

Supporting Information

Synthesis of (2*H*)-Indazoles through Rh(III)-Catalyzed Annulation Reaction of Azobenzenes with Sulfoxonium Ylides

Hyunjung Oh,[†] Sangil Han,[†] Ashok Kumar Pandey, Sang Hoon Han, Neeraj Kumar Mishra, Saegun Kim, Rina Chun, Hyung Sik Kim, Jihye Park,^{*} and In Su Kim^{*}

School of Pharmacy, Sungkyunkwan University, Suwon 440-746, Republic of Korea

^{*} Corresponding author. Tel.: +82-31-290-7788; fax: +82-31-292-8800; e-mail: pjh3548@skku.edu (J.P), insukim@skku.edu (I.S.K)

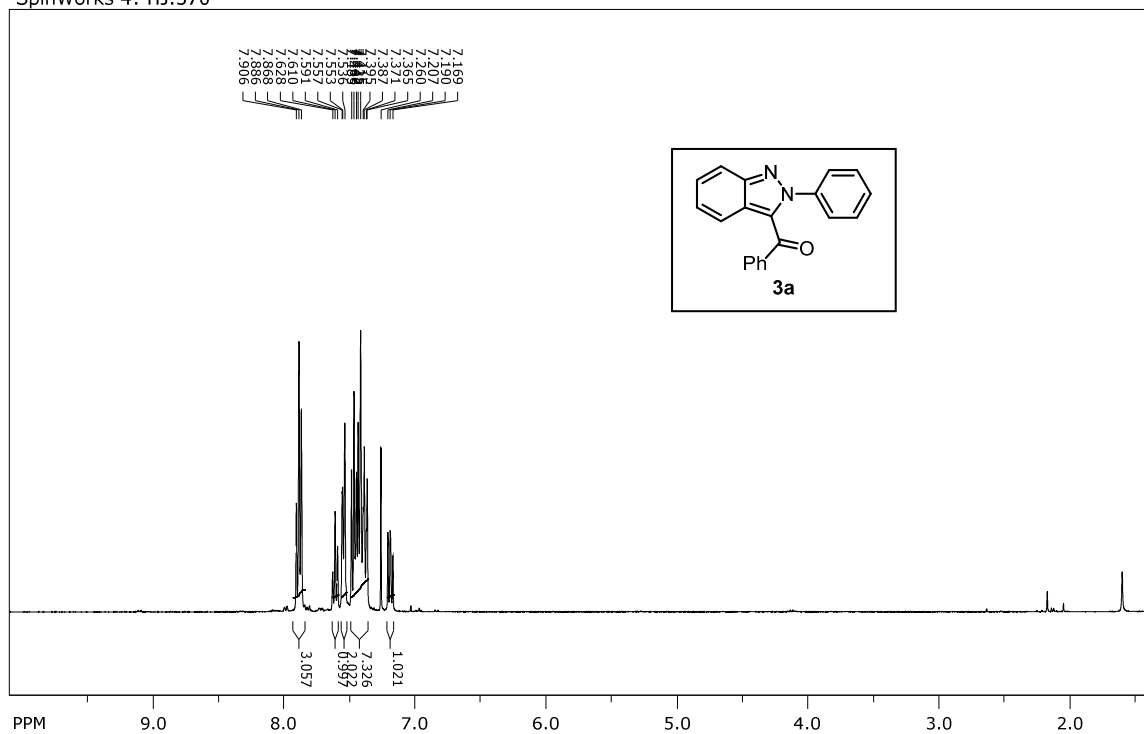
[†] These authors equally contributed.

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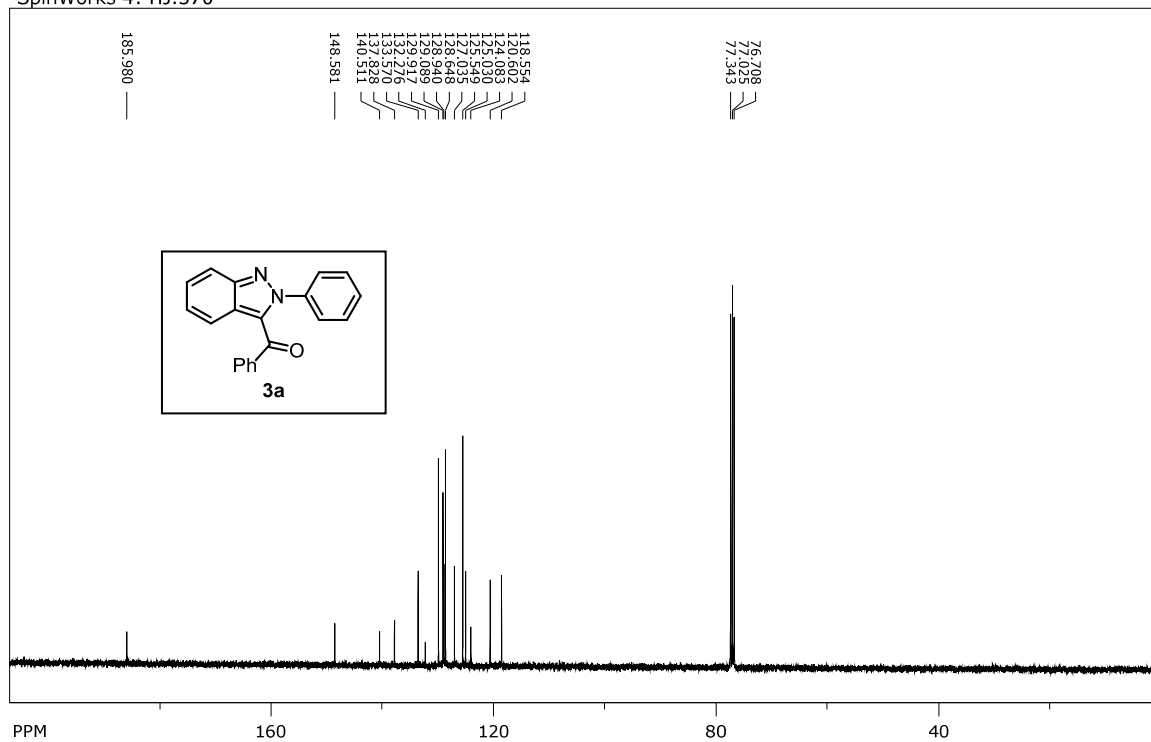
¹H NMR and ¹³C NMR spectra of all compounds ----- S2–S45

