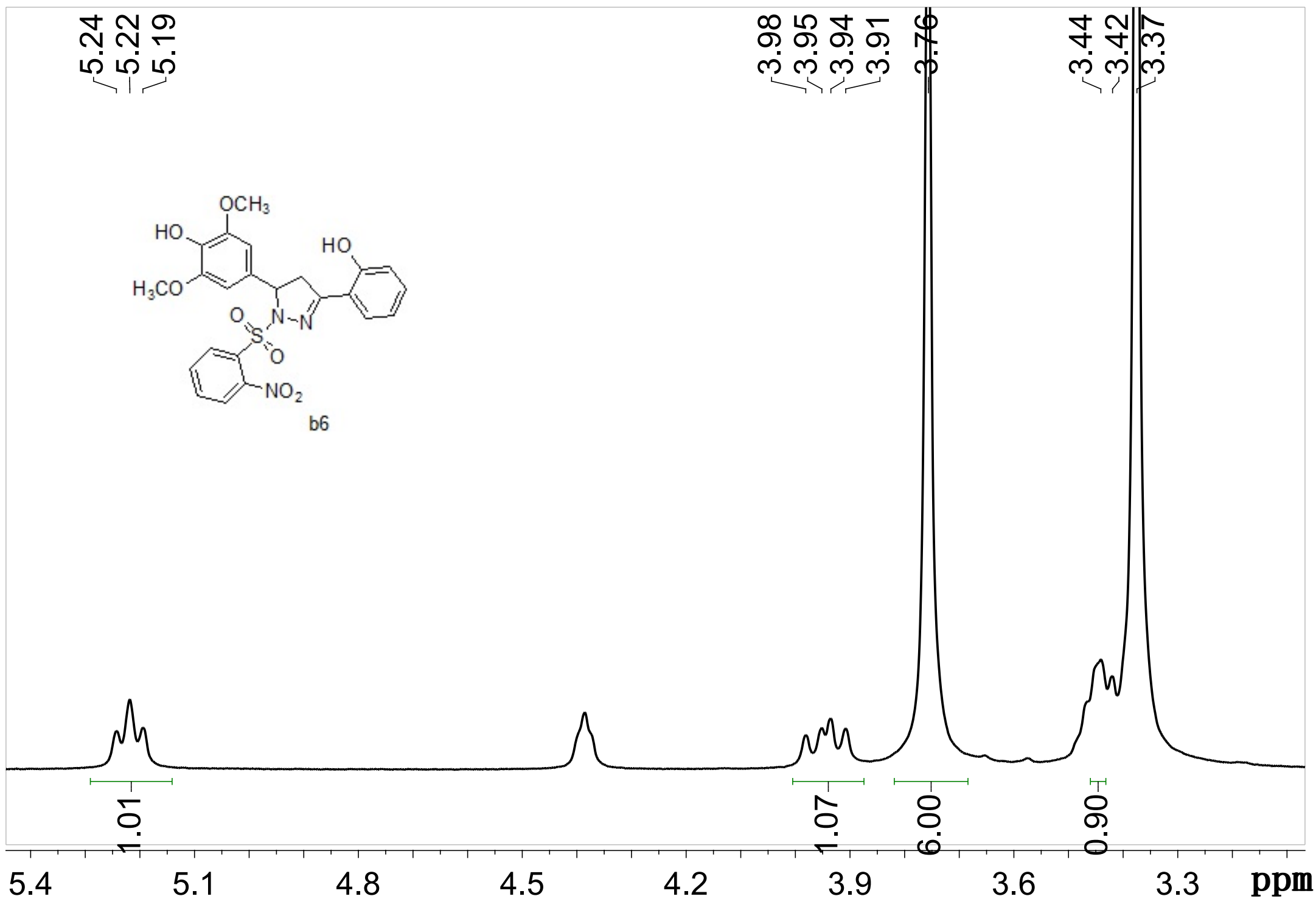
0.0

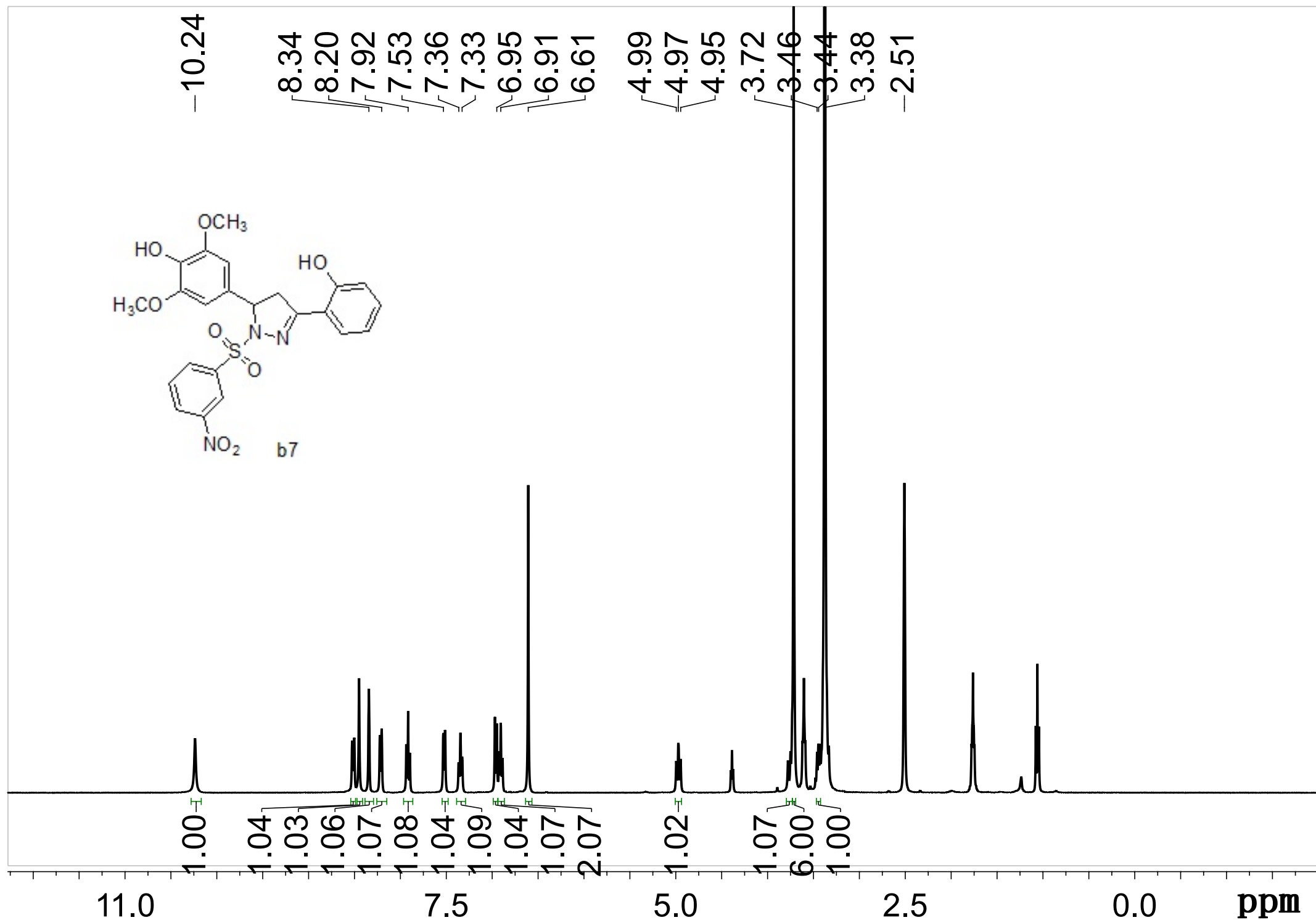
ppm

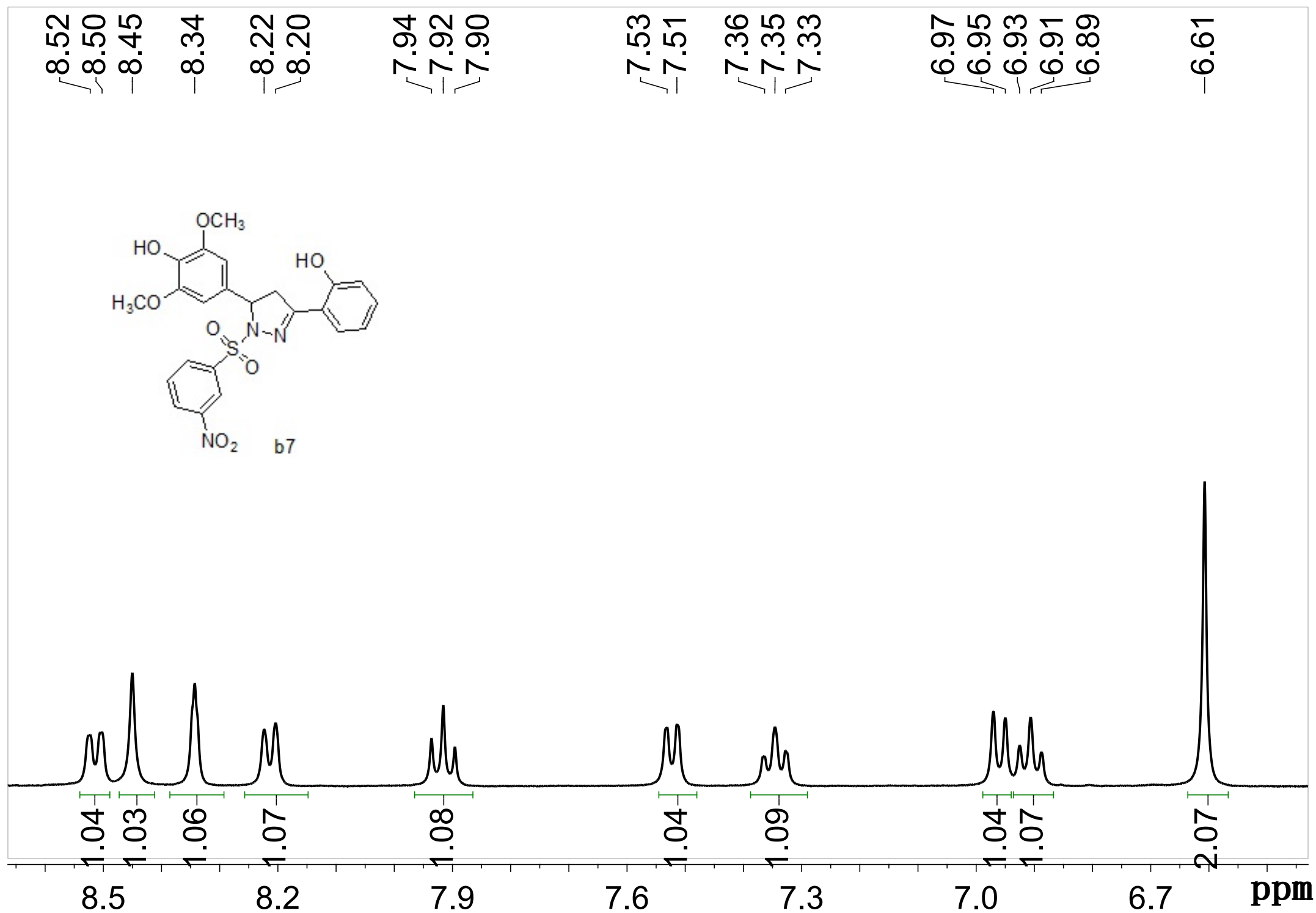


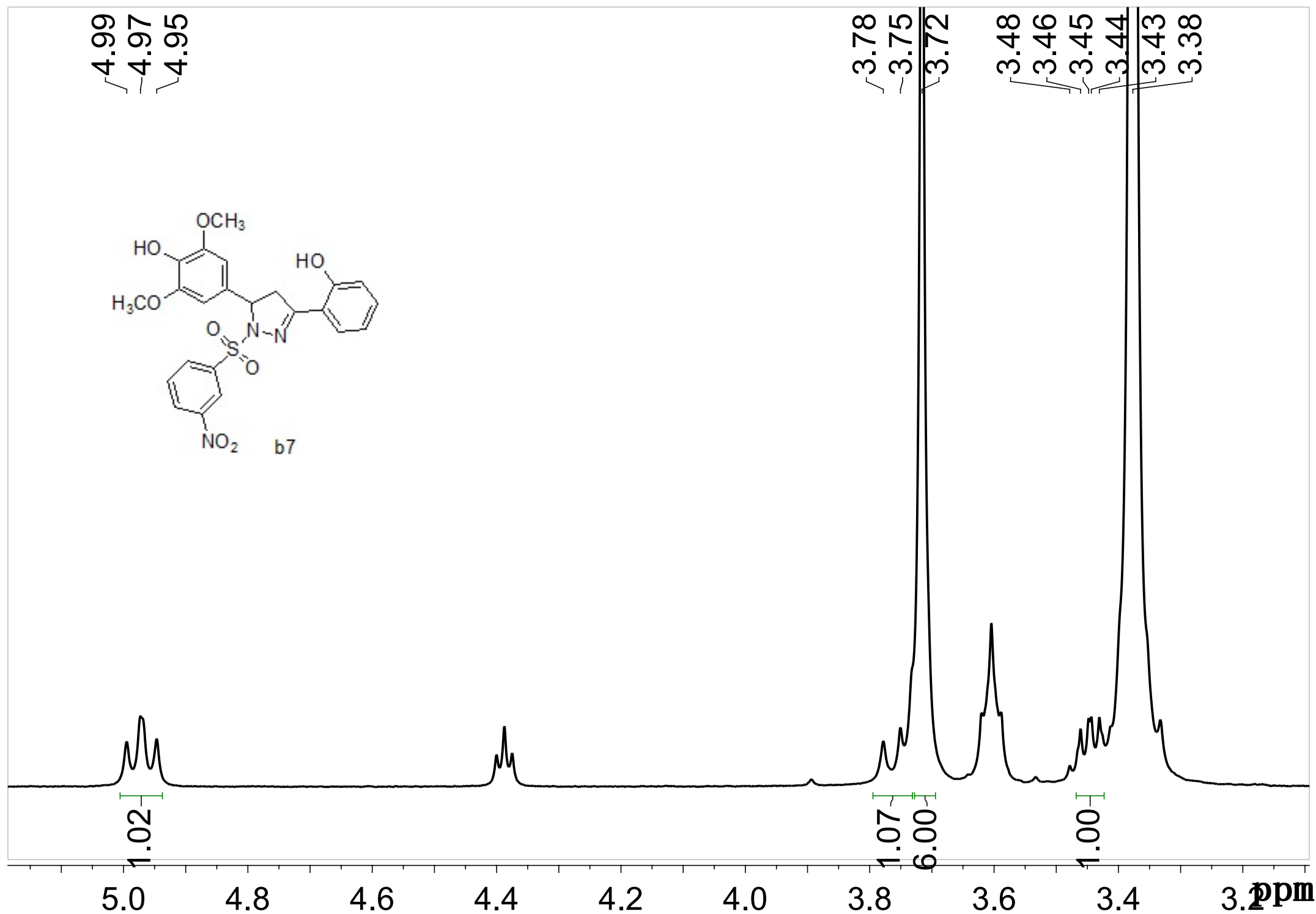


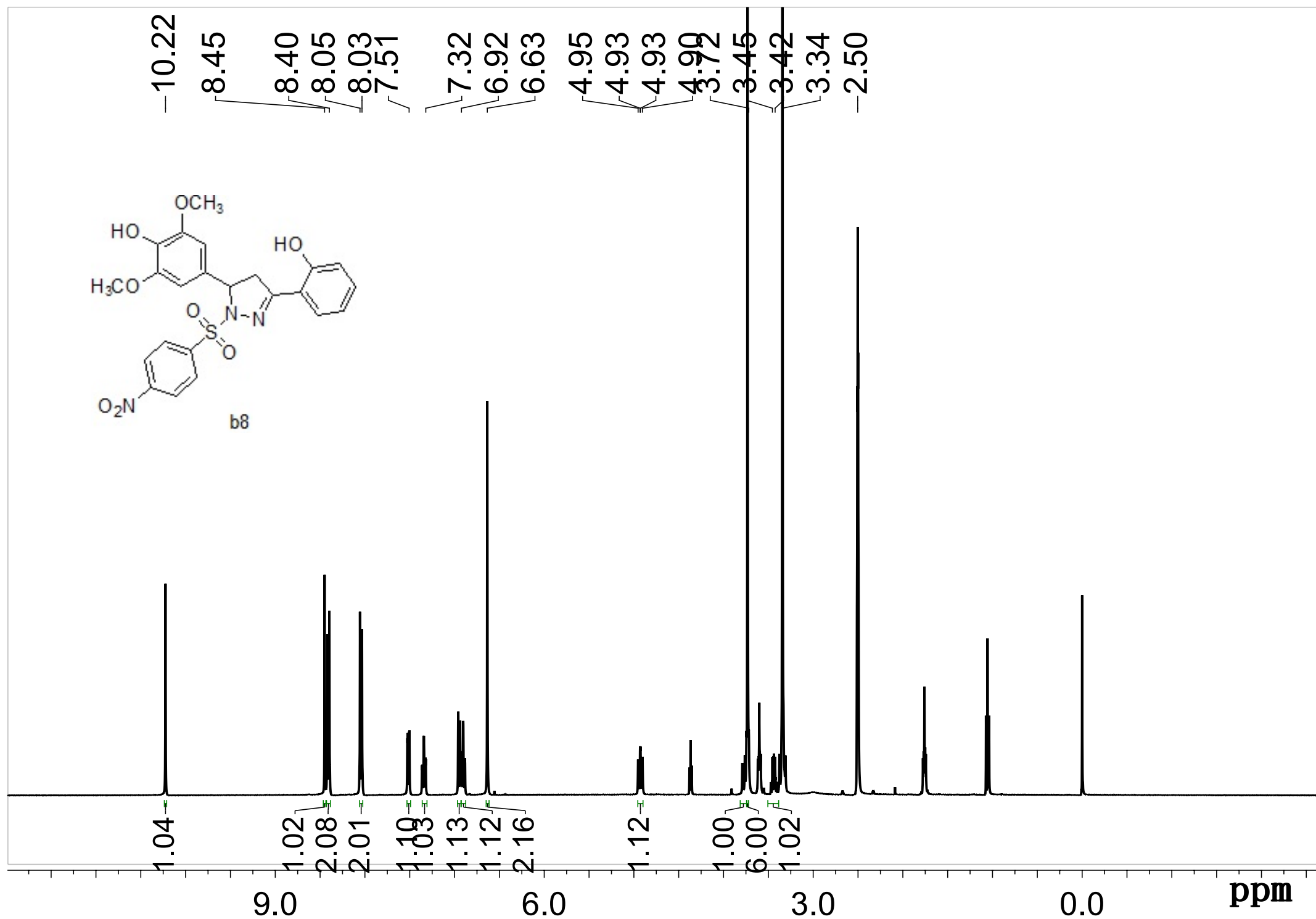
b6