¹H and ¹³C NMR spectra of all compounds

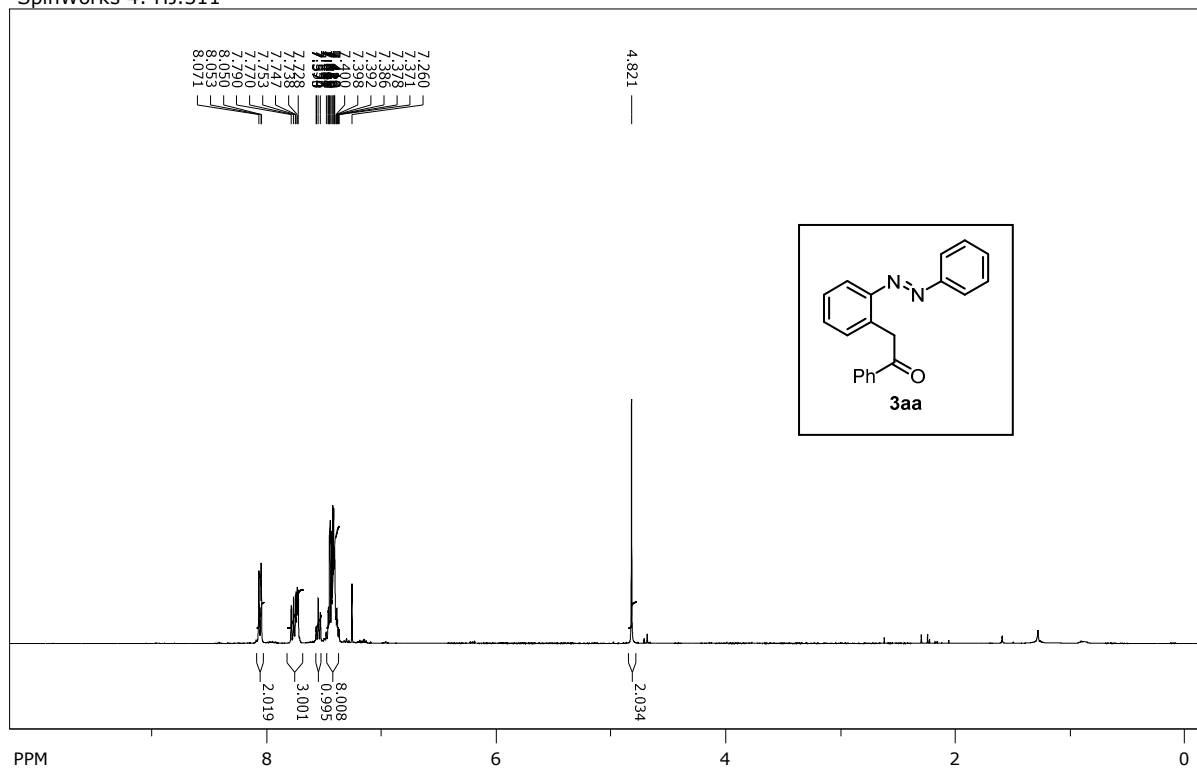
SpinWorks 4: HJ.370



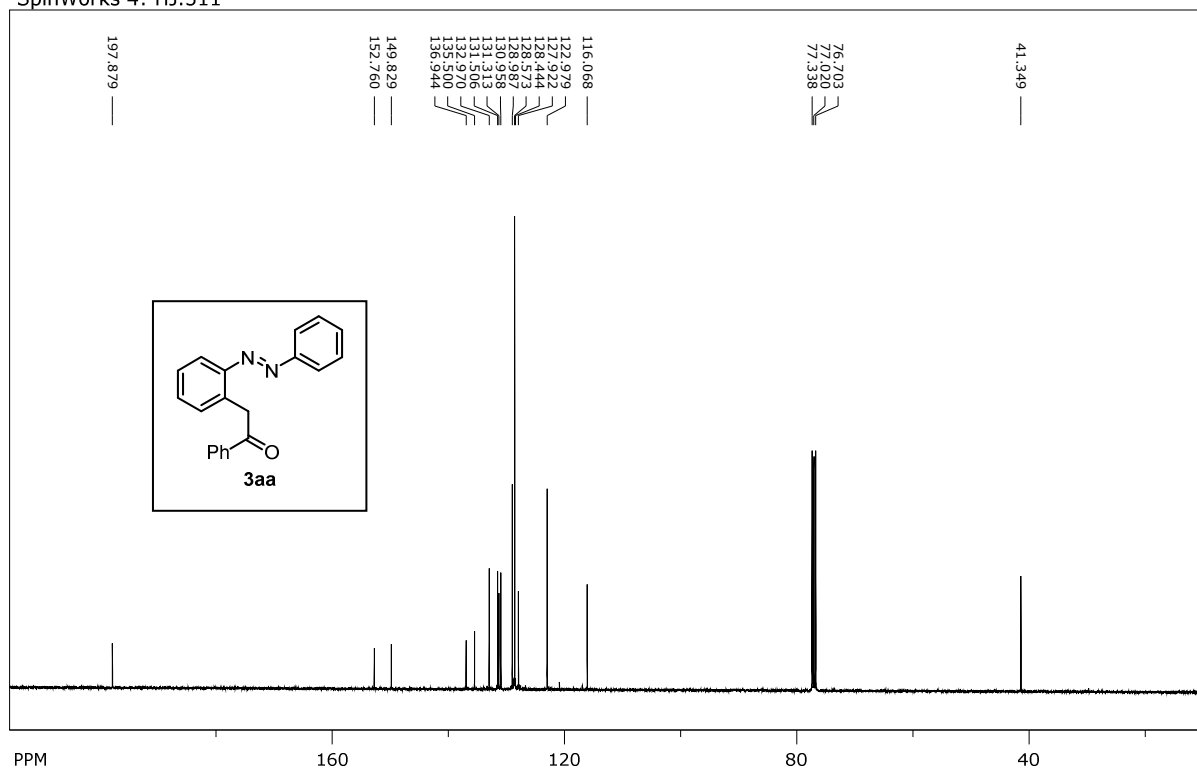
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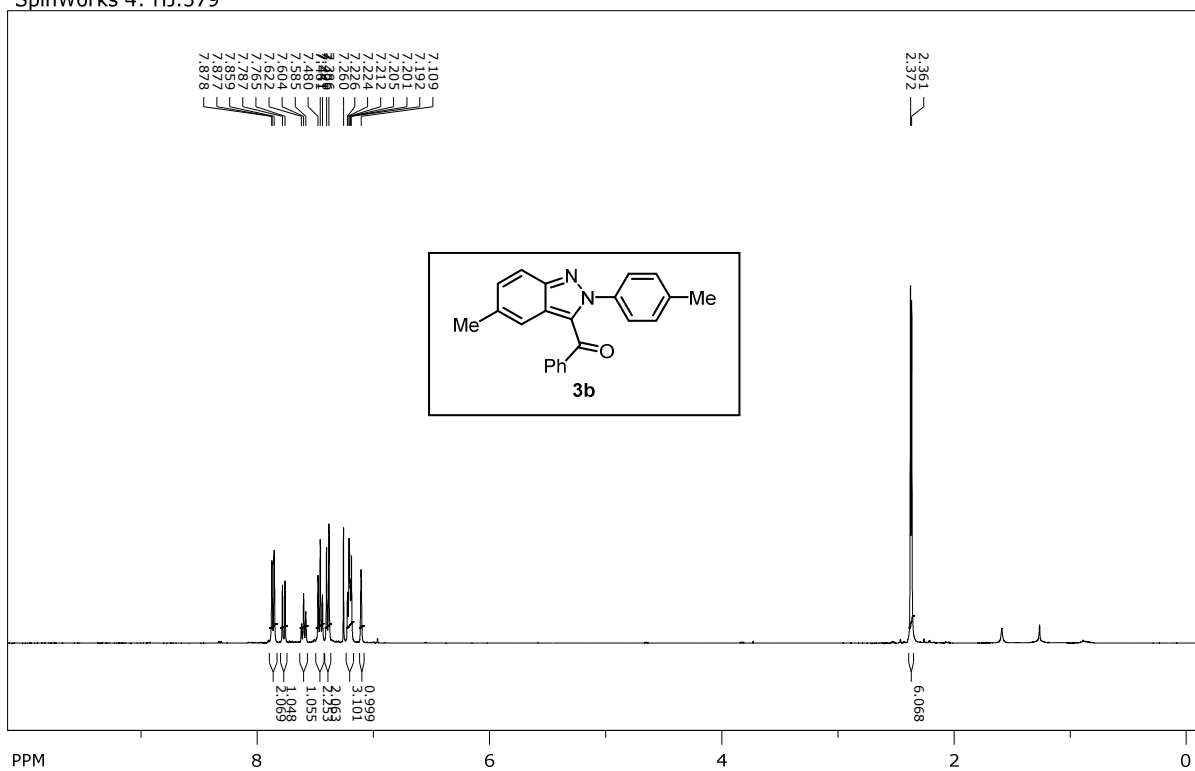
SpinWorks 4: HJ.311



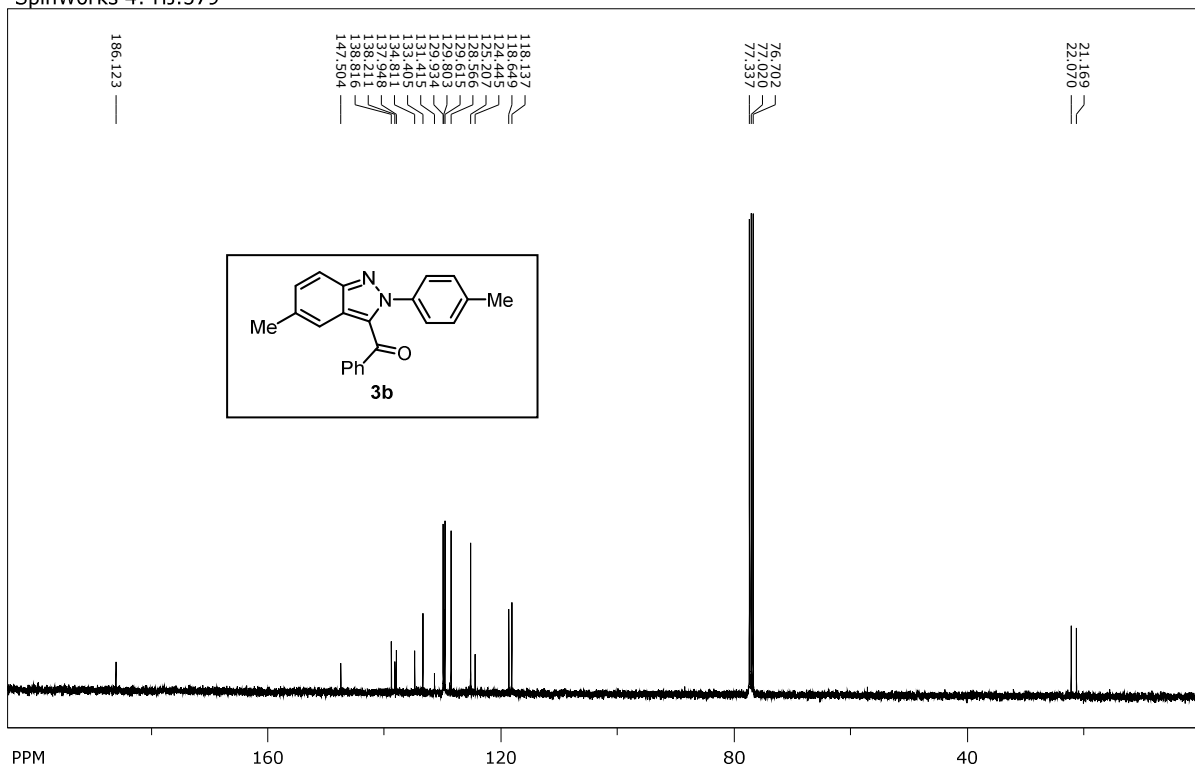
SpinWorks 4: HJ.311



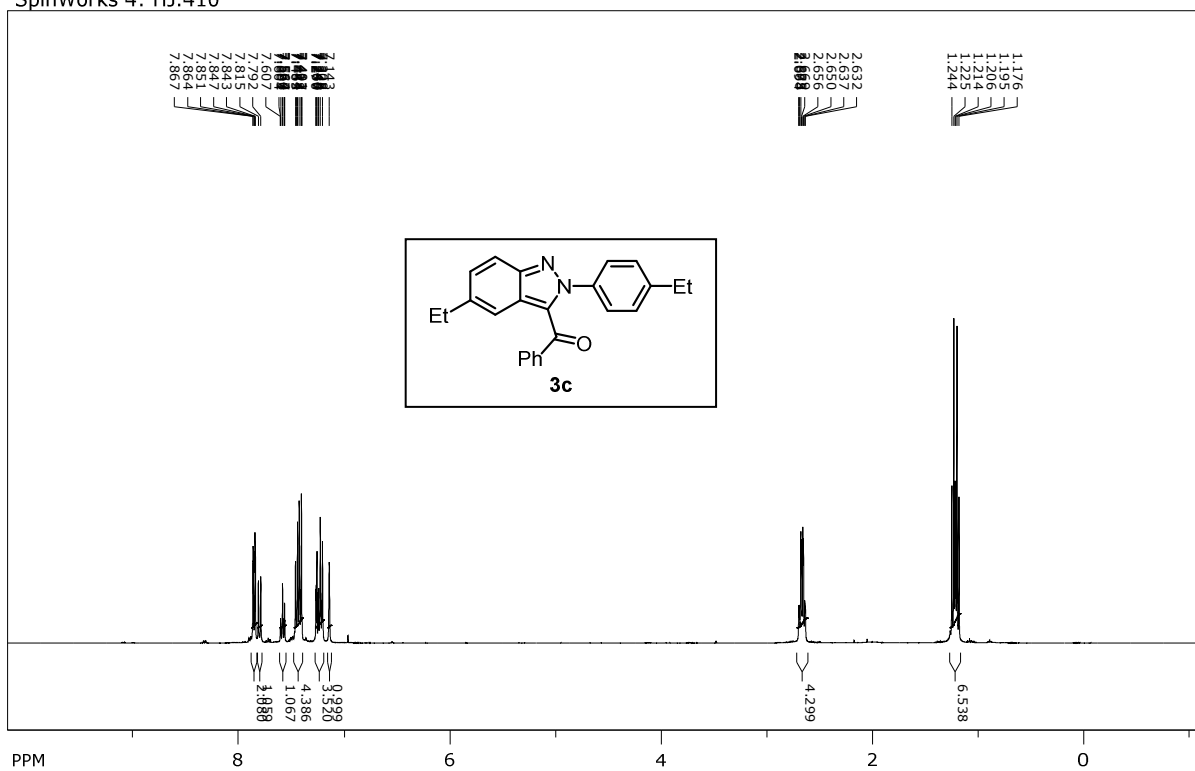
SpinWorks 4: HJ.379



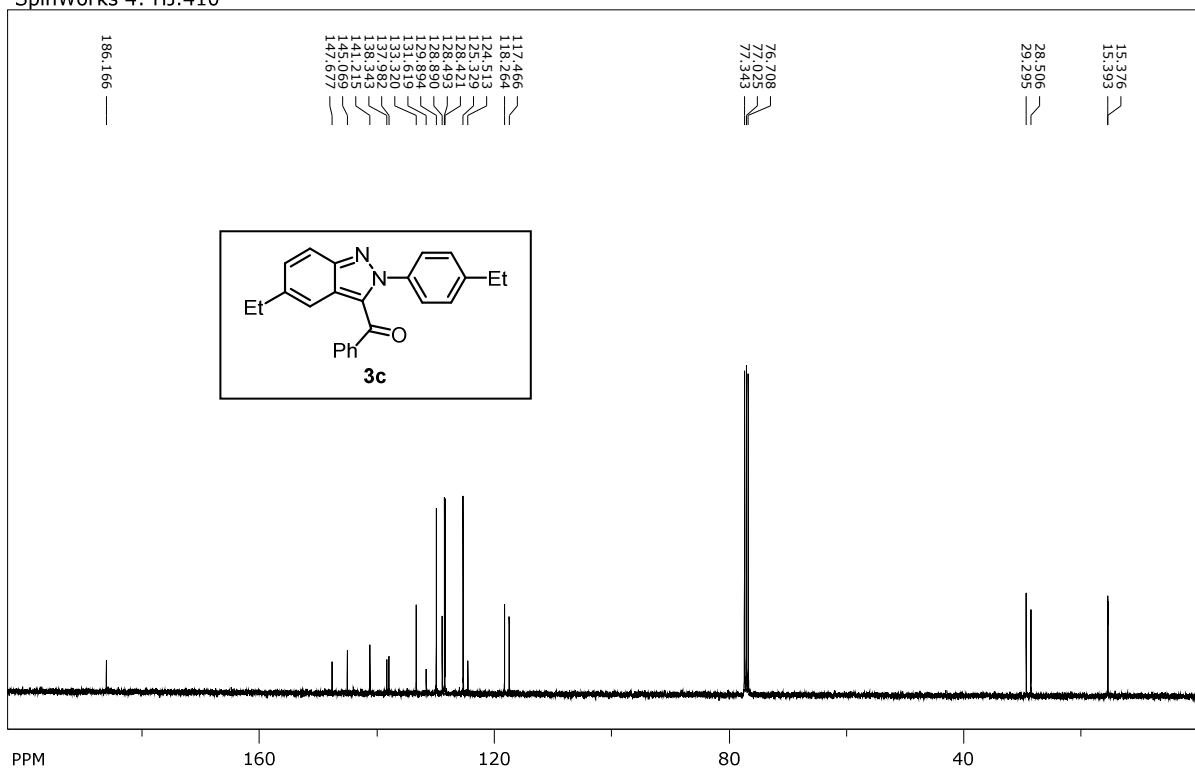
SpinWorks 4: HJ.379



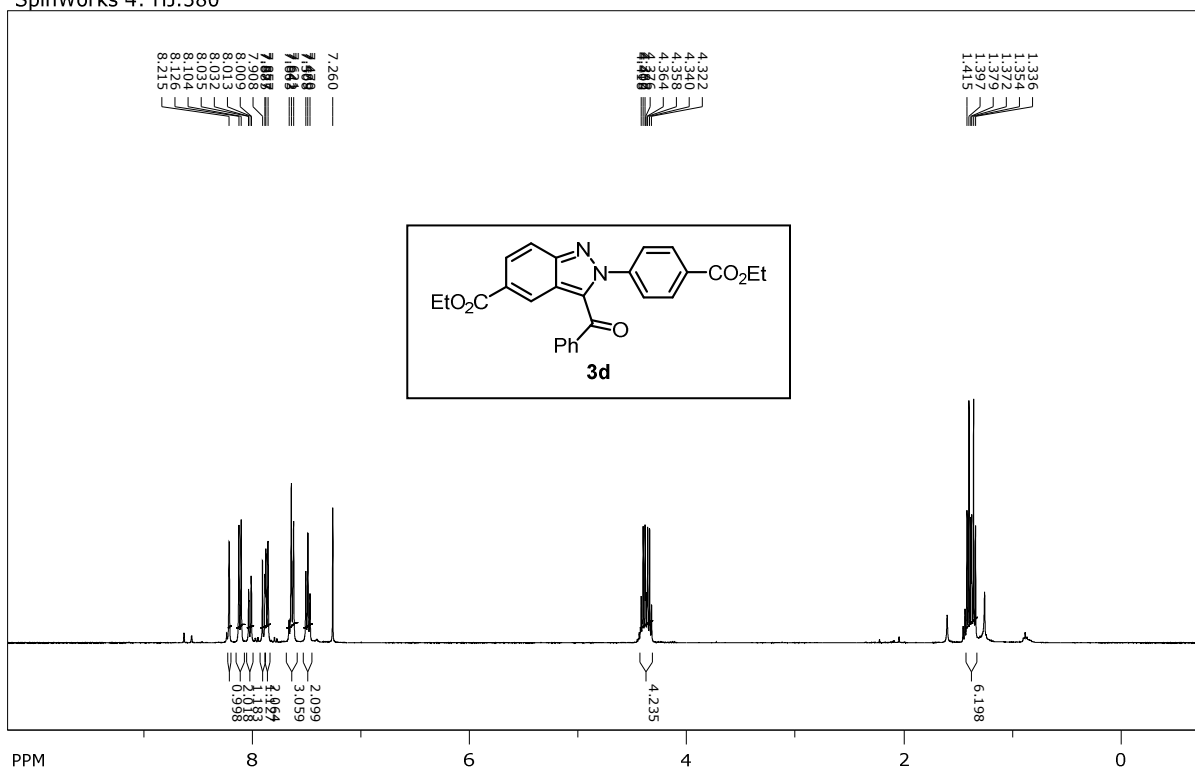