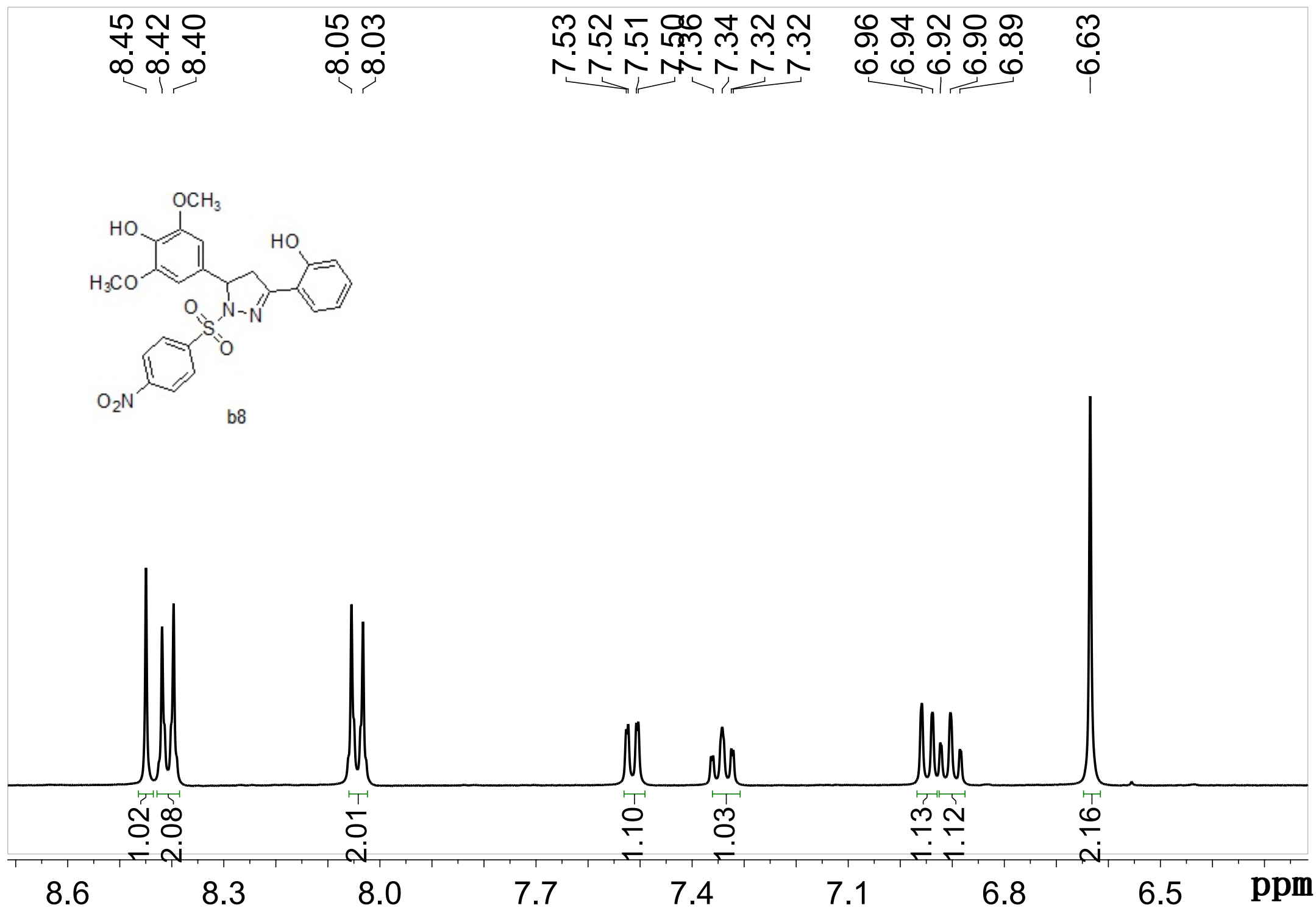


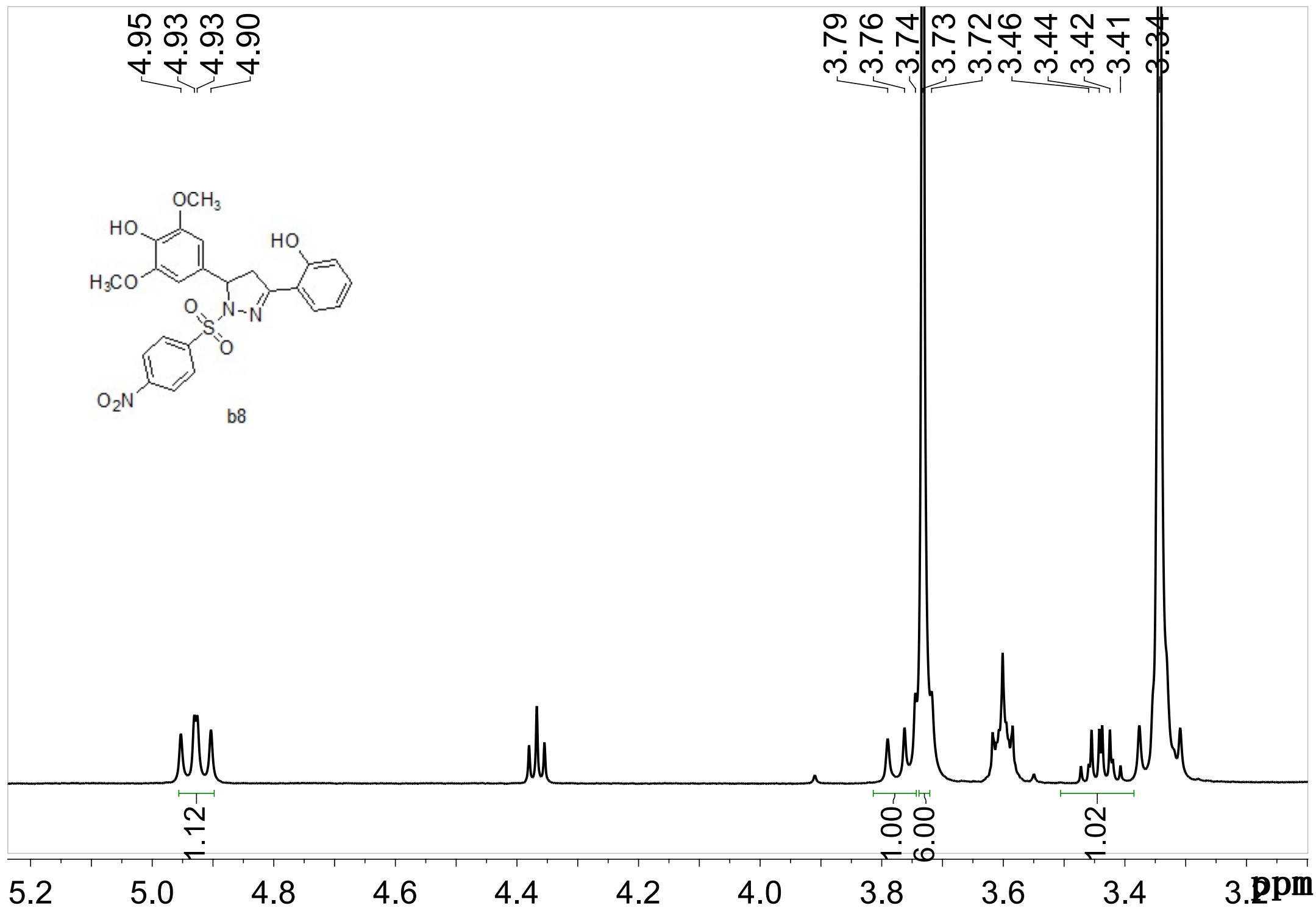


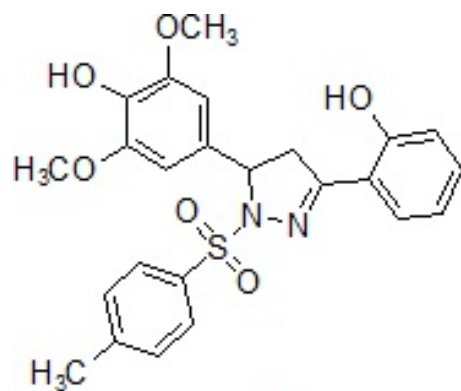




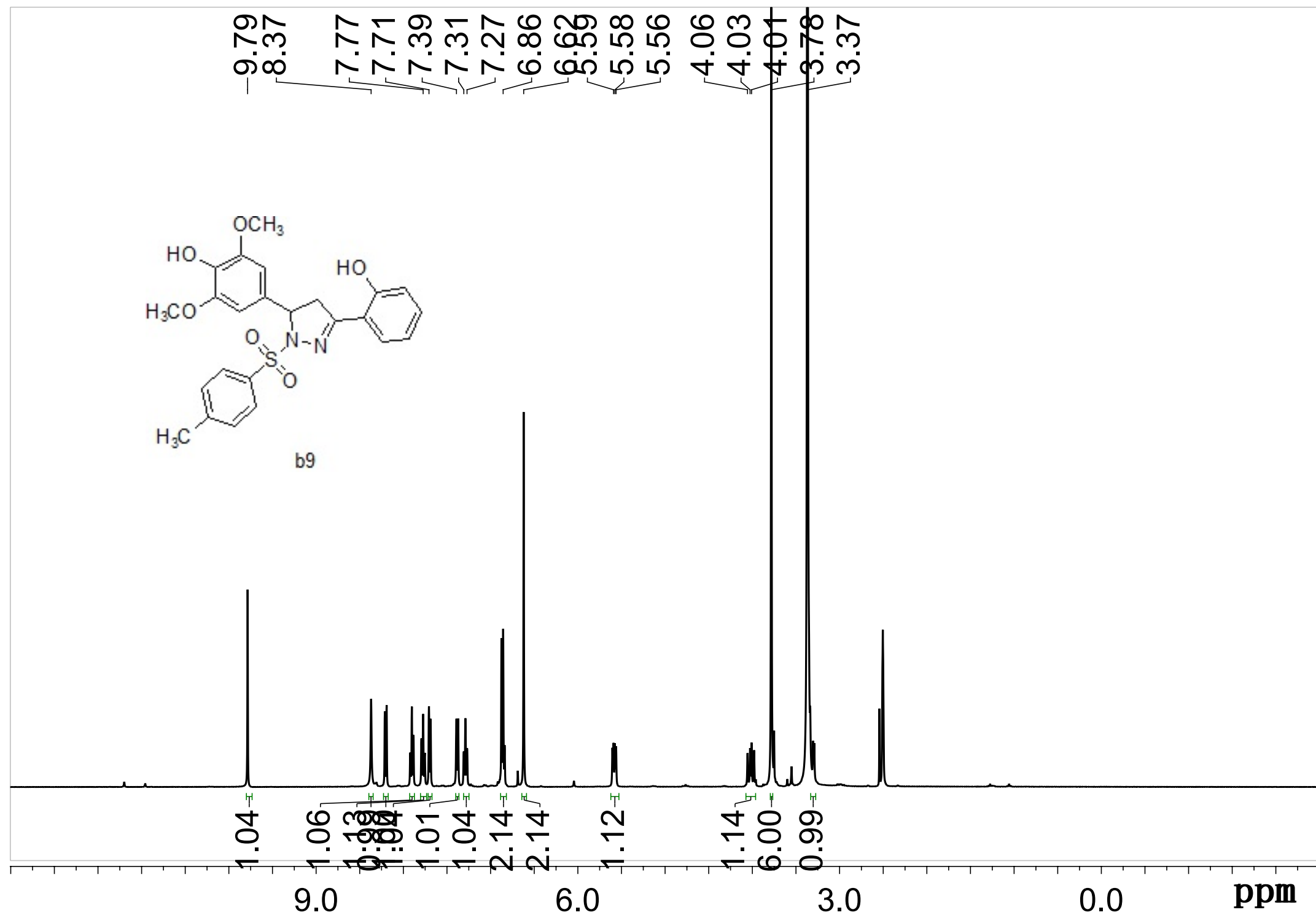


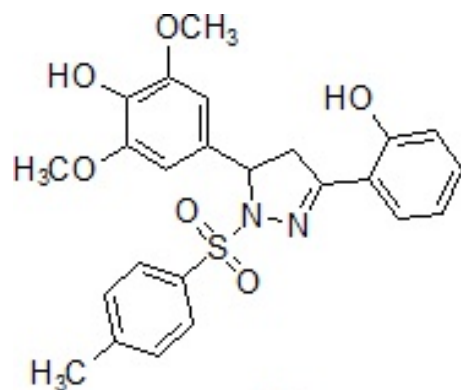




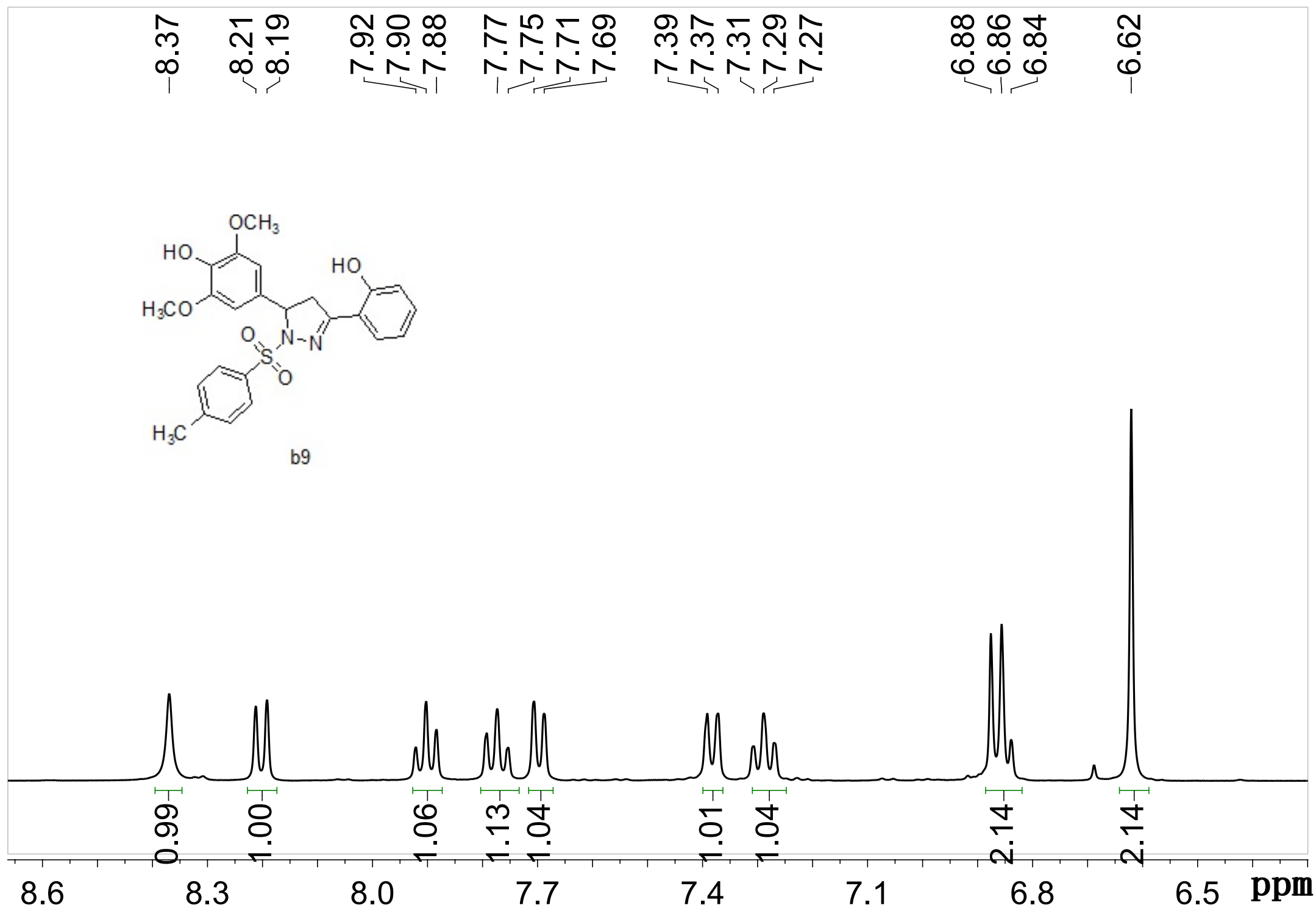