SpinWorks 4: HJ.410



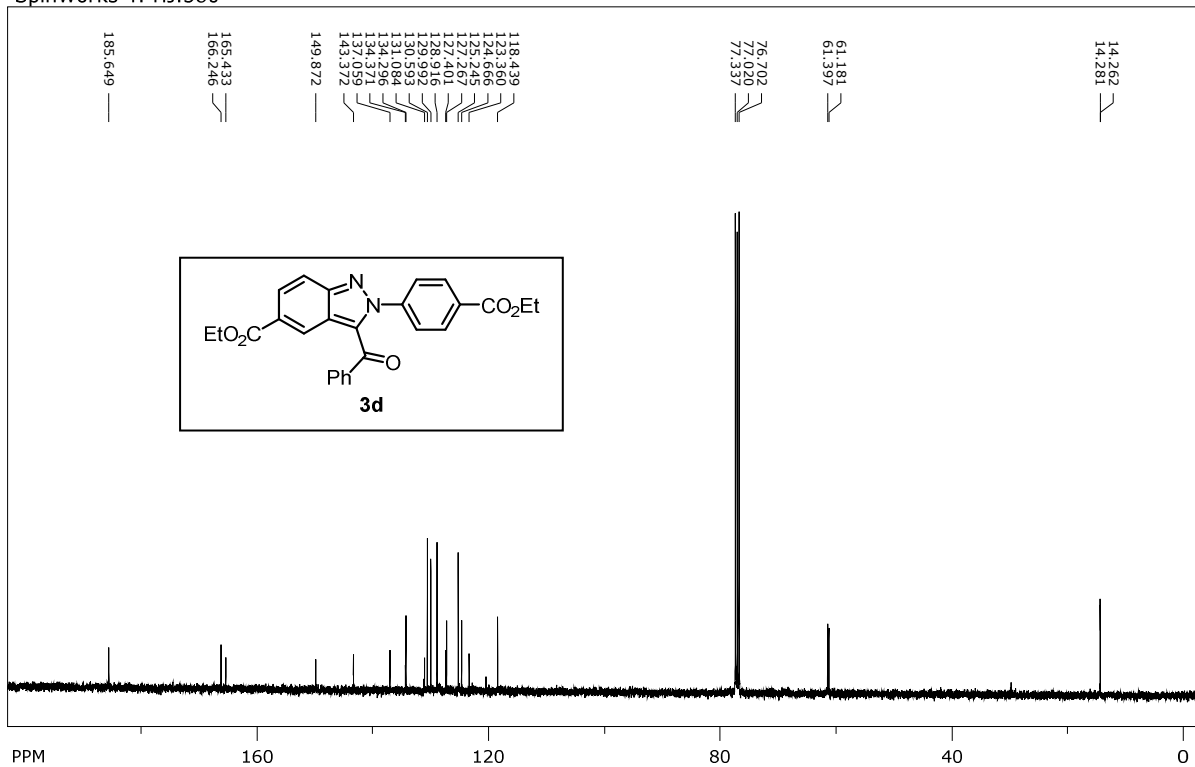
SpinWorks 4: HJ.410



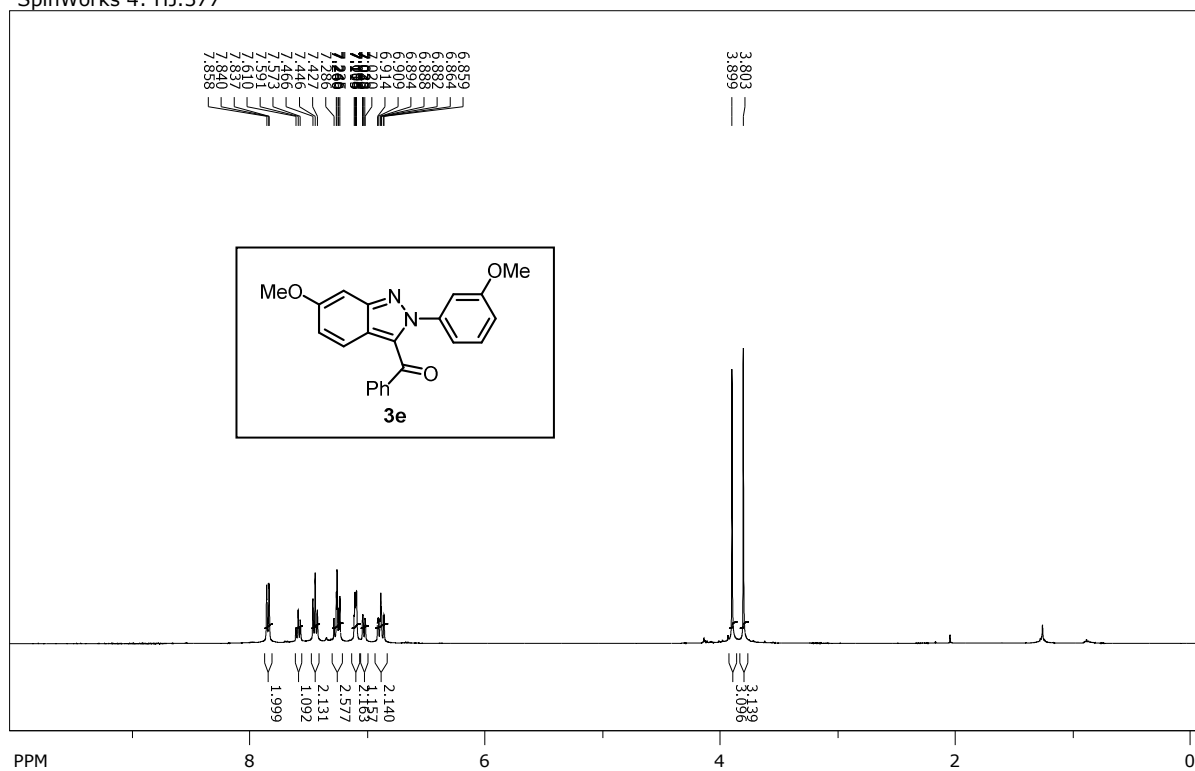
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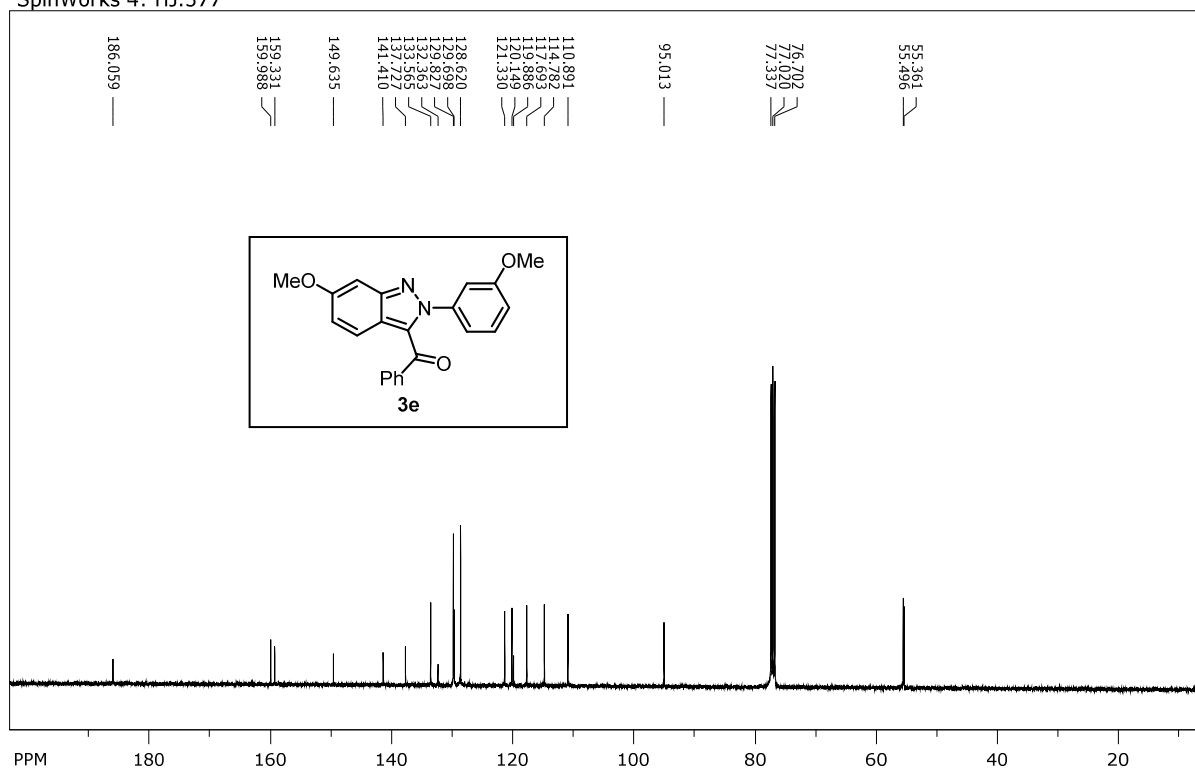
SpinWorks 4: HJ.380



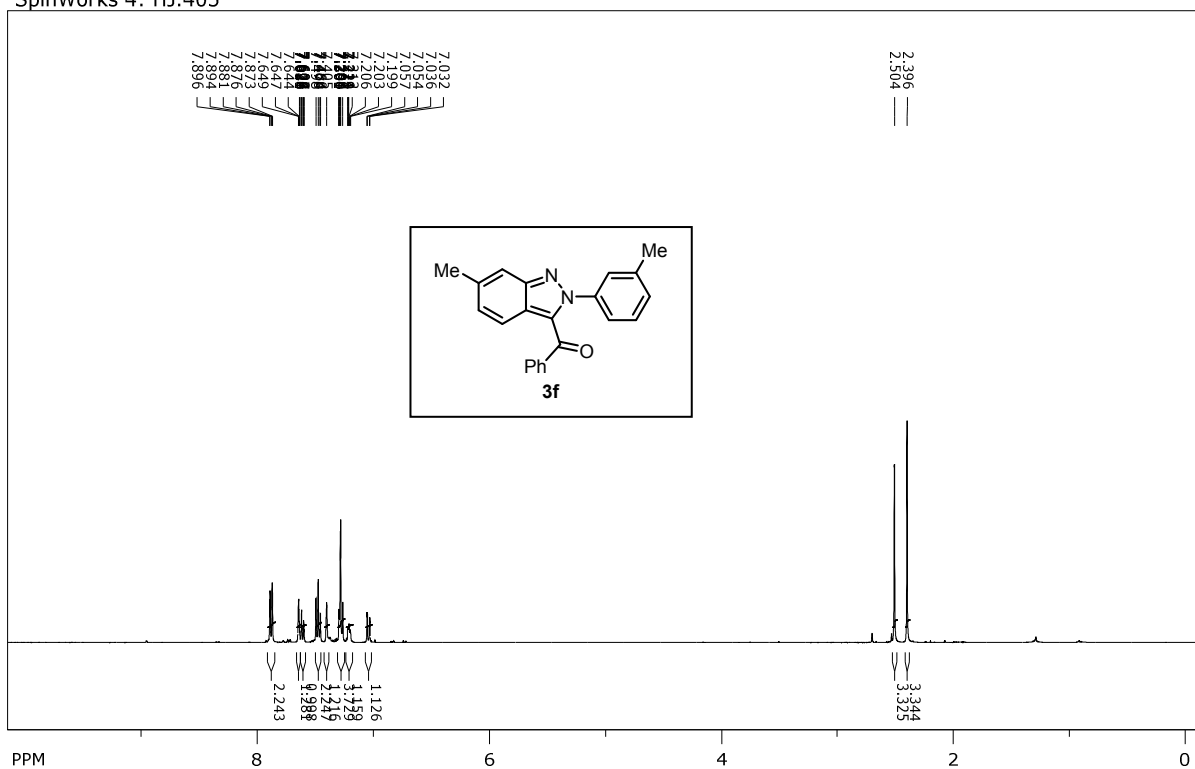
SpinWorks 4: HJ.377



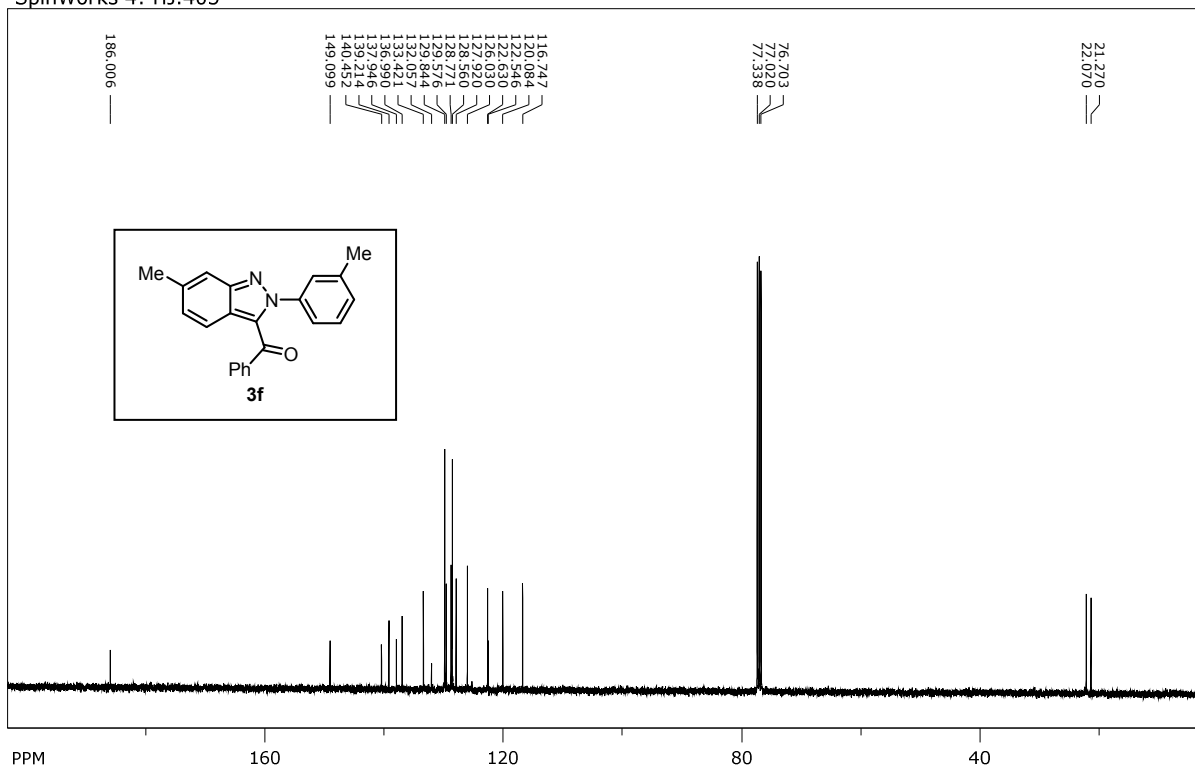
SpinWorks 4: HJ.377



SpinWorks 4: HJ.405



SpinWorks 4: HJ.405



[illegible]

SpinWorks 4: 15.462

Chemical structure of **3g** is shown in the top left:

O=C(c1ccc(Br)cc1)c2ccc(Br)cc2

The spectrum displays the following peak list (PPM):

Peak List (PPM)
185.407
149.238
141.122
137.329
134.164
132.540
129.066
129.840
130.251
132.286
128.658
128.658
128.658
124.215
122.619
122.540
121.477
121.477
120.839

Peak list (PPM) for the solvent region:

Peak List (PPM)
76.839
77.020
77.201

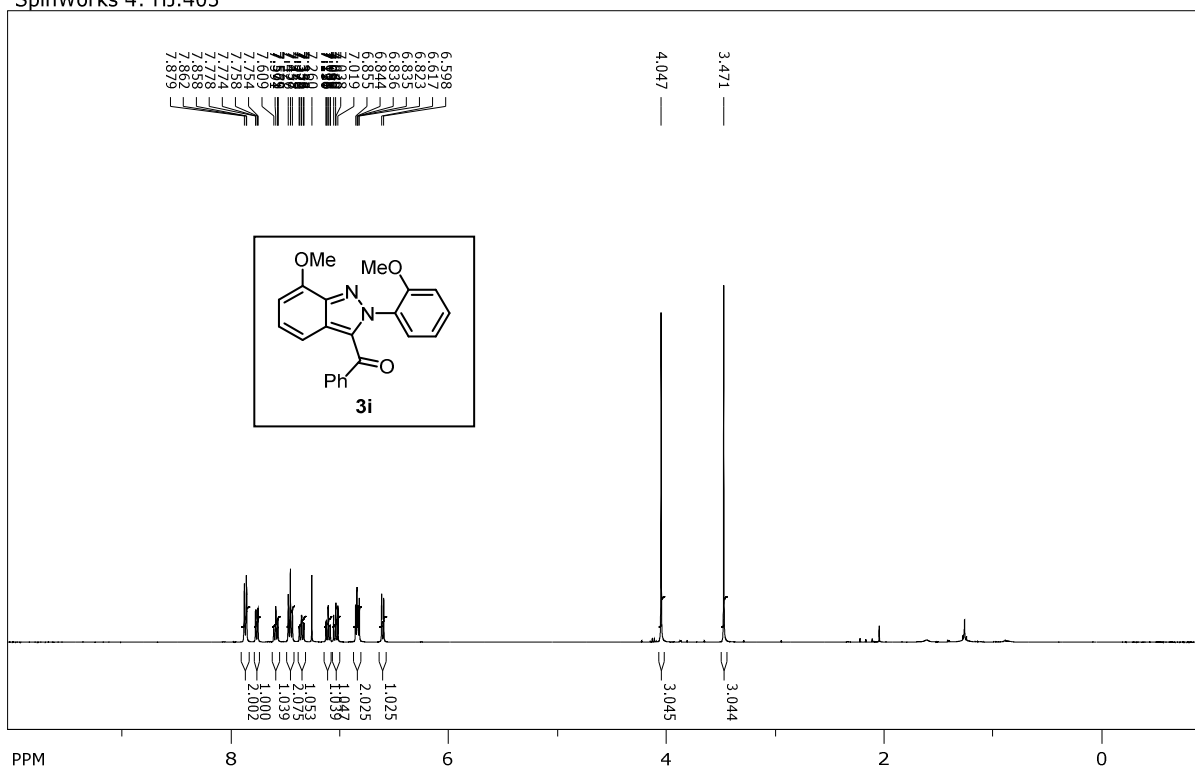
Chemical structure of 3h: O=C(c1ccc(C(F)(F)F)cc1)c2cc(F)ccn2

¹H NMR spectrum (CDCl₃):

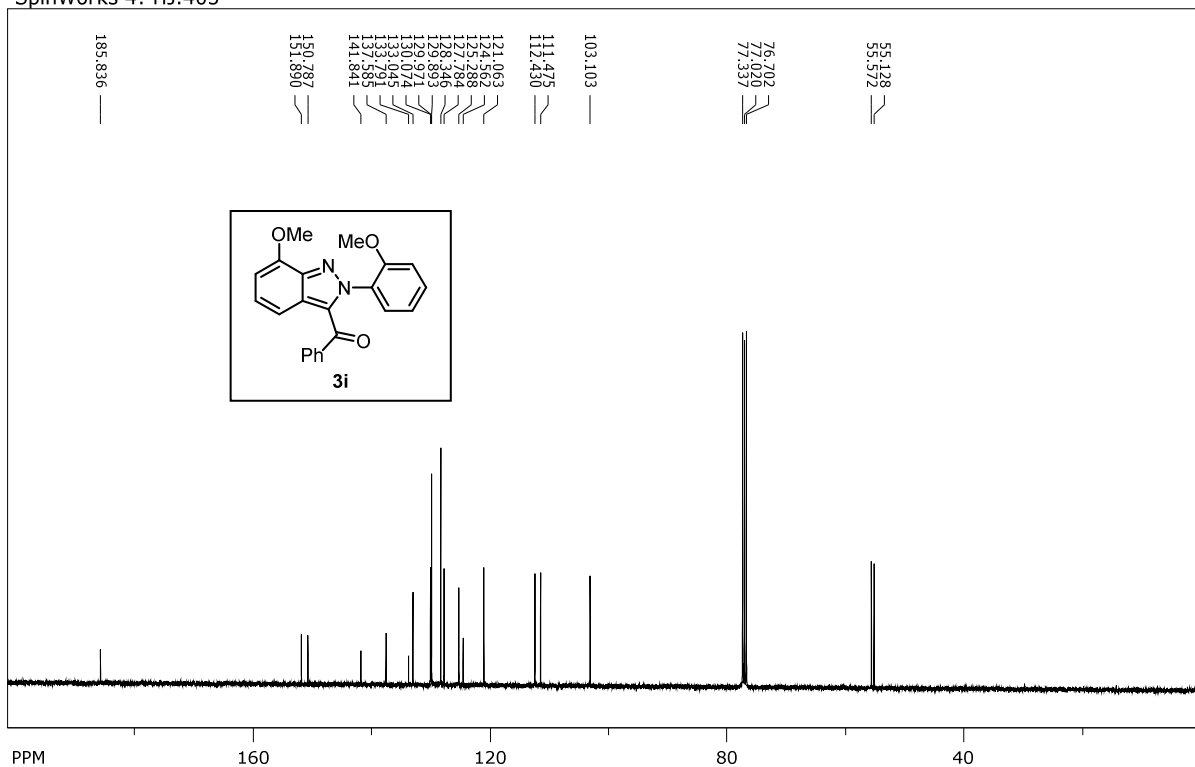
Chemical Shift (ppm)	Integration
7.887	2.002
7.869	2.075
7.753	2.062
7.635	4.186
7.617	1.034
7.464	1.000
7.435	
7.402	
7.382	
7.362	
7.149	
7.124	
7.103	
7.099	
6.837	
6.818	
6.791	

Chemical structure of **3h** is shown in the inset. The ¹³C NMR spectrum (CDCl₃) shows peaks at 186.664, 163.860, 156.144, 153.592, 150.559, 145.624, 141.027, 137.921, 134.374, 131.405, 129.611, 127.571, 125.721, 121.020, 120.866, 116.464, 116.253, 114.582, 114.532, 113.502, 110.799, 77.338, 77.020, and 76.703 ppm.