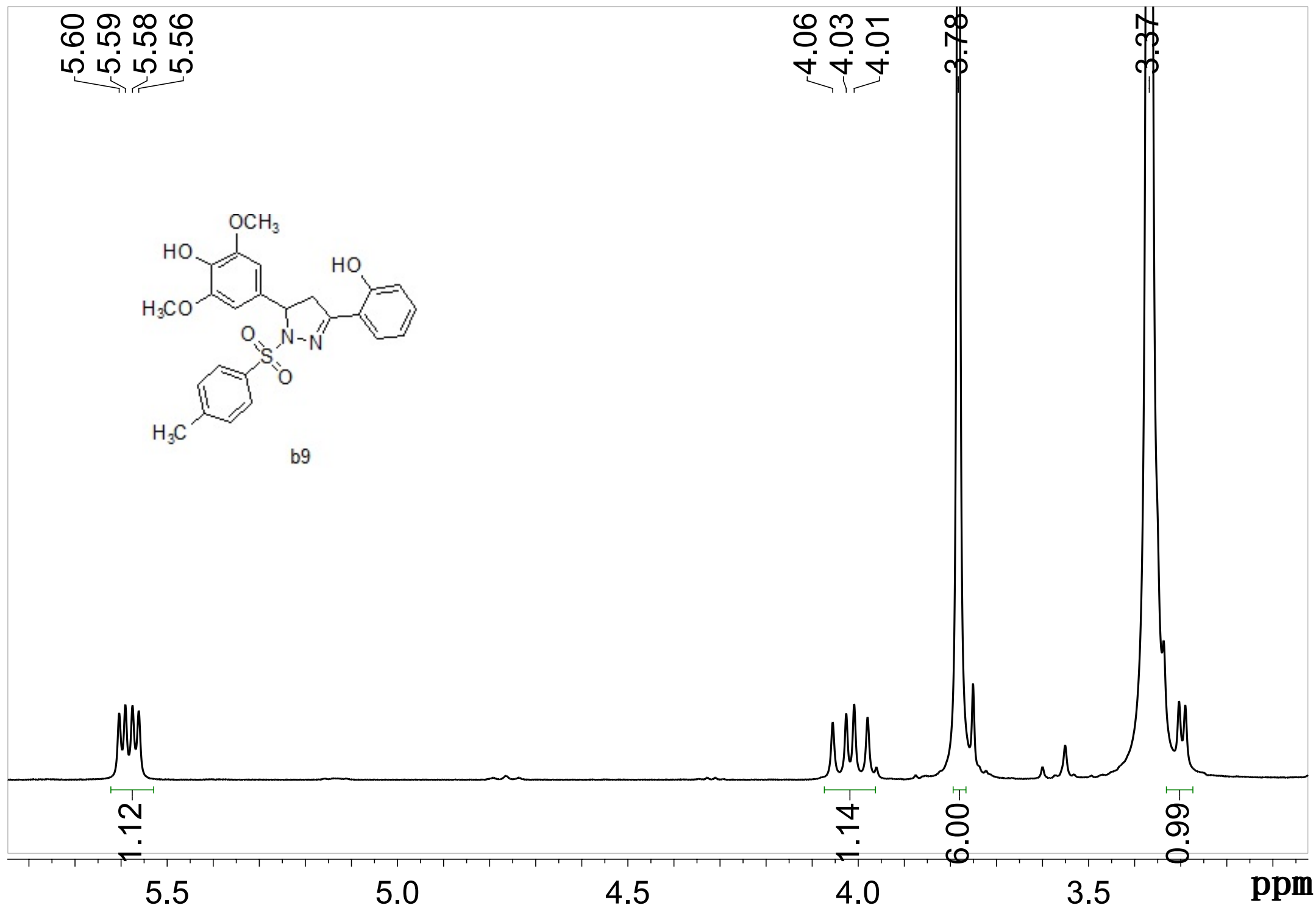
b9

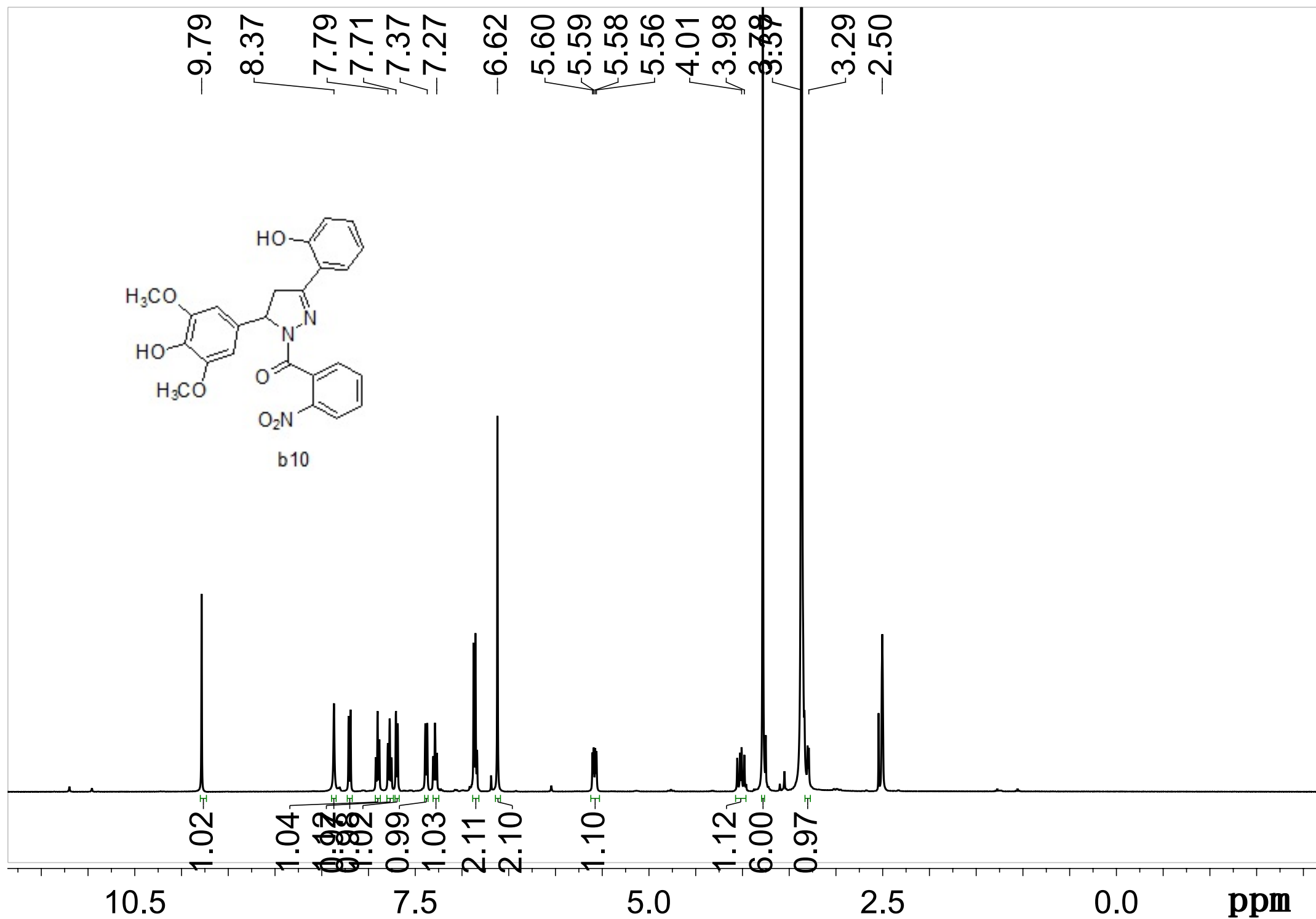
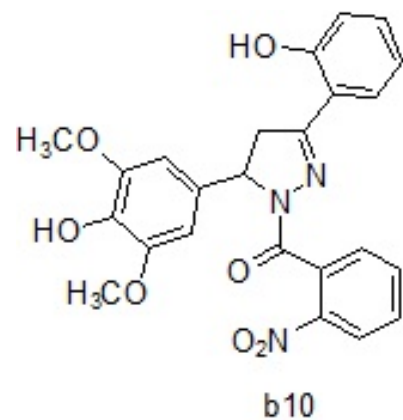


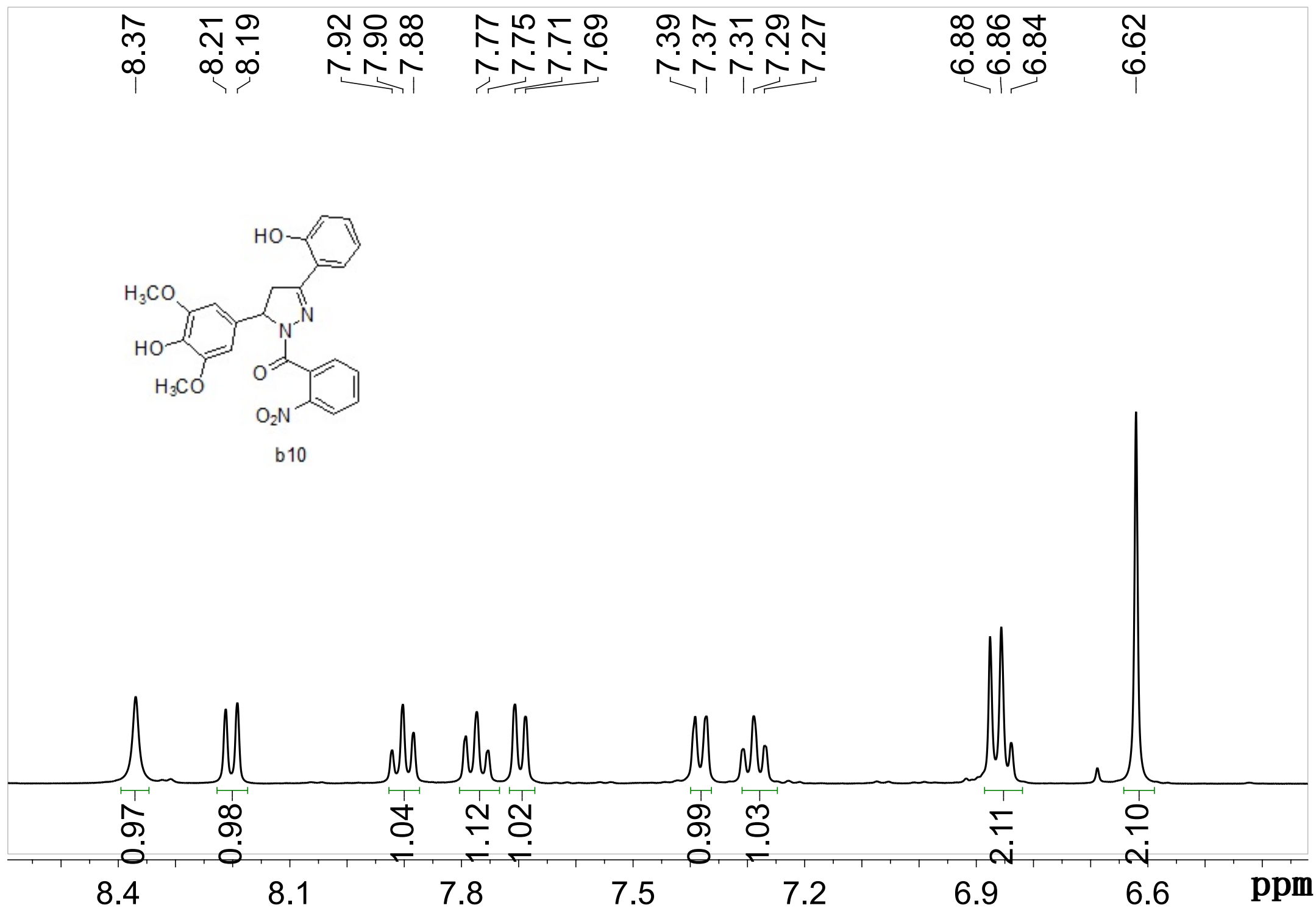
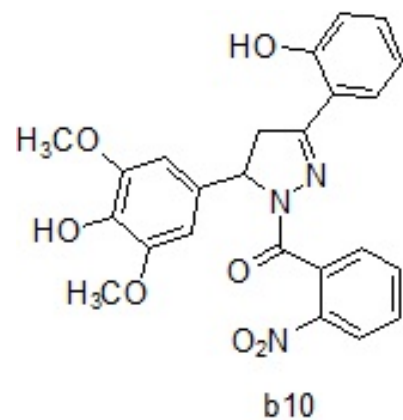