SpinWorks 4: HJ.403



SpinWorks 4: HJ.403



Spinworks 4.15.575

Chemical structure of **3j** is shown in the inset:

Cc1ccc(cc1)n2c(c3ccccc3C(=O)c24ccccc44)c5ccccc51

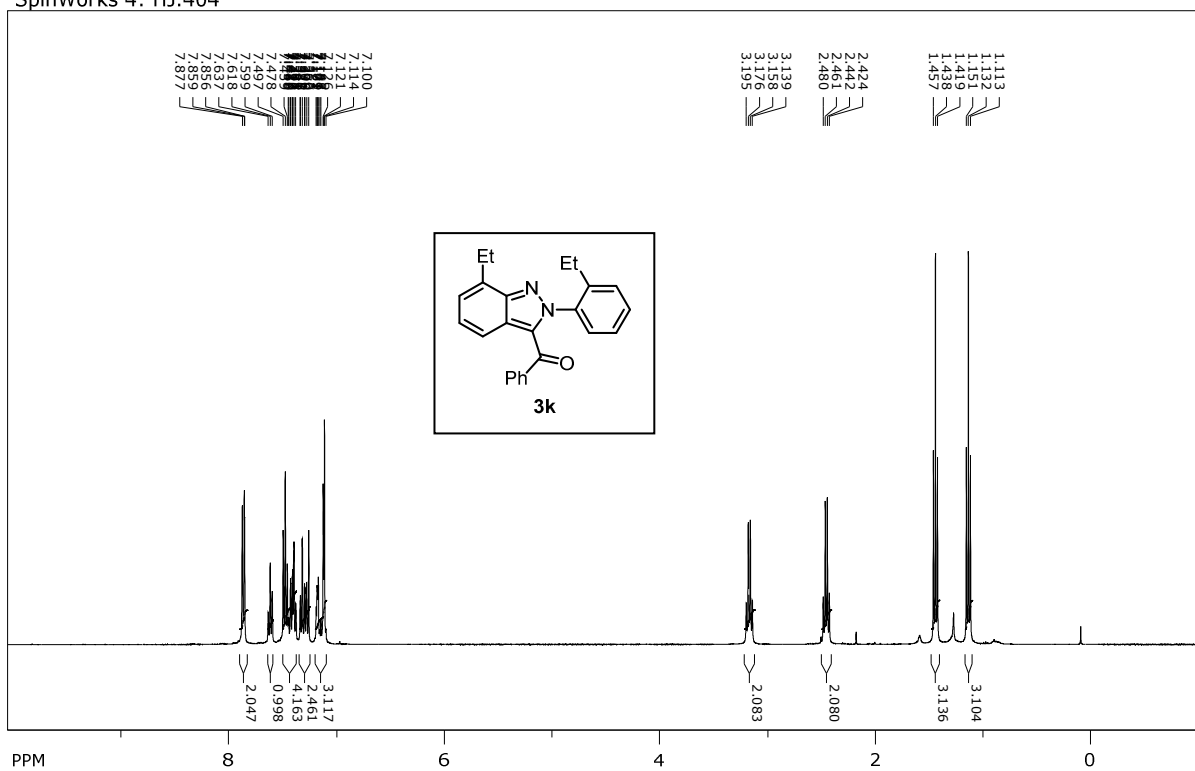
¹H NMR spectrum (CDCl₃) of compound **3j** is displayed. The spectrum shows peaks at 2.129, 2.725, and a multiplet between 7.0 and 7.9 ppm. Integration values are 3.069, 3.208, and 2.080, 0.966, 2.169, 4.154, 3.278 respectively.

13C NMR spectrum of compound **3j**. The chemical structure of **3j** is shown in the top left: 1-(2-methylphenyl)-2-phenyl-1H-indole-3-carboxamide.

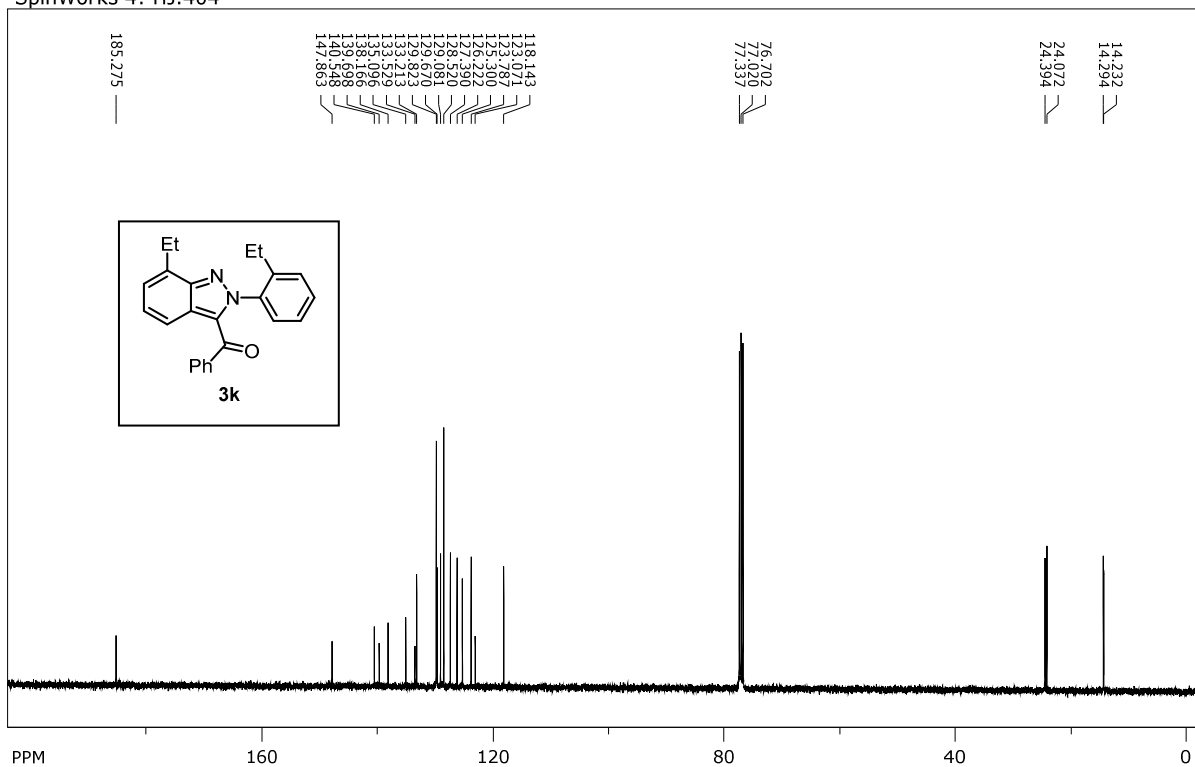
The spectrum displays the following chemical shifts (ppm):

- 185.344
- 148.691
- 140.227
- 137.995
- 134.815
- 133.462
- 130.850
- 129.500
- 129.100
- 126.274
- 126.116
- 125.809
- 125.298
- 122.432
- 118.140
- 77.704
- 77.339
- 17.113
- 17.605