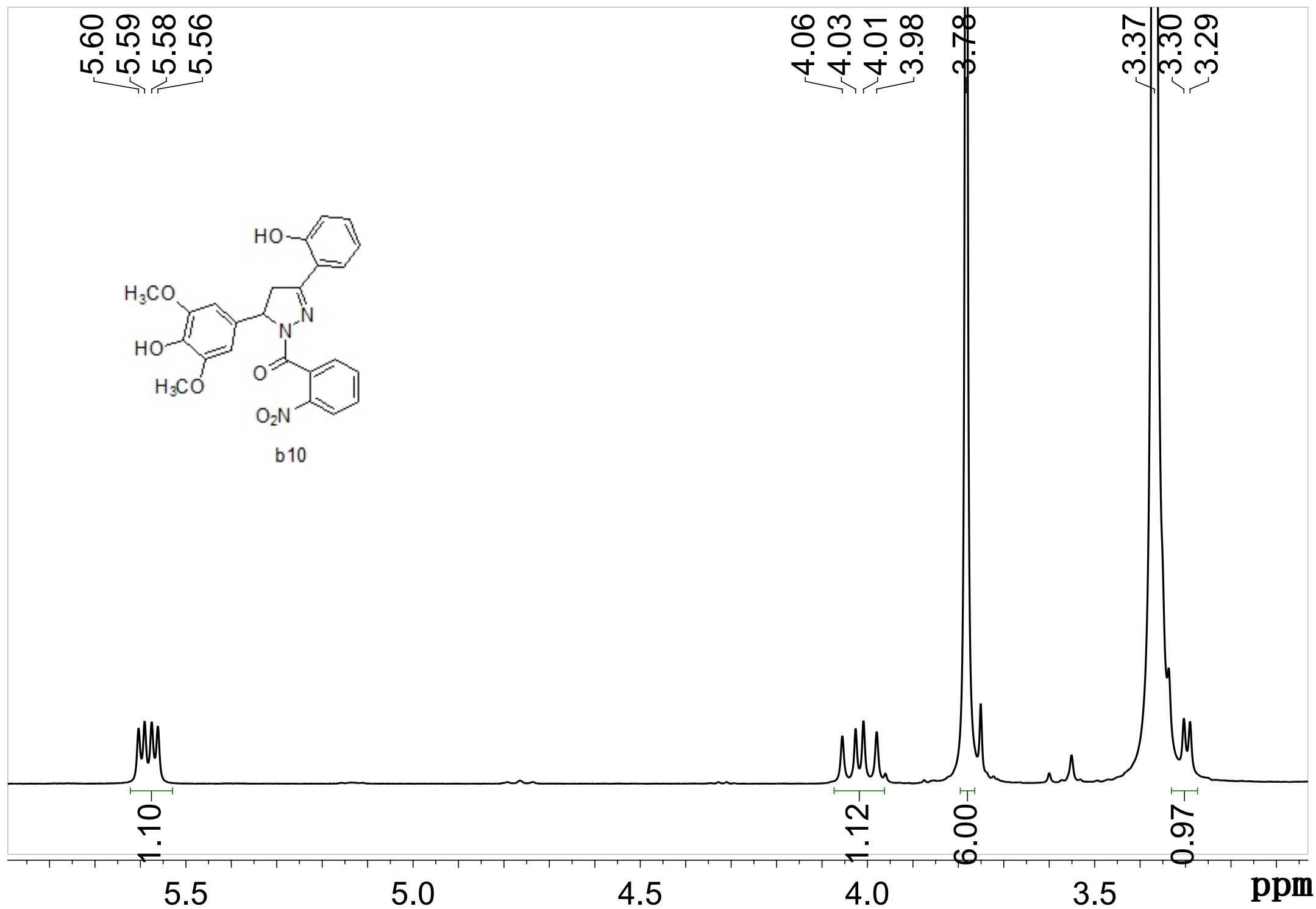
b9

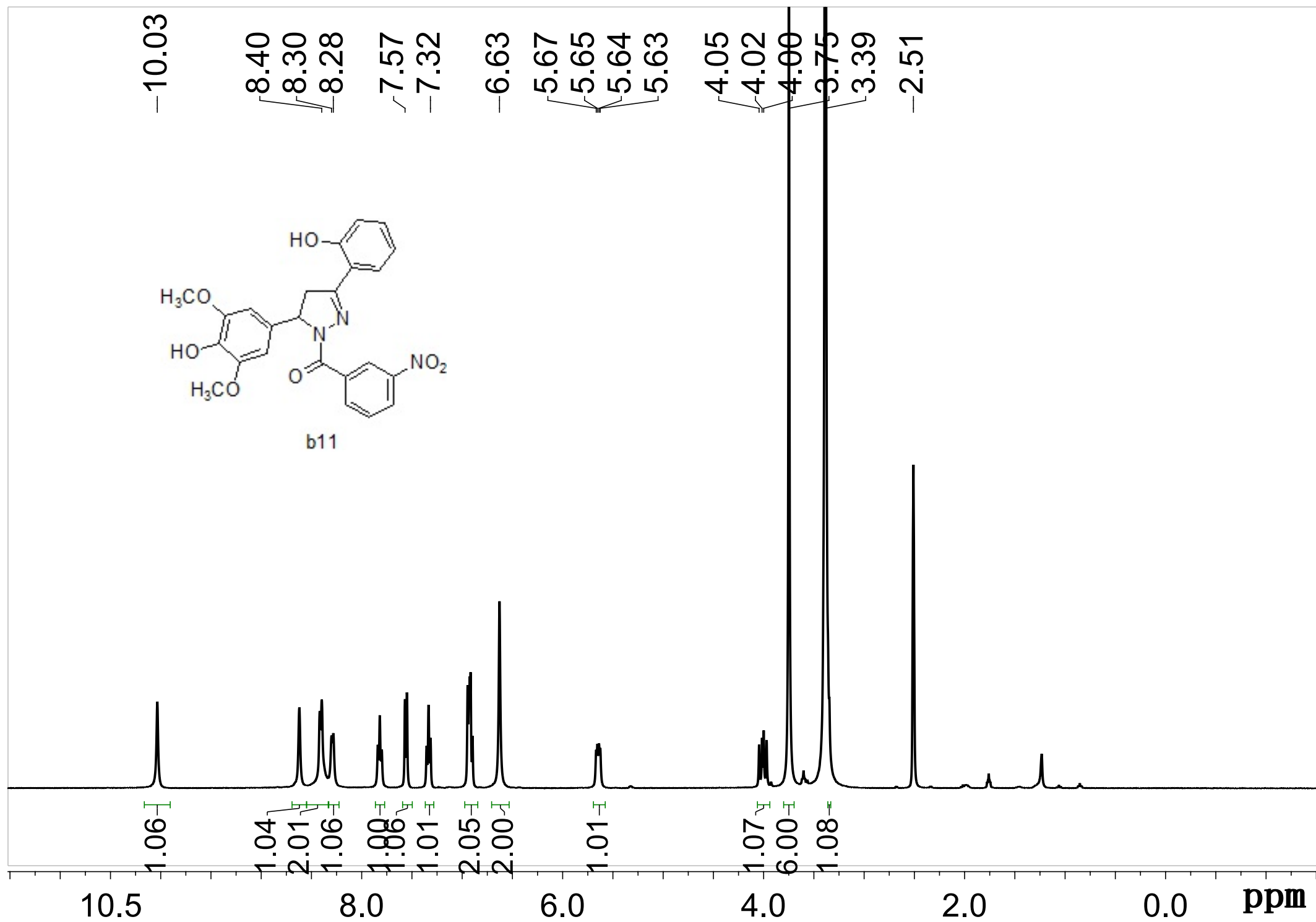


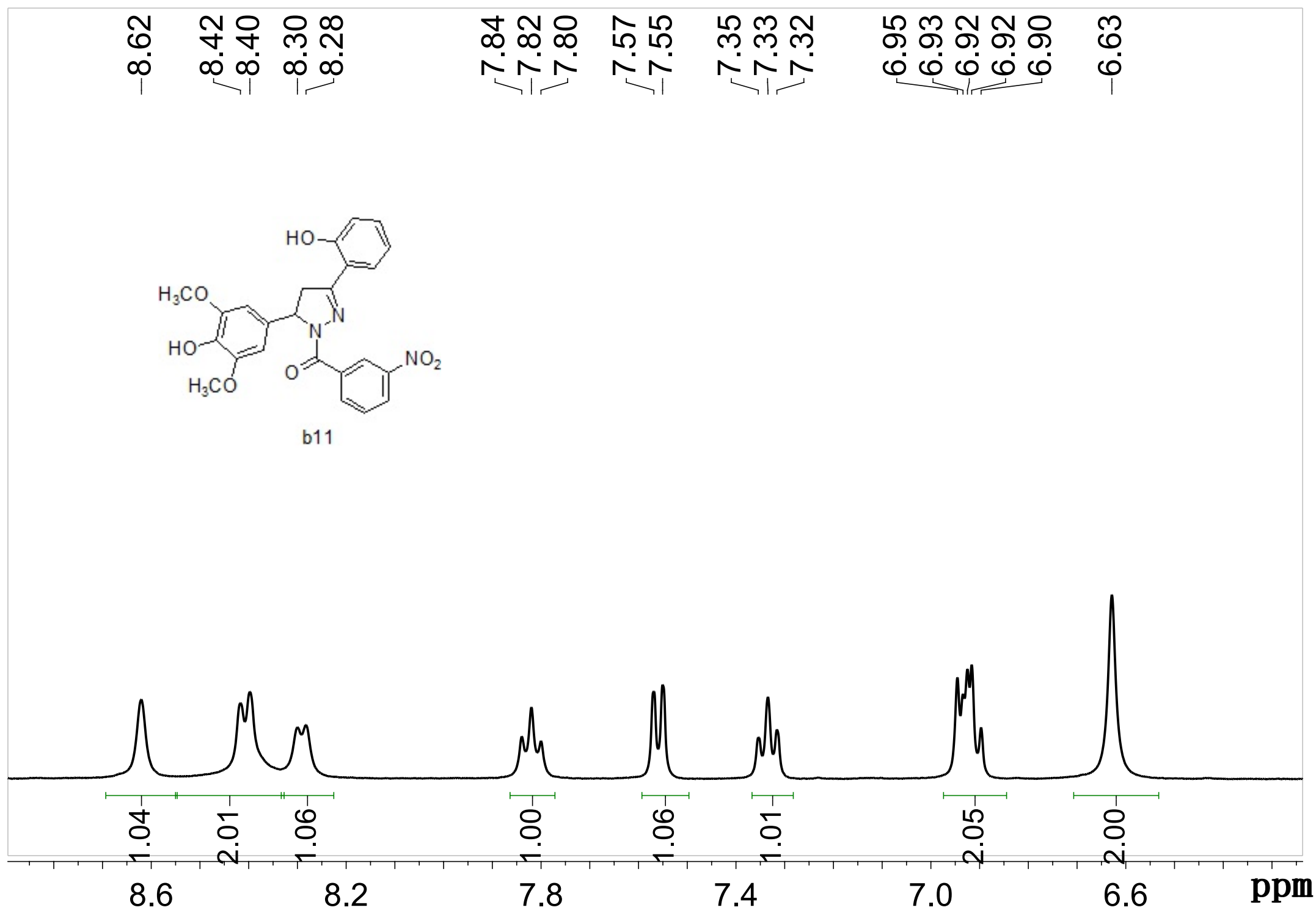


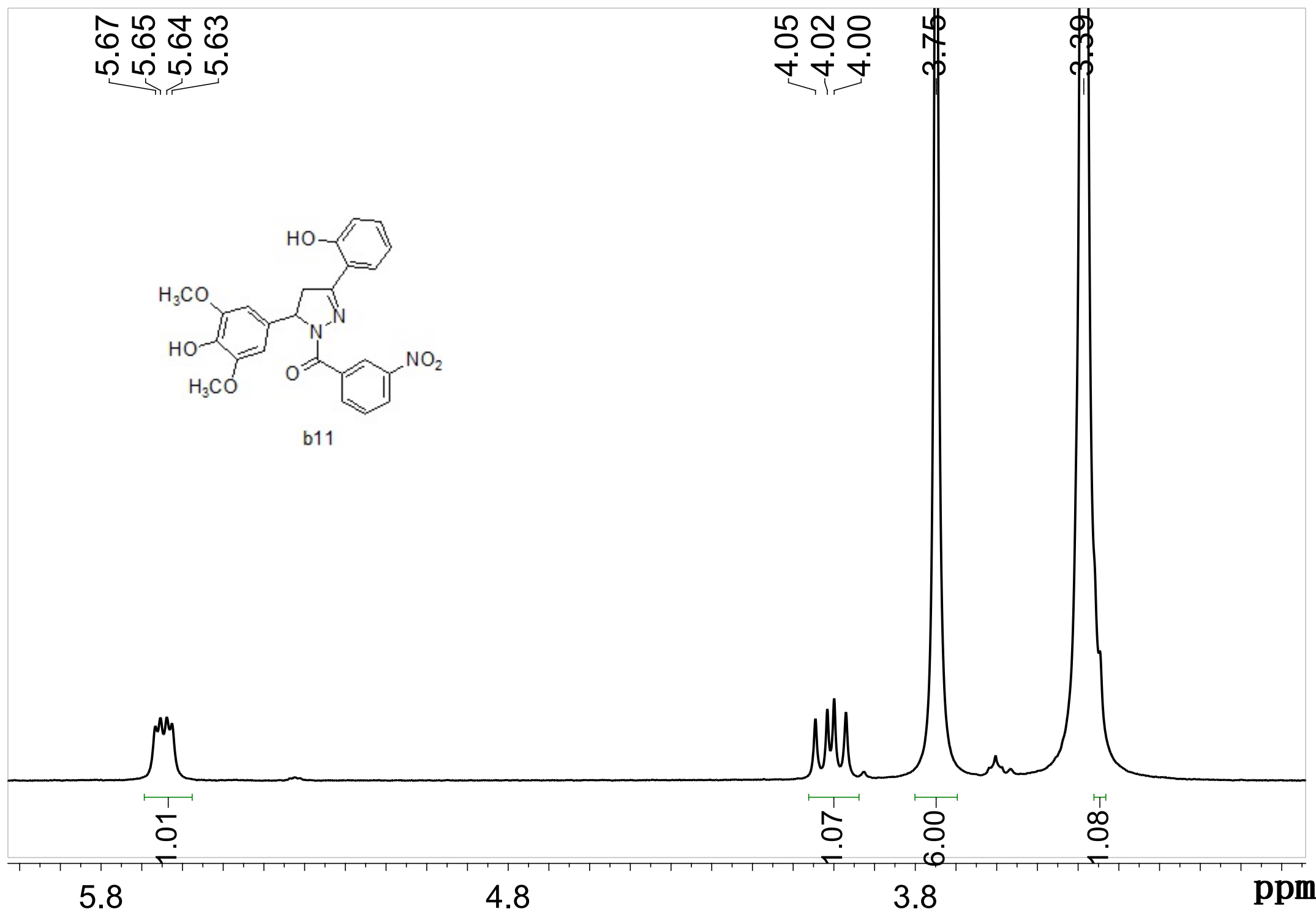


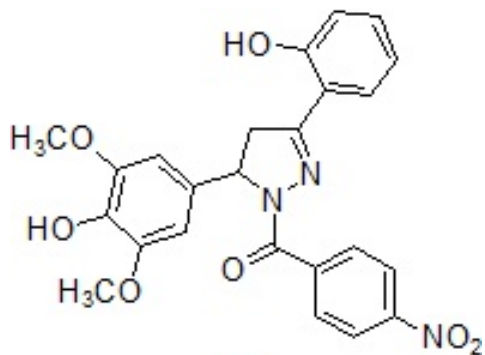












b12

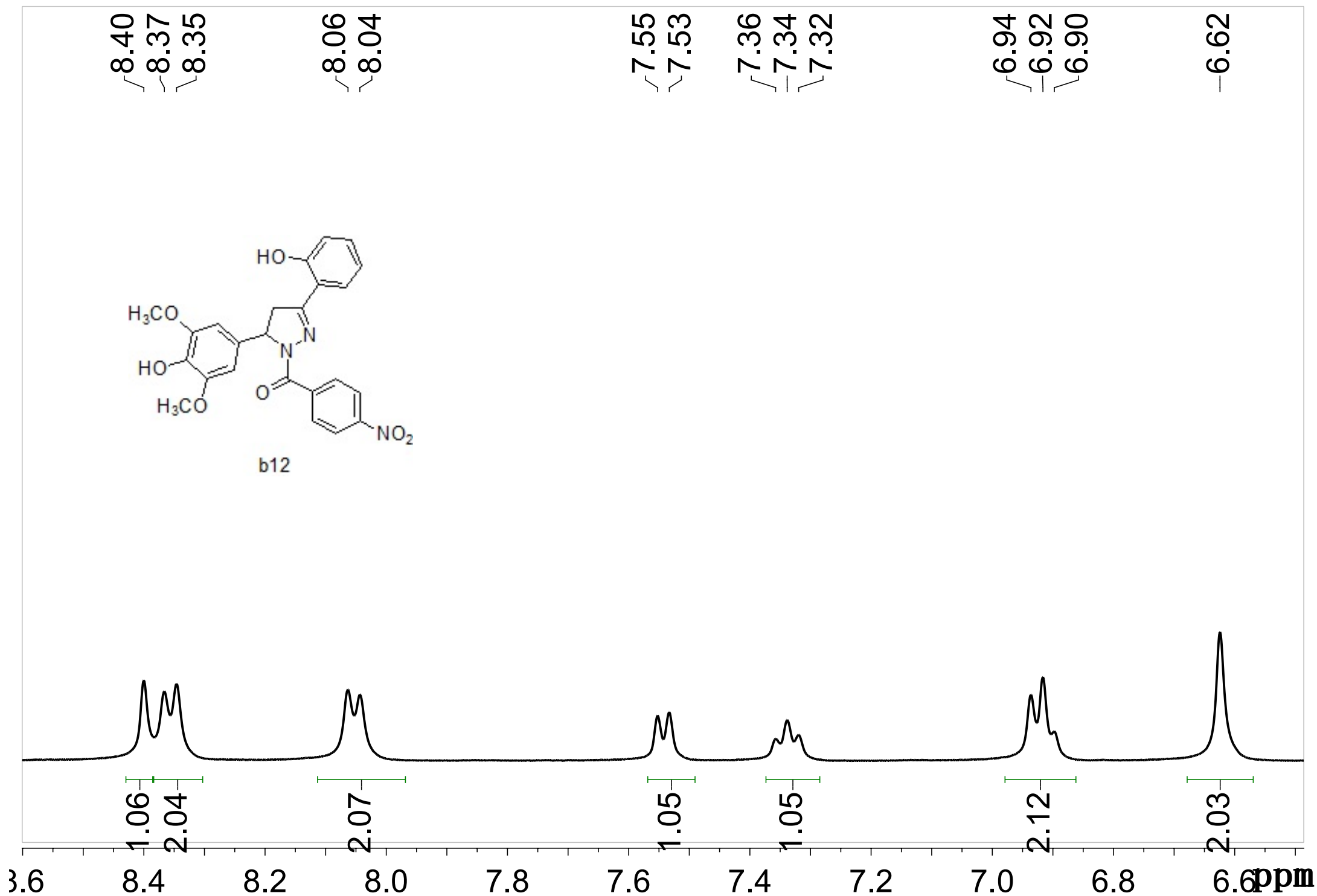
Chemical shift values (ppm) for peaks in the spectrum:

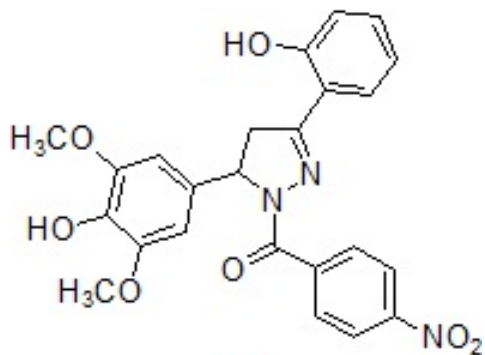
- 10.02
- 8.40
- 8.06
- 8.04
- 7.36
- 6.94
- 6.62
- 5.66
- 5.65
- 4.02
- 4.01
- 3.98
- 3.76
- 3.37
- 2.51

Integration values for peaks in the spectrum:

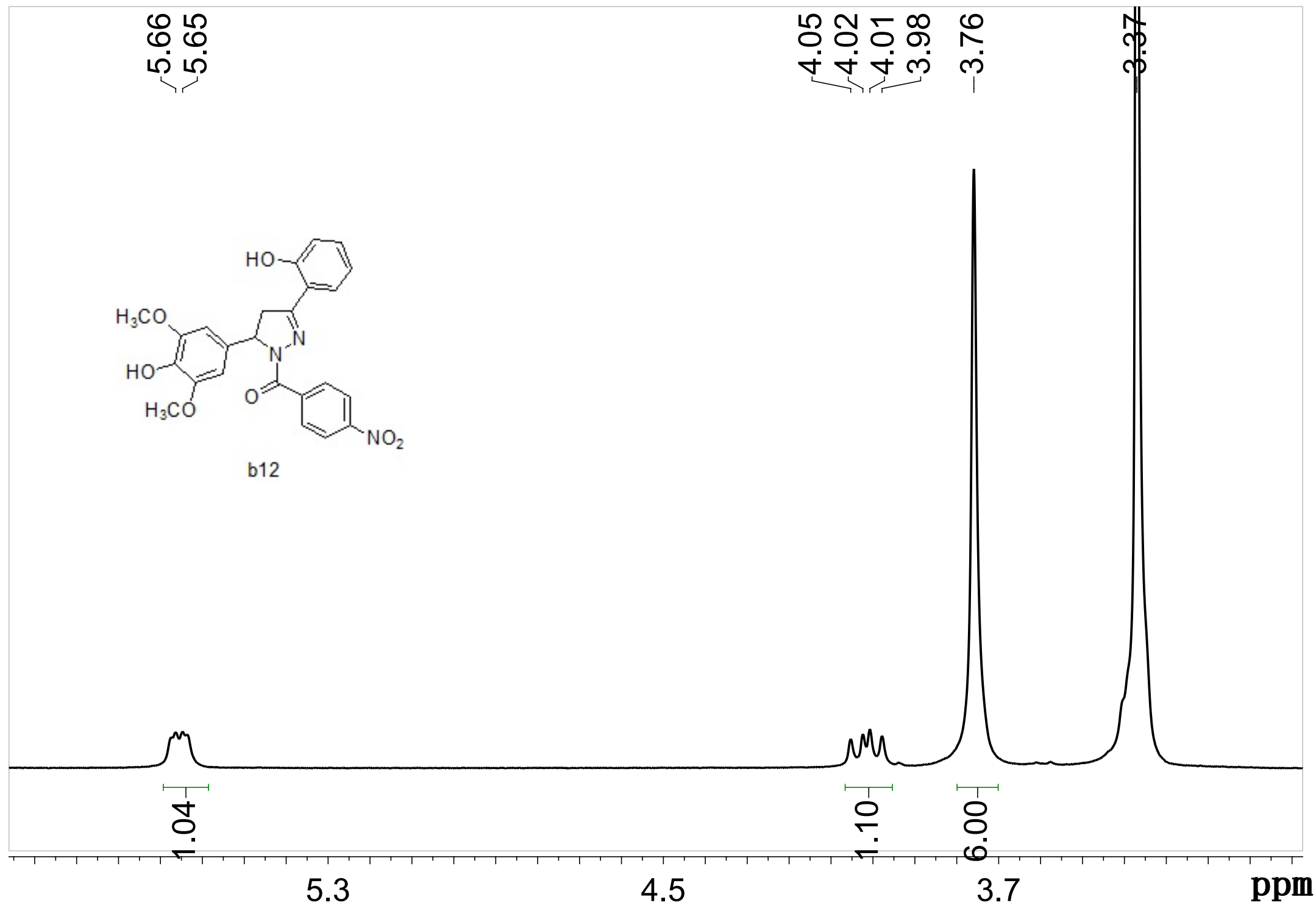
- 1.05
- 1.06
- 2.04
- 2.07
- 1.85
- 1.85
- 2.12
- 2.03
- 1.04
- 1.10
- 6.00

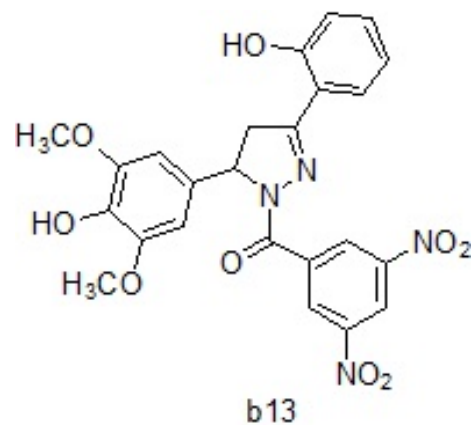
ppm





b12





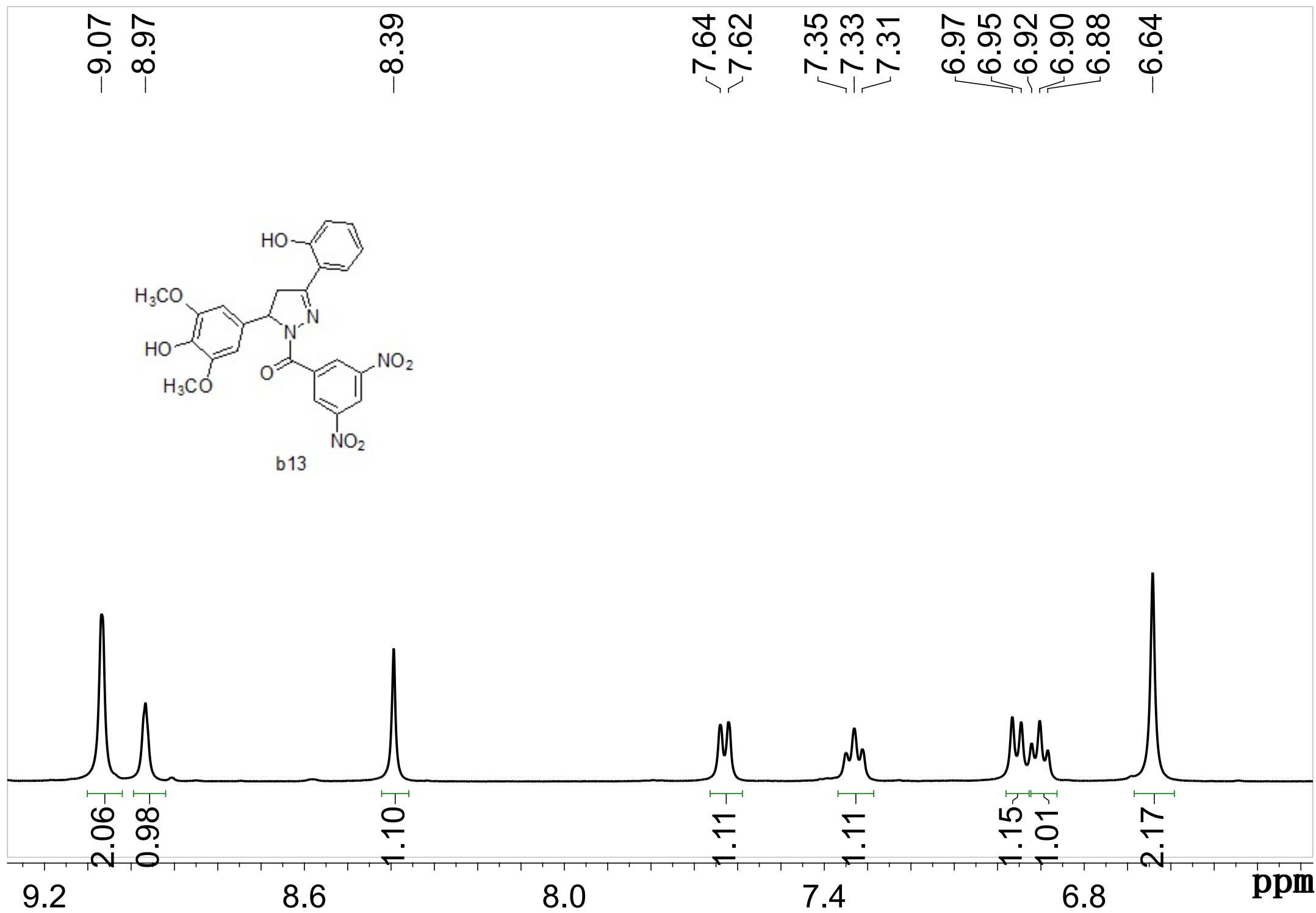
Chemical shifts (ppm):

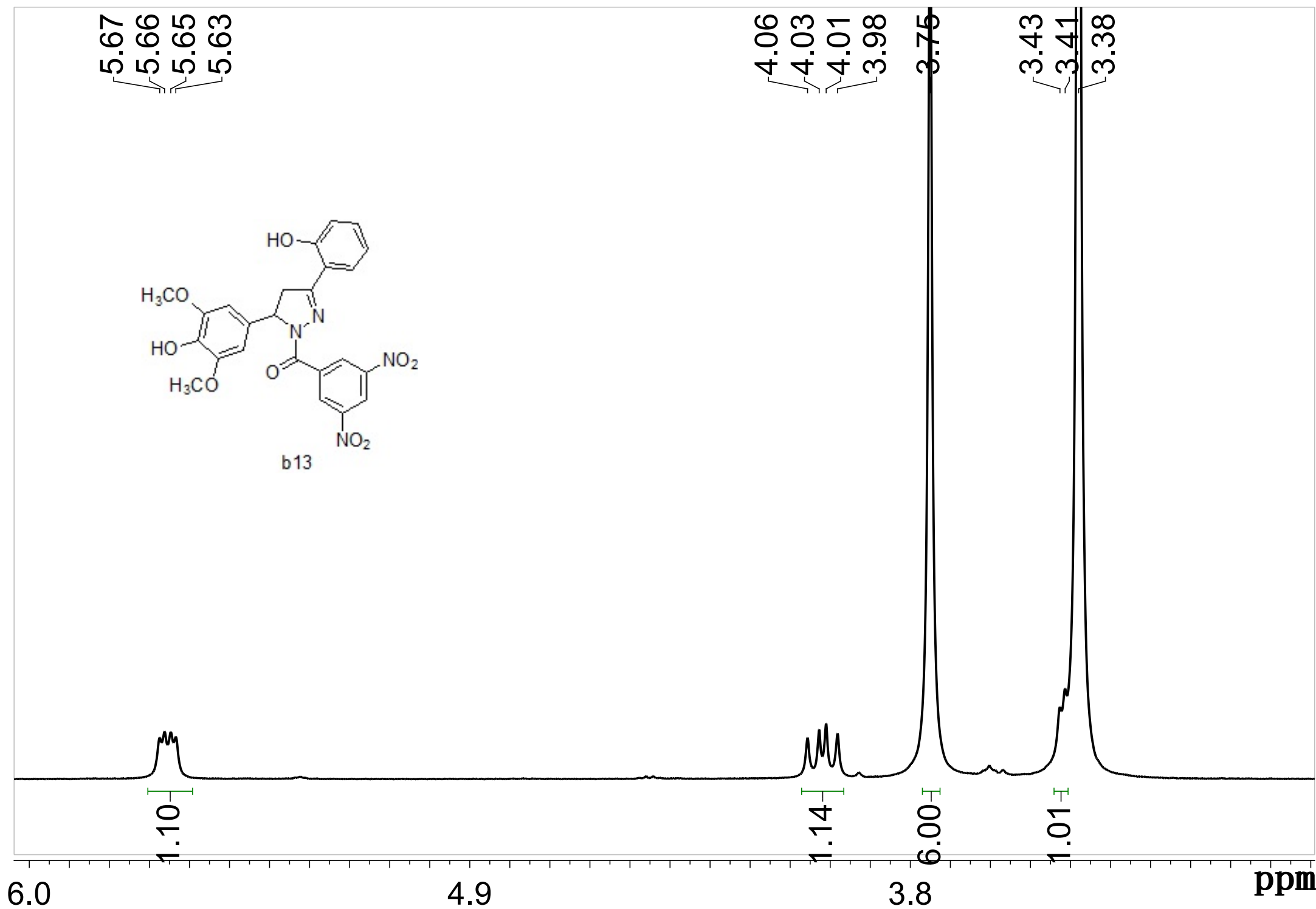
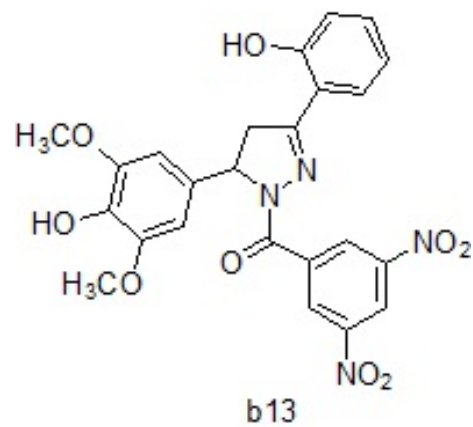
- 10.06
- 9.07
- 8.97
- 8.39
- 7.62
- 6.92
- 6.64
- 5.67
- 5.66
- 5.65
- 4.63
- 4.61
- 3.75
- 3.43
- 3.41
- 3.38
- 2.51

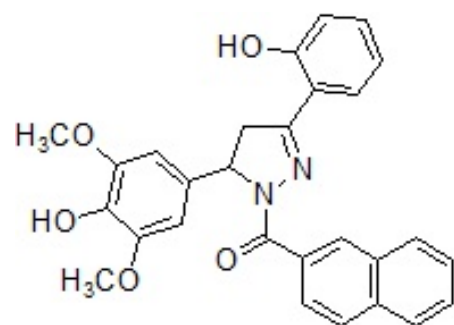
Integration values:

- 1.07
- 2.06
- 0.98
- 1.10
- 1.11
- 1.11
- 1.15
- 1.01
- 2.17
- 1.10
- 1.14
- 6.00
- 1.01

ppm







b14

Chemical shift values (ppm) for compound b14:

- 10.13
- 8.38
- 7.67
- 7.63
- 7.51
- 7.34
- 6.94
- 6.65
- 5.71
- 5.70
- 5.68
- 5.67
- 3.75
- 3.41
- 3.39
- 3.36
- 3.34
- 2.50

Integration values for compound b14:

- 1.10
- 1.02
- 1.02
- 2.10
- 0.86
- 2.18
- 1.09
- 1.13
- 2.12
- 1.93
- 1.04
- 1.08
- 6.00
- 1.04

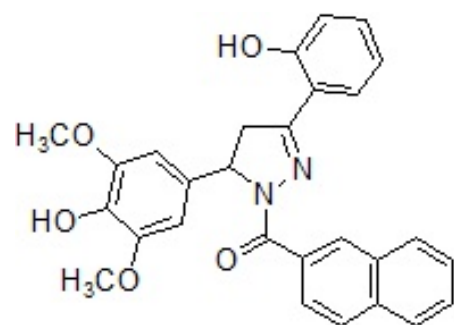
9.0

6.0

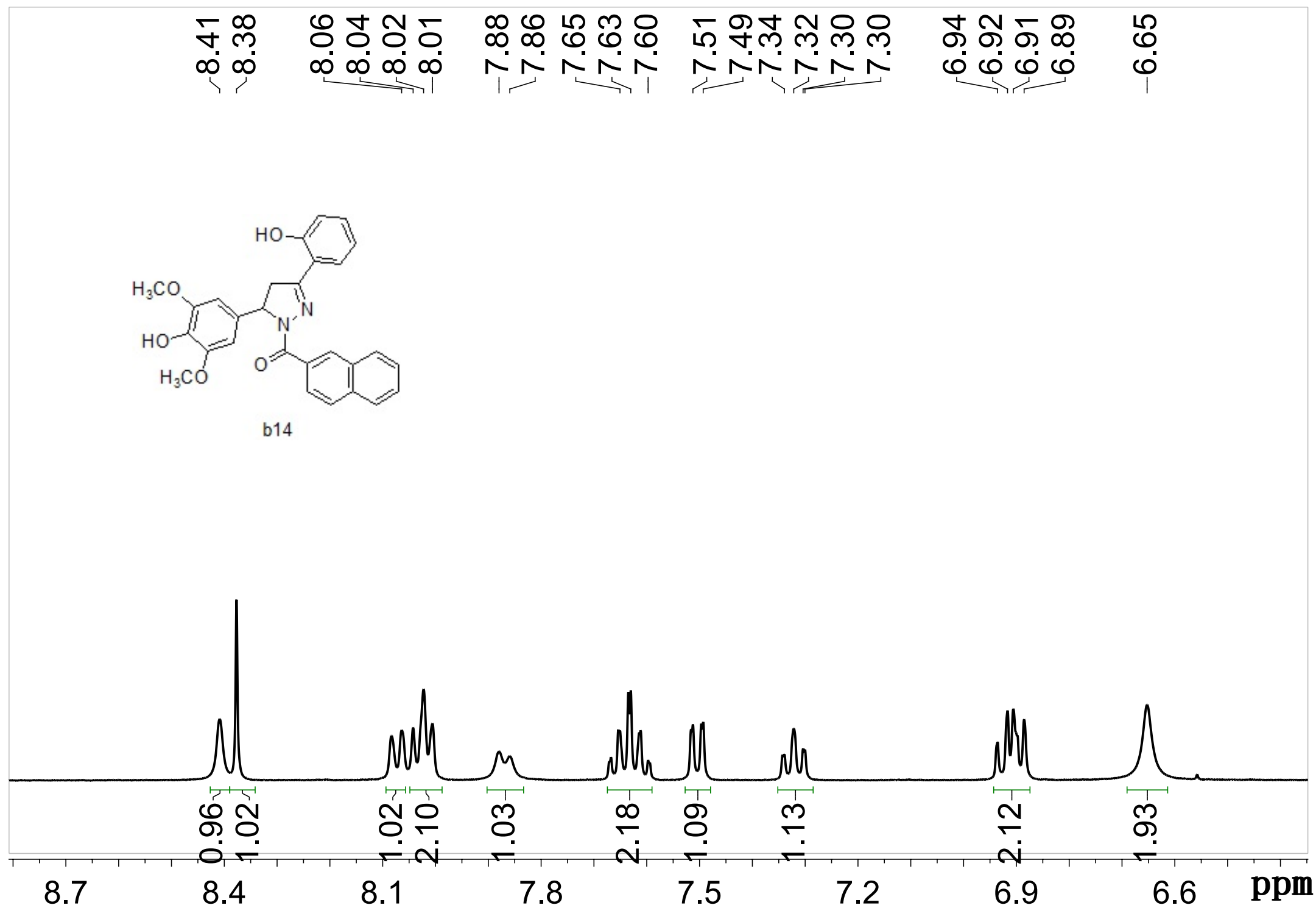
3.0

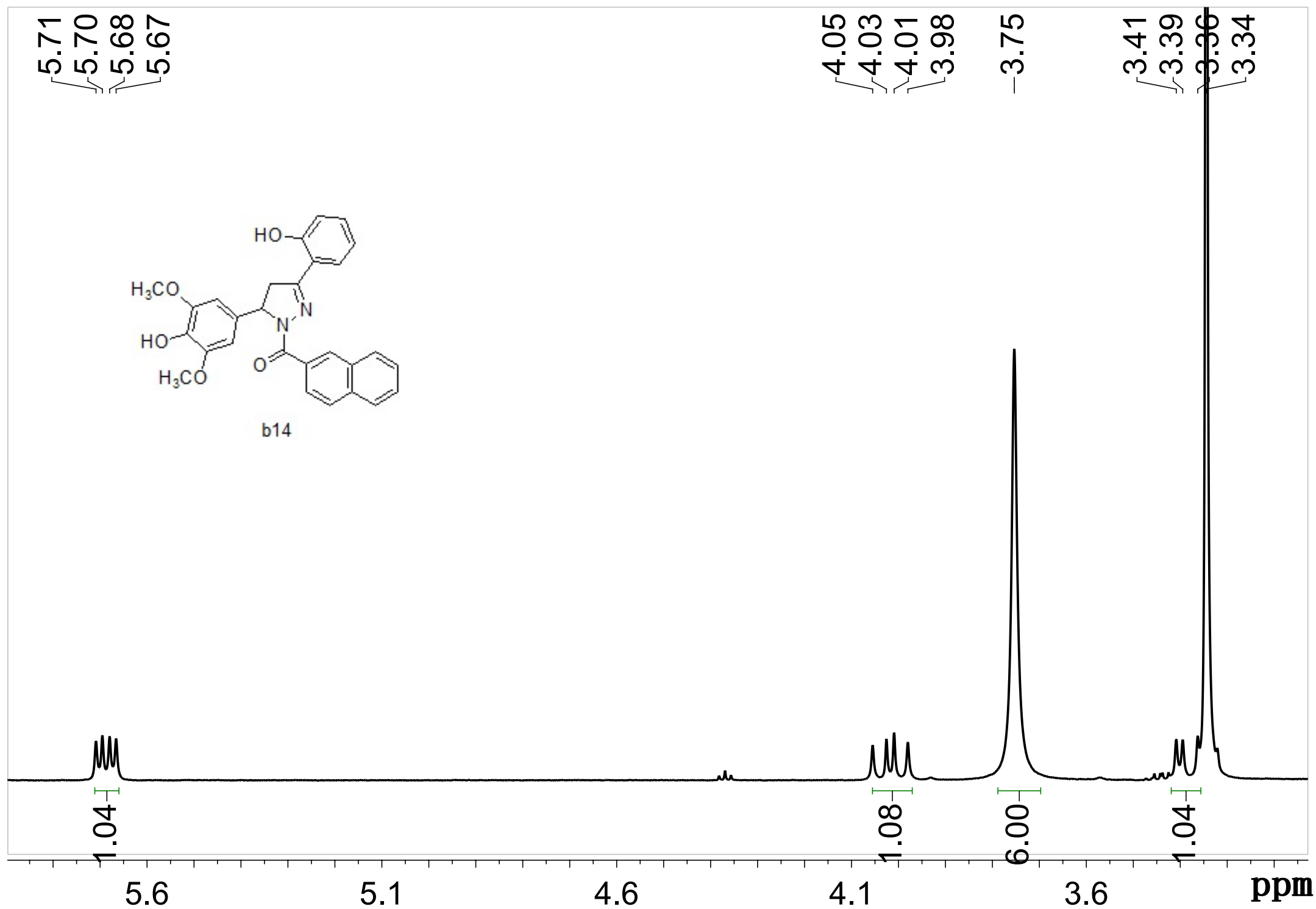
0.0

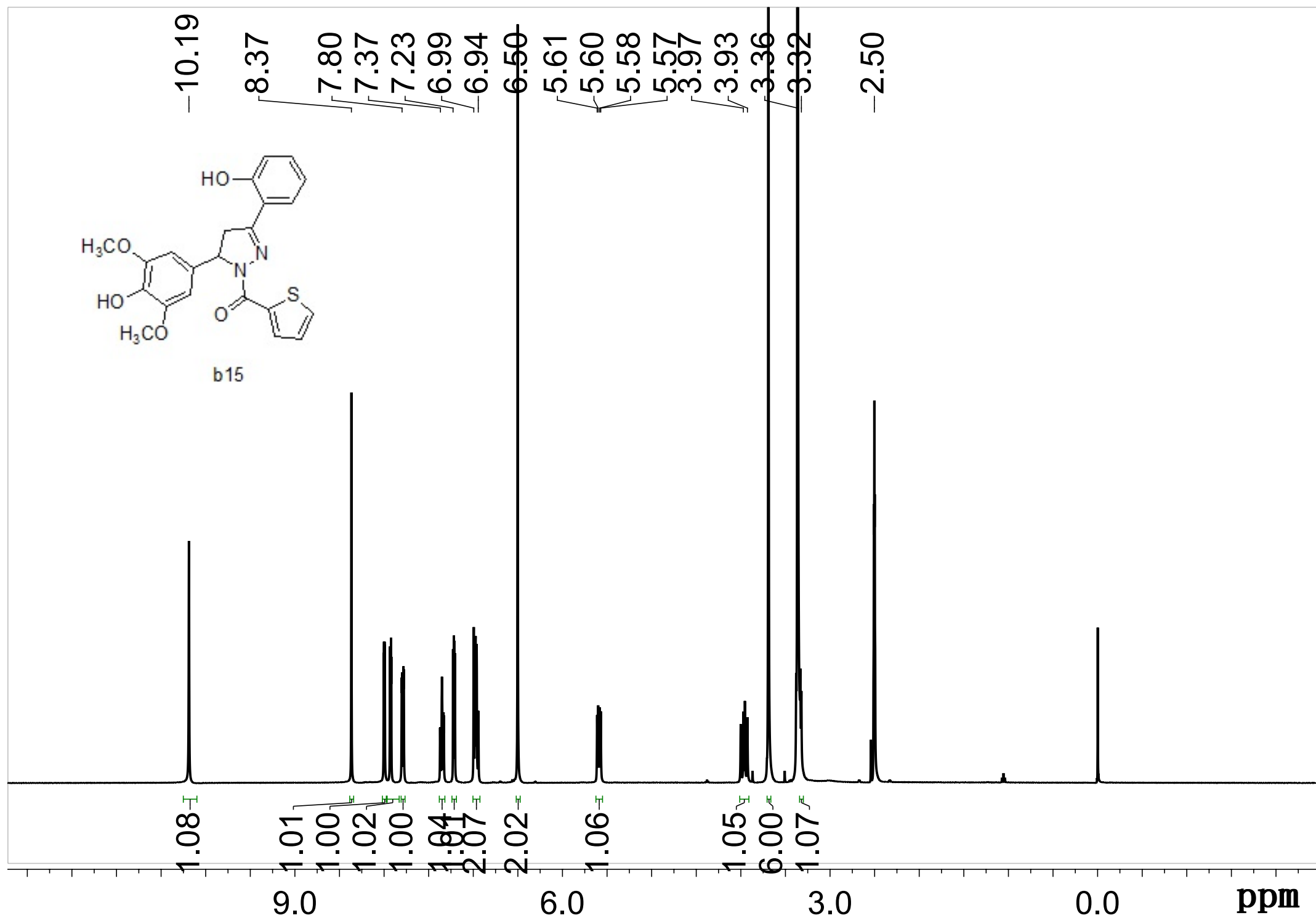
ppm

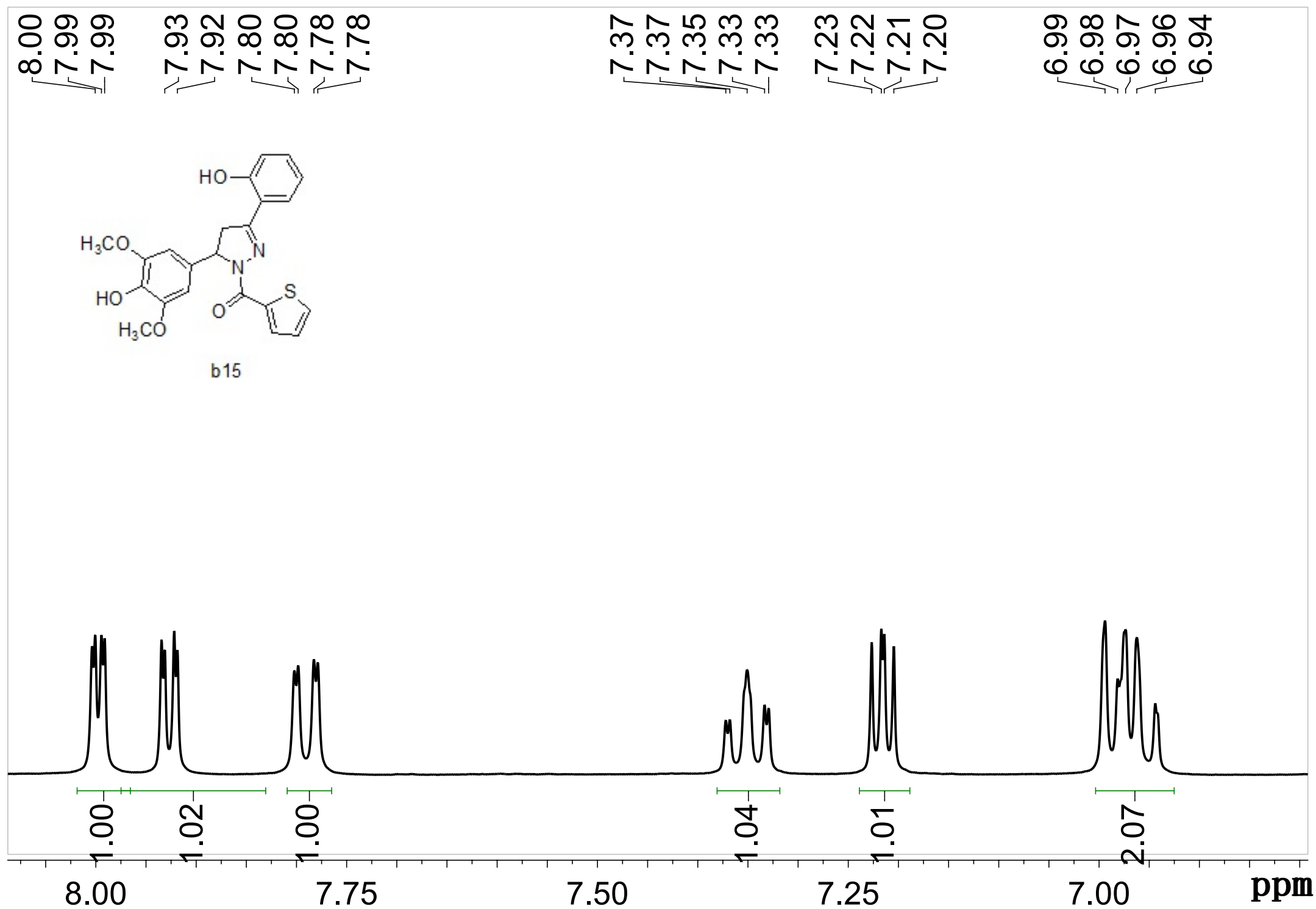


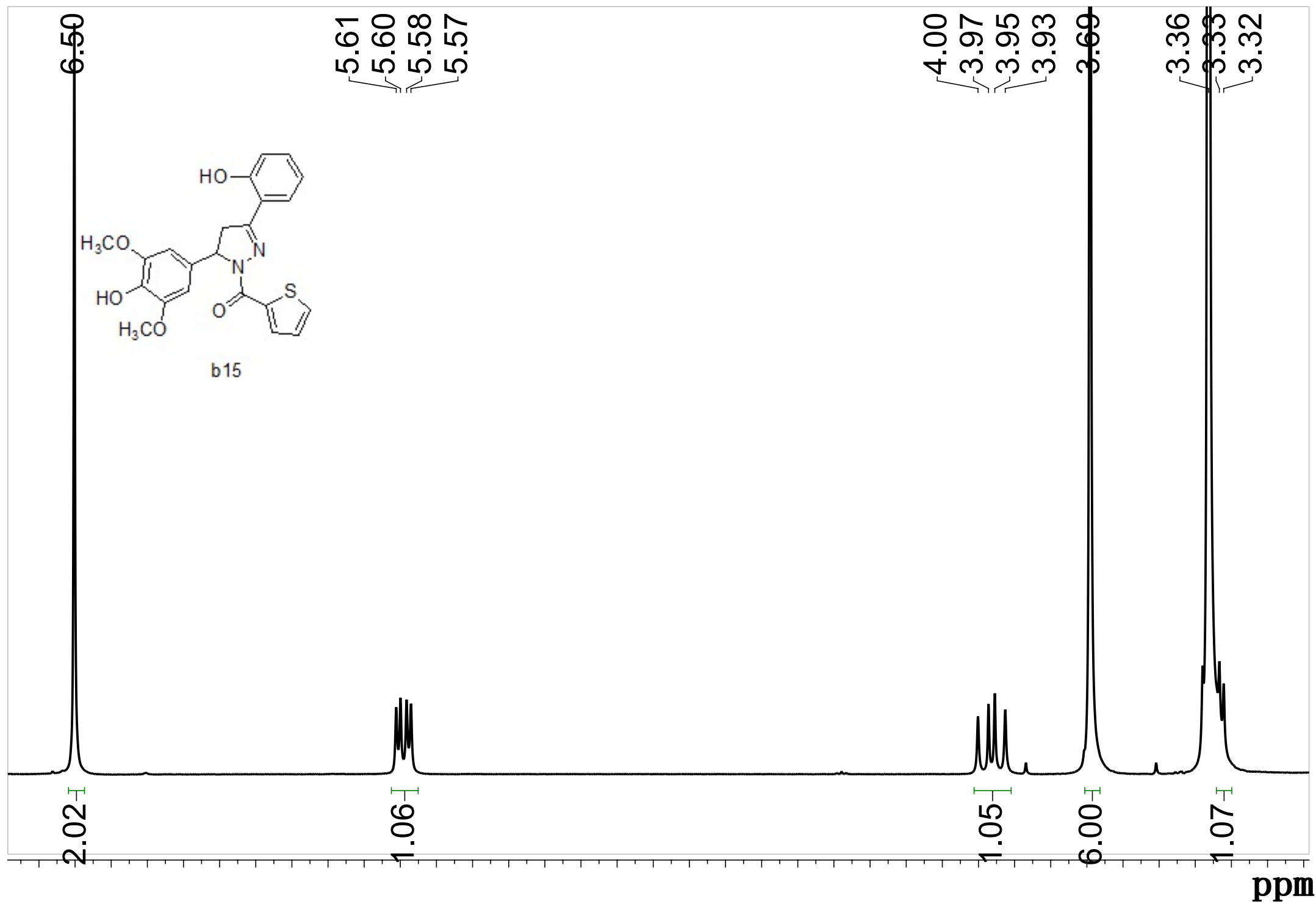
b14

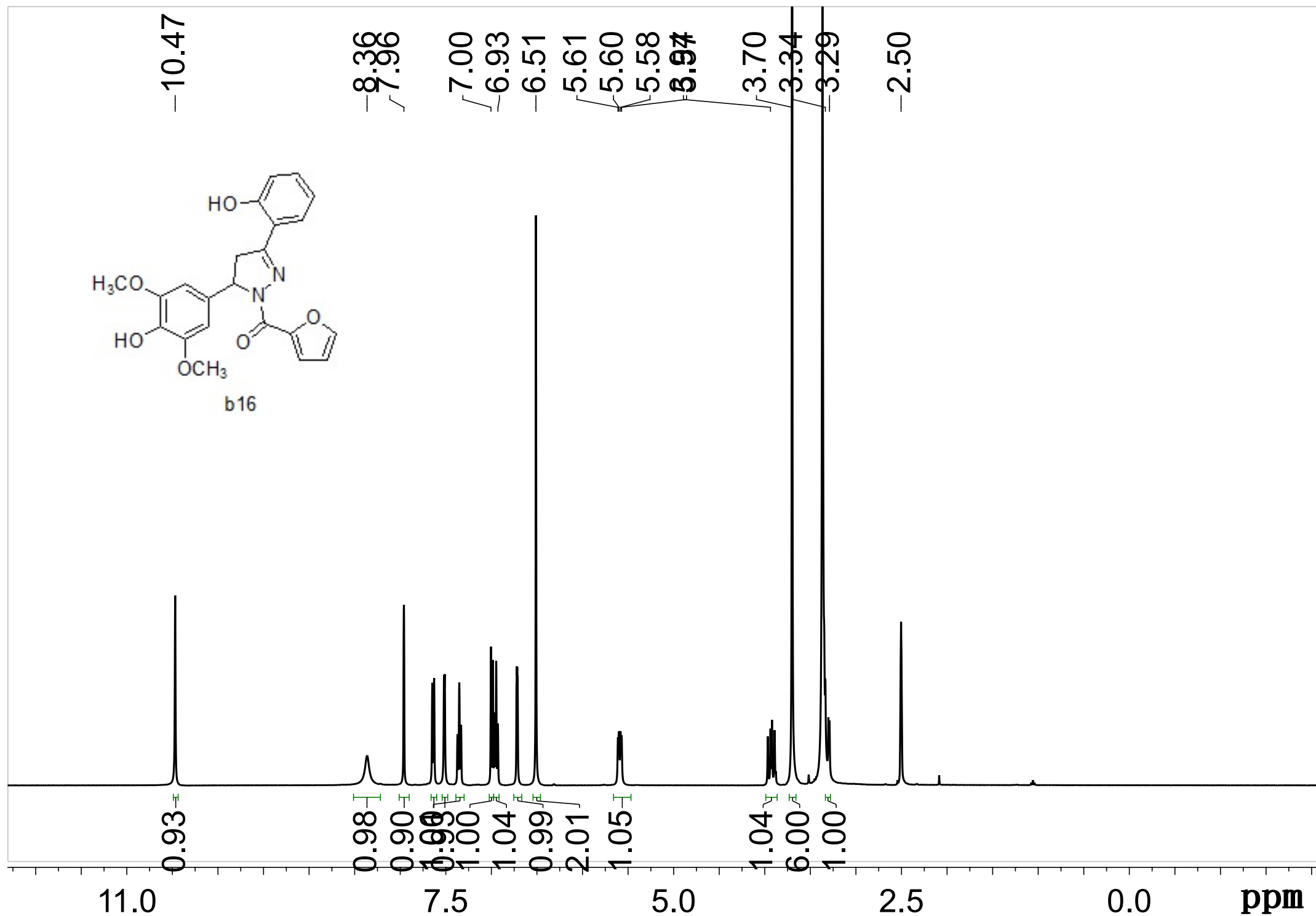
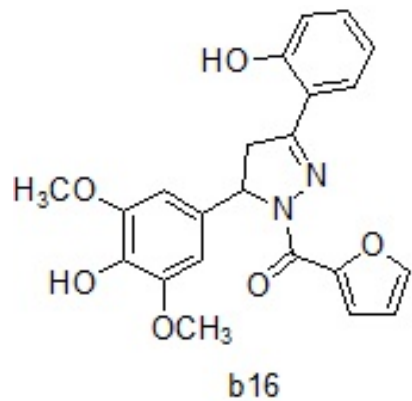