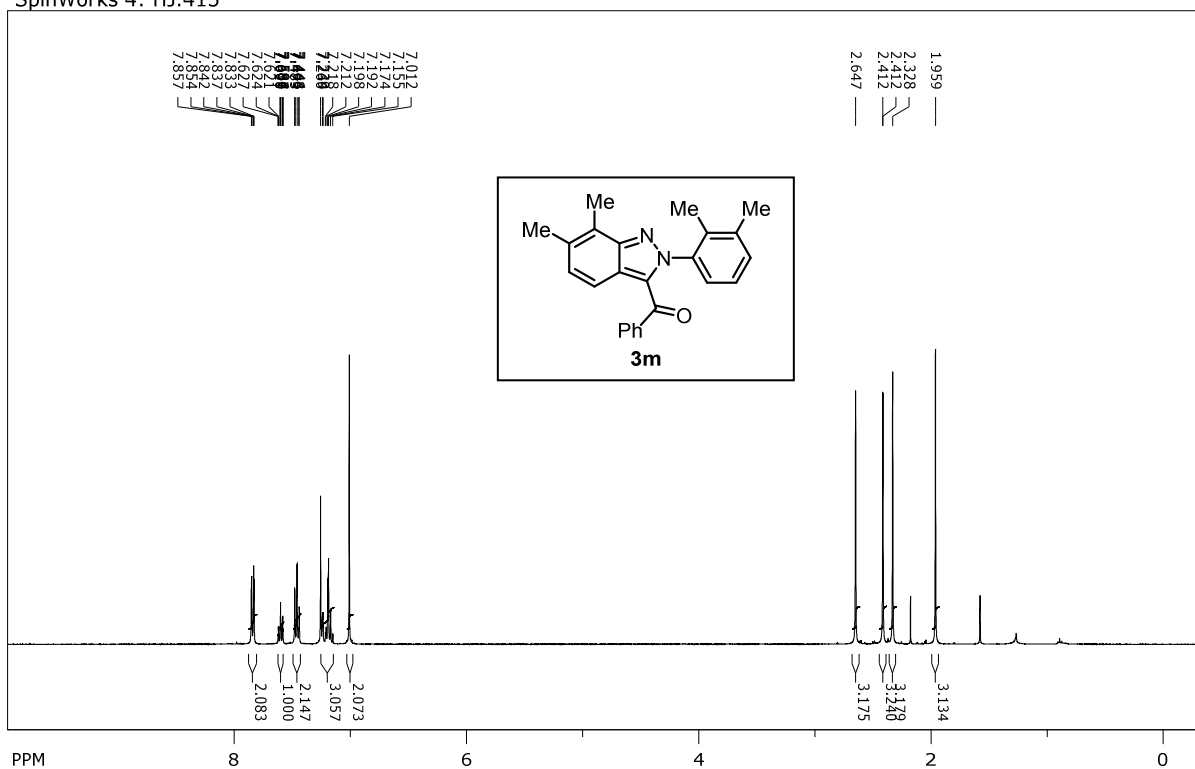
SpinWorks 4: HJ.404



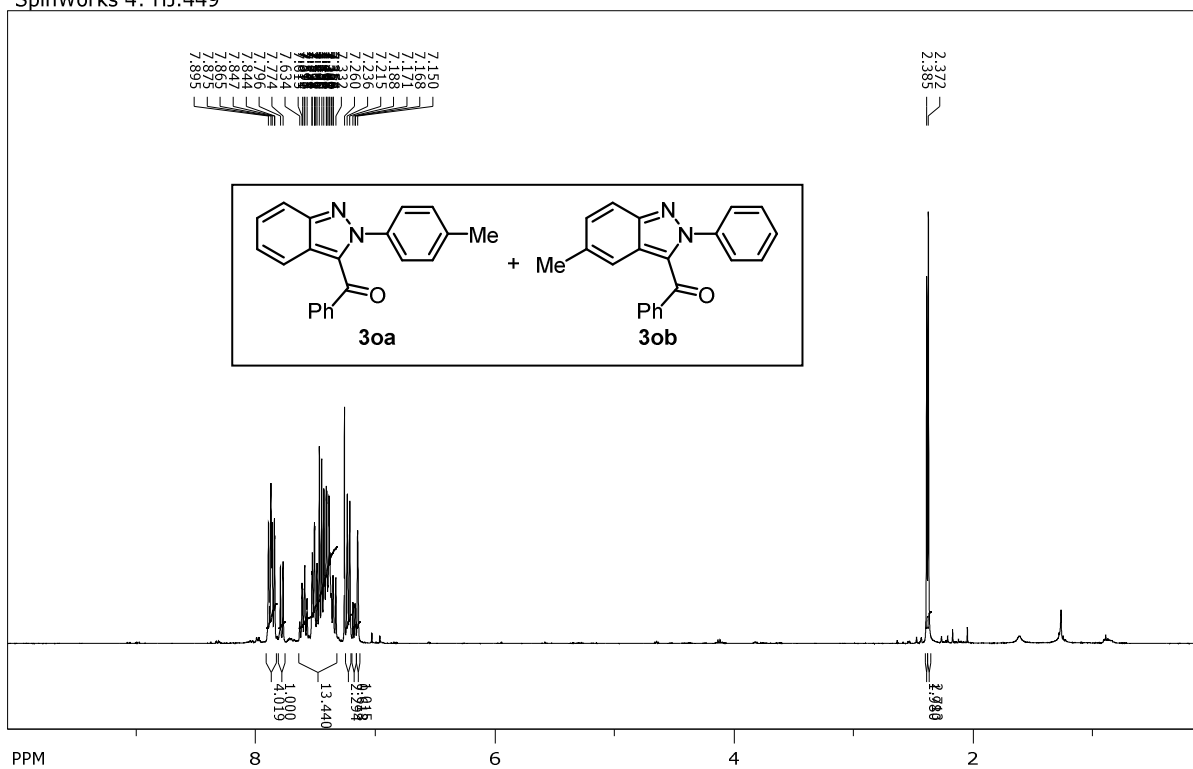
SpinWorks 4: HJ.404



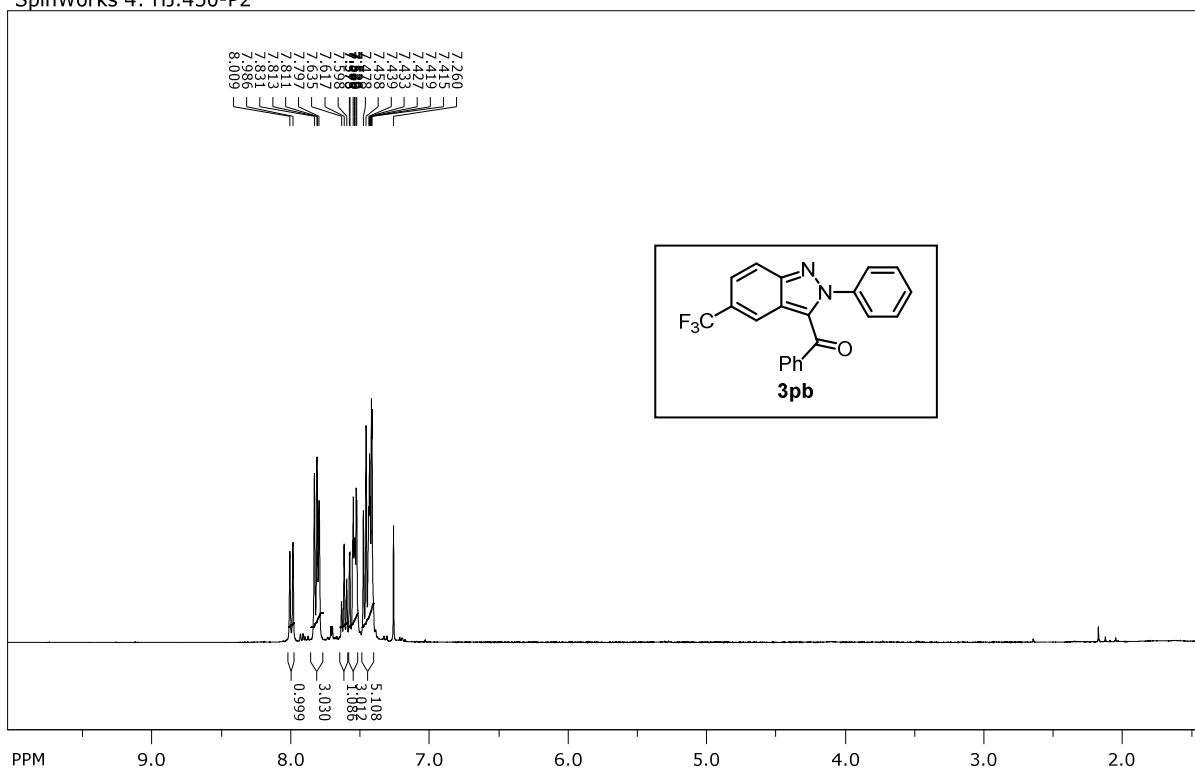
SpinWorks 4: HJ.415



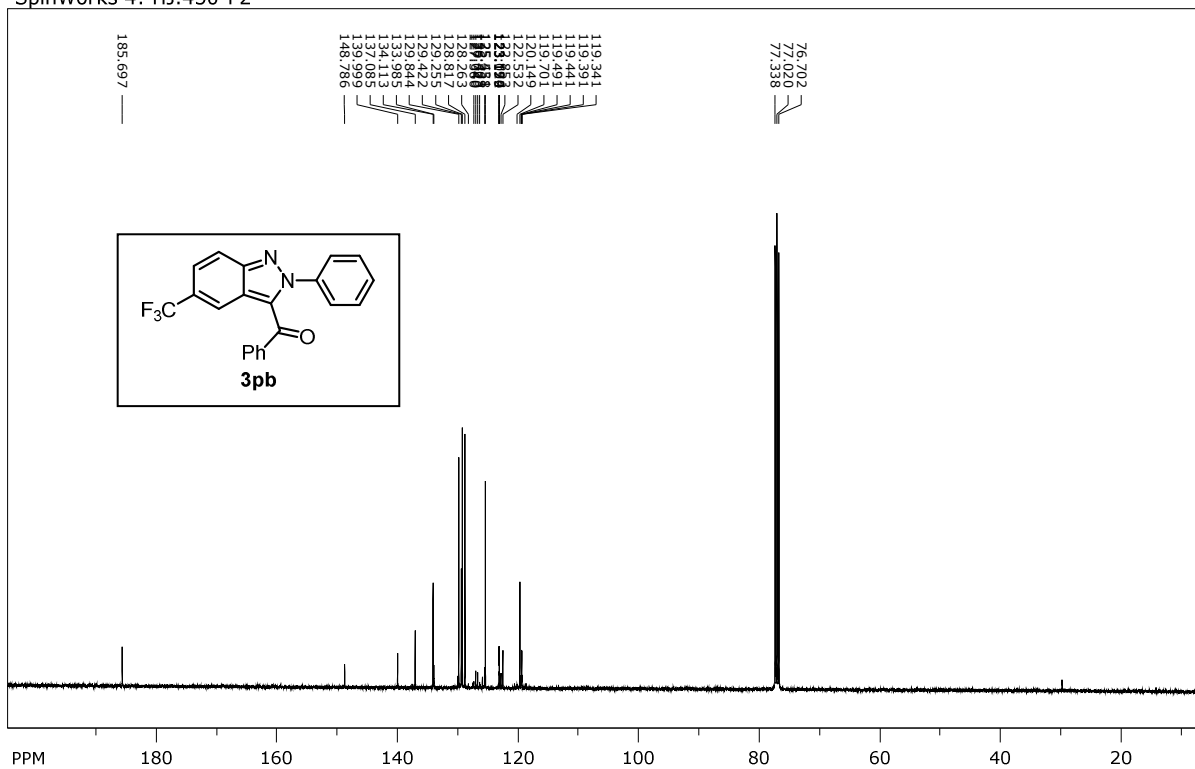