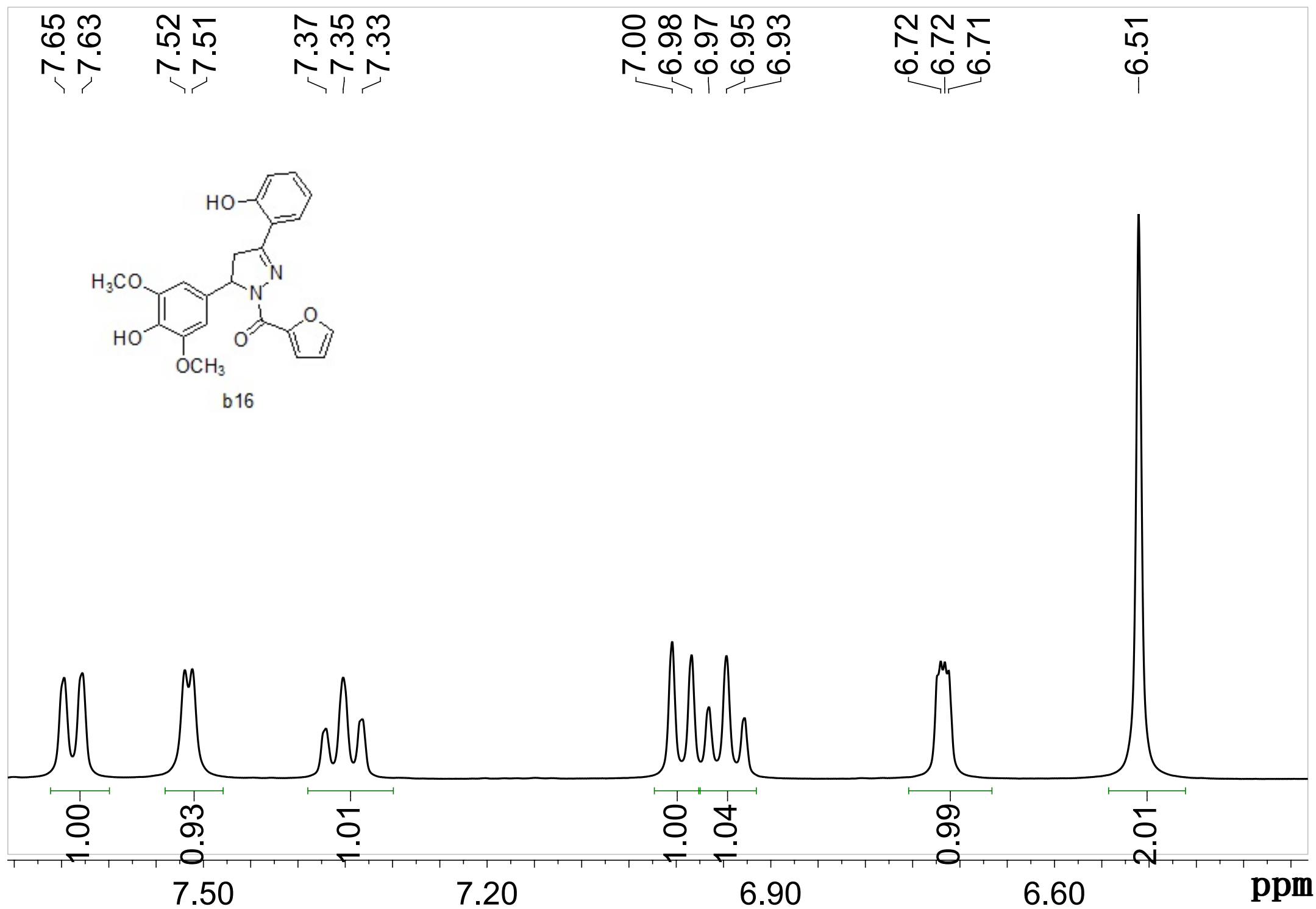


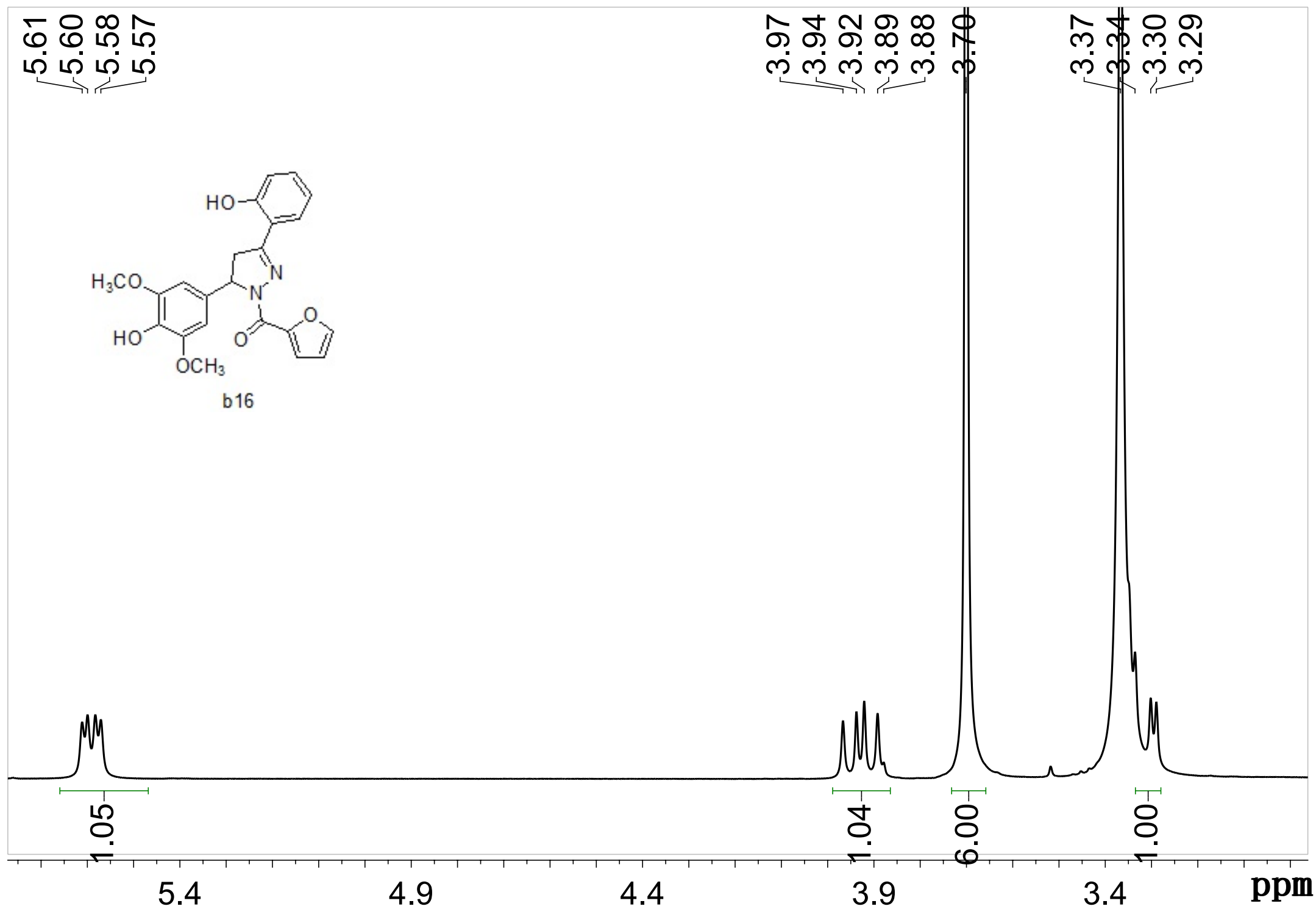


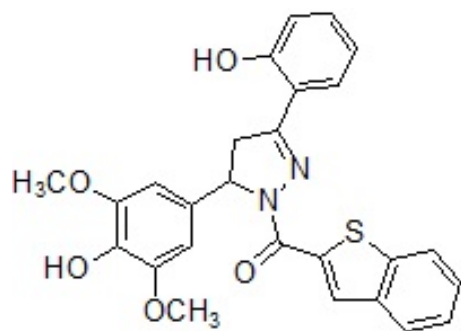




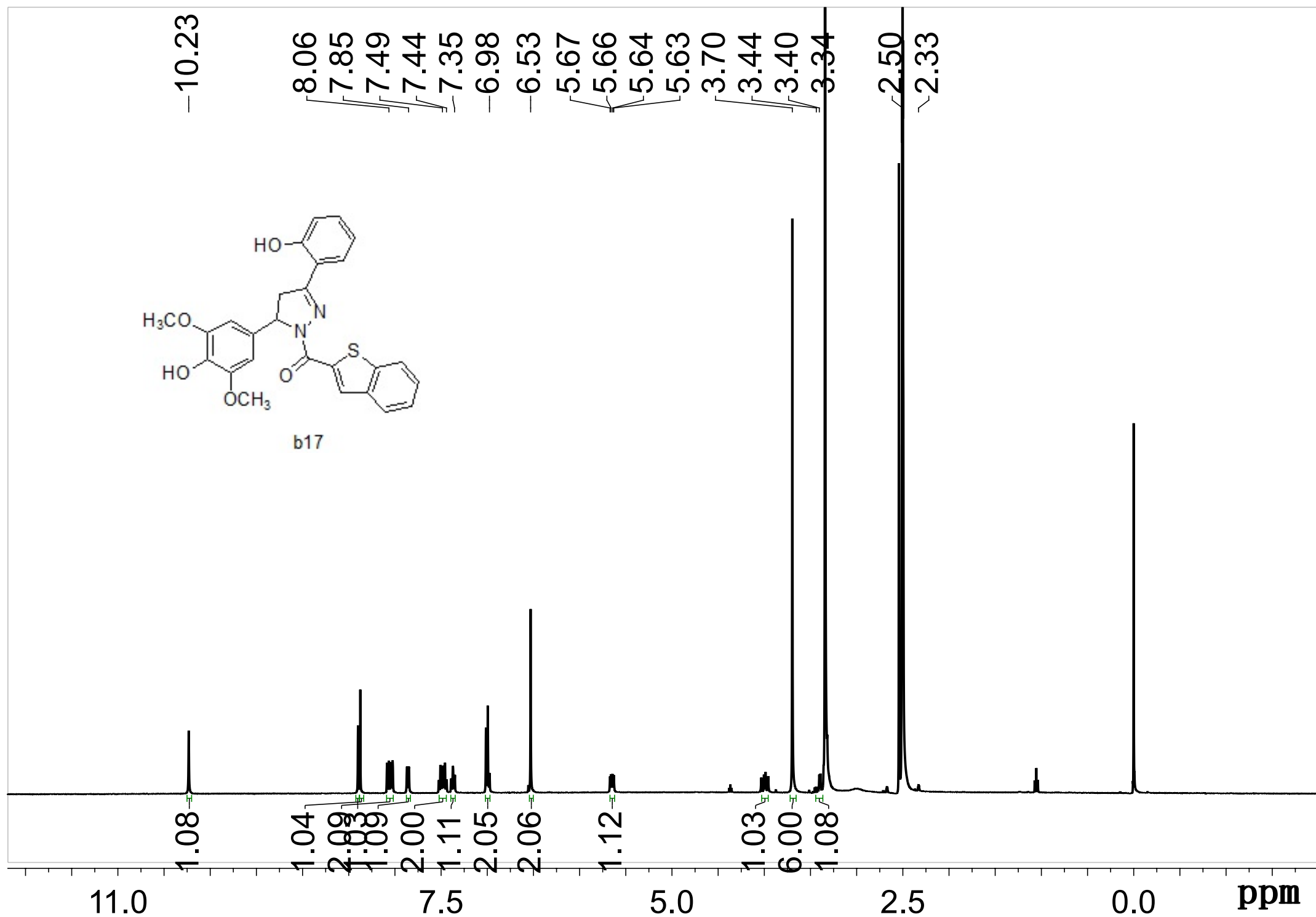


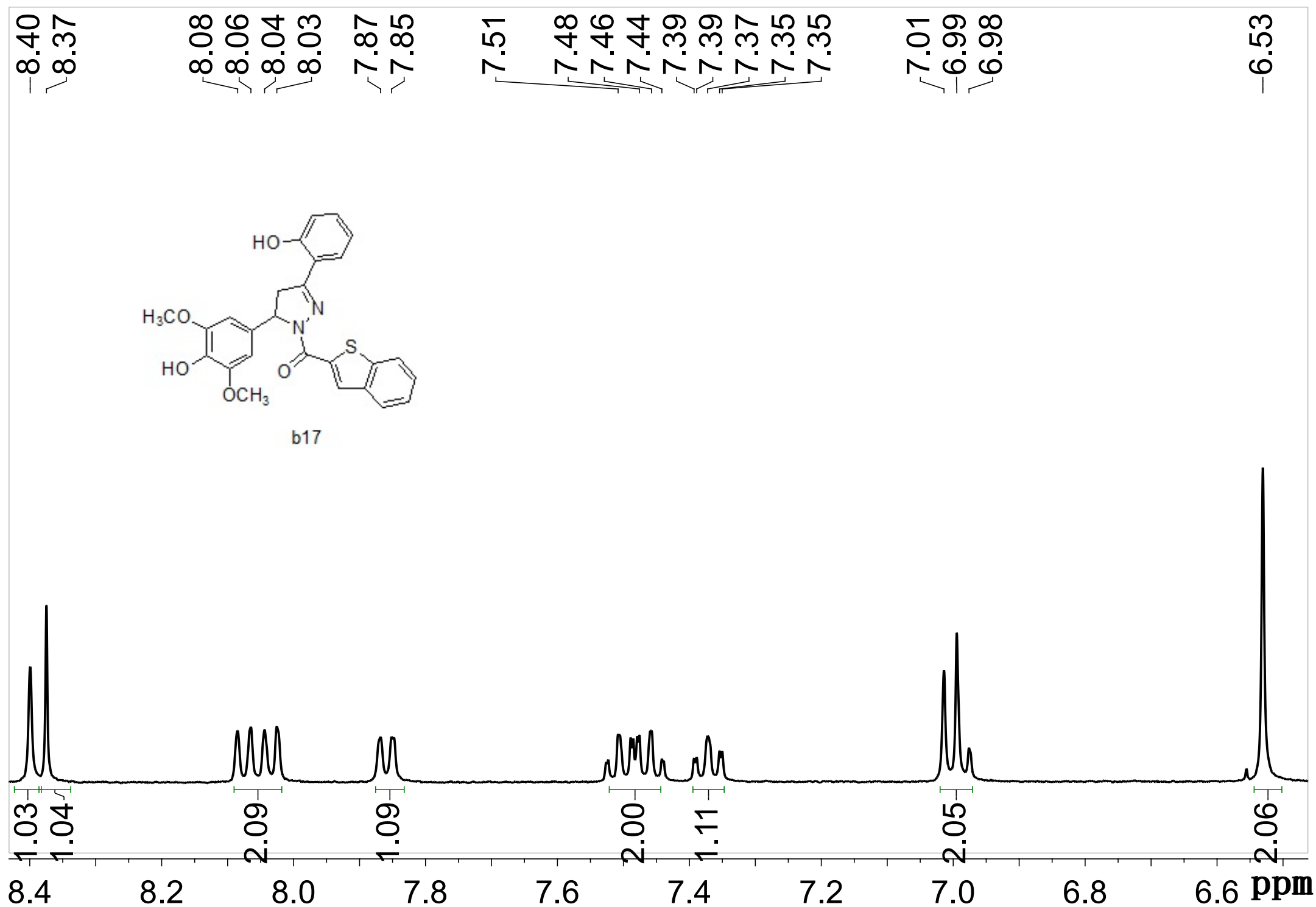


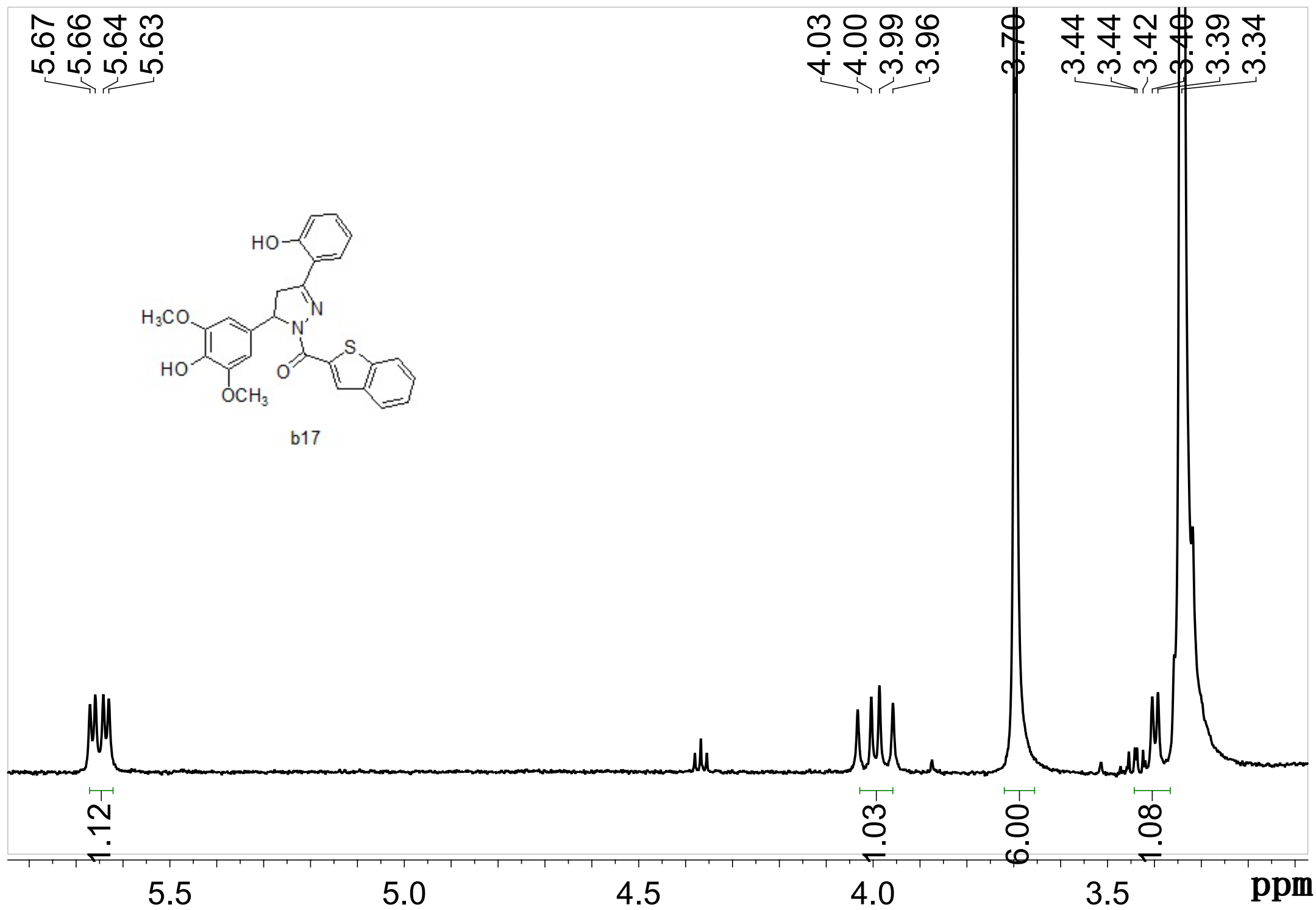


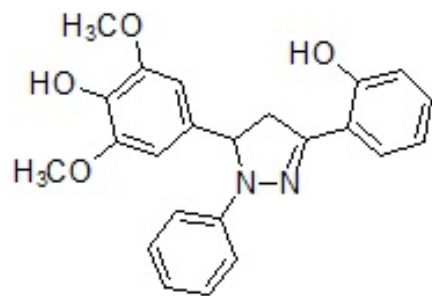


b17

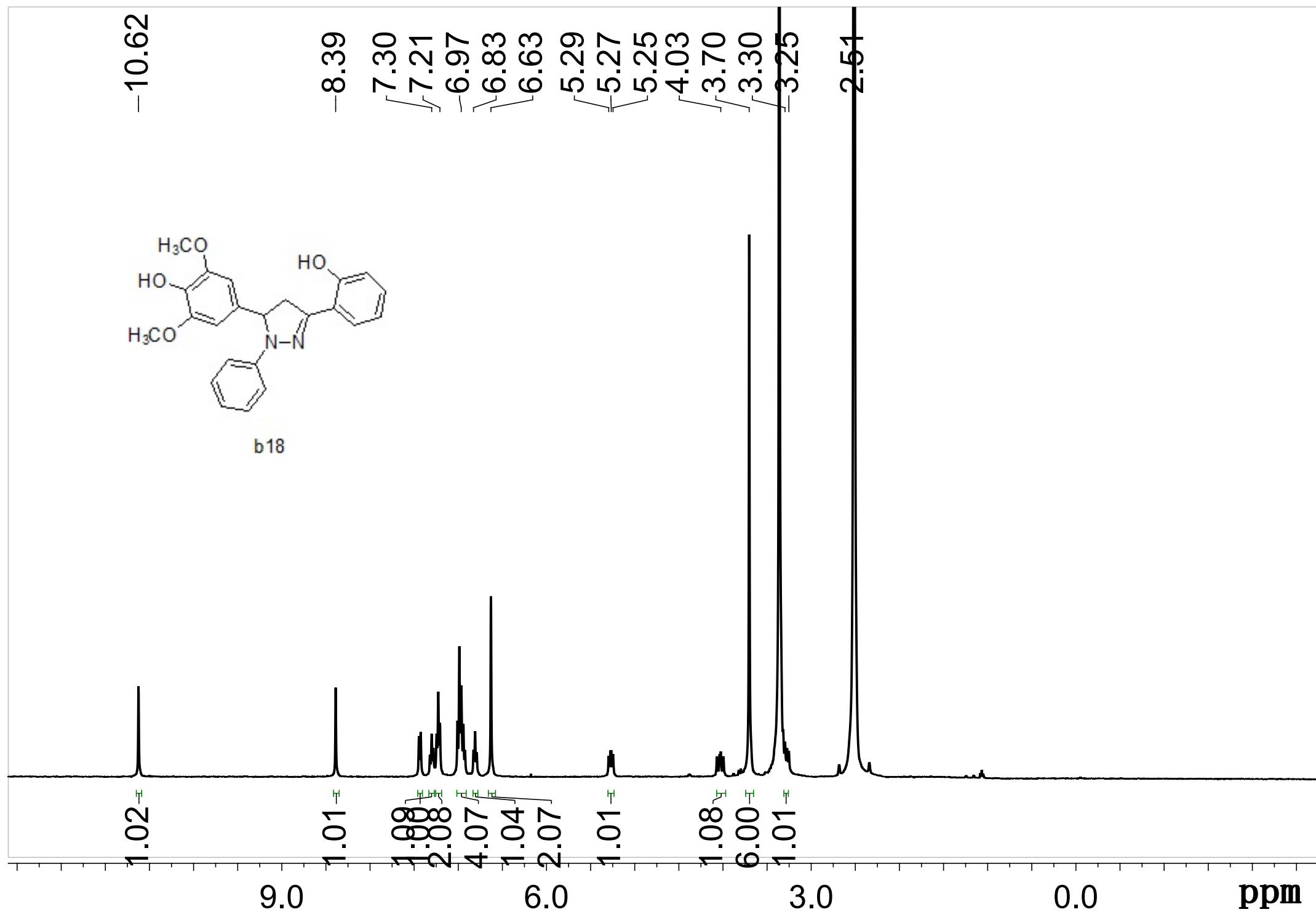


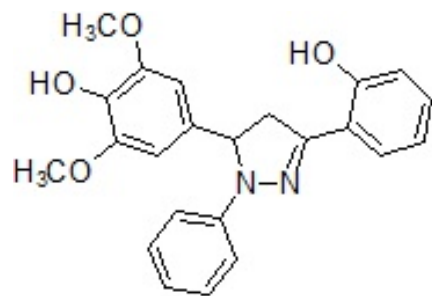






b18





b18