SpinWorks 4: HJ.449



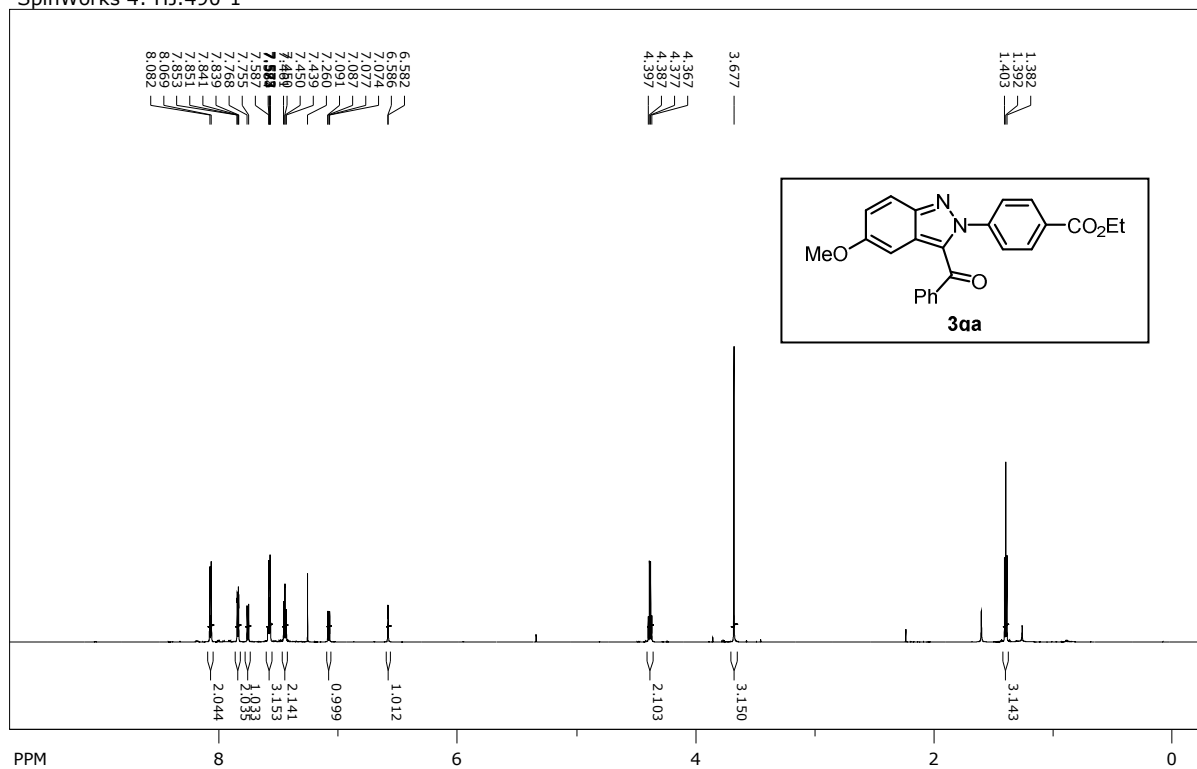
SpinWorks 4: HJ.450-P2



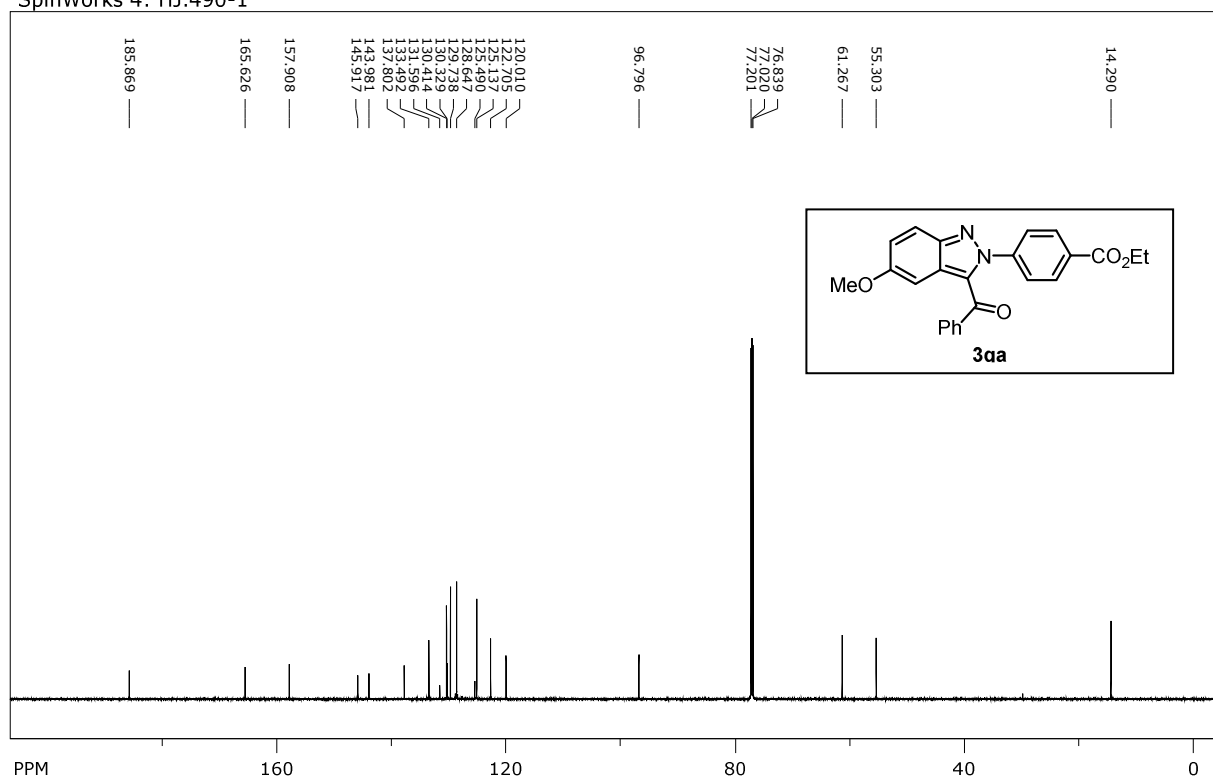
SpinWorks 4: HJ.450-P2



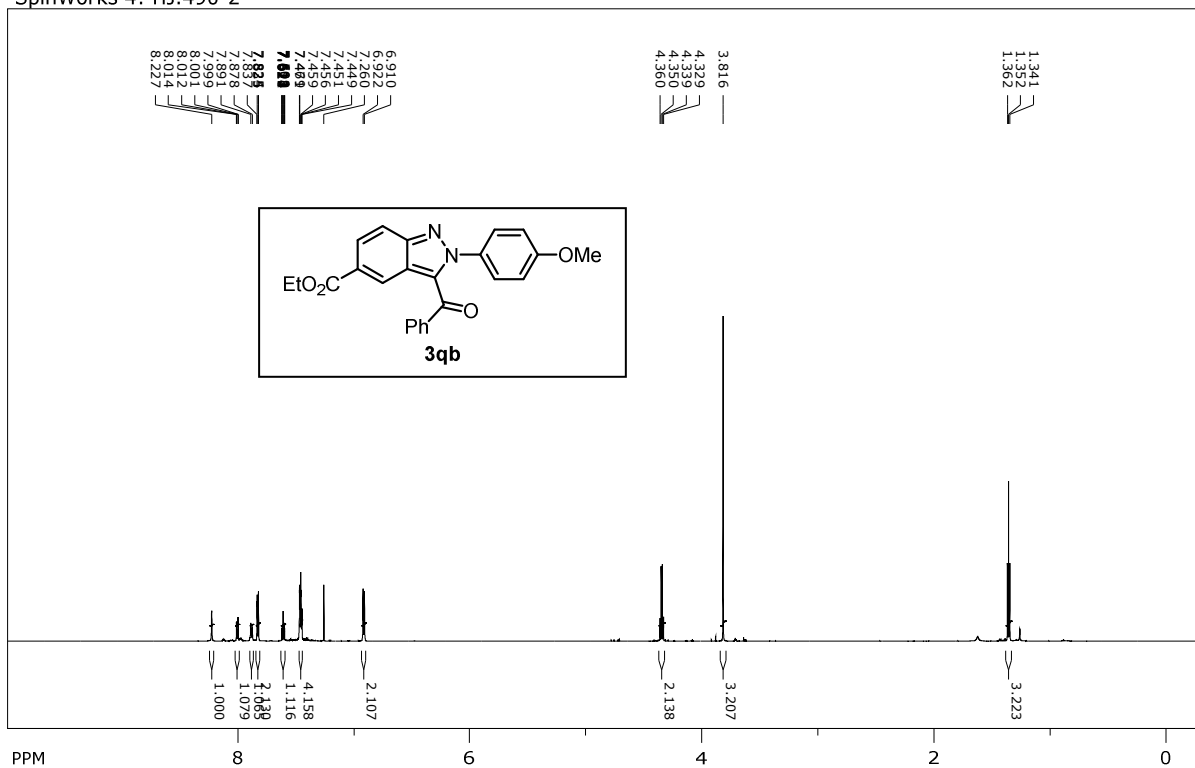
SpinWorks 4: HJ.490-1



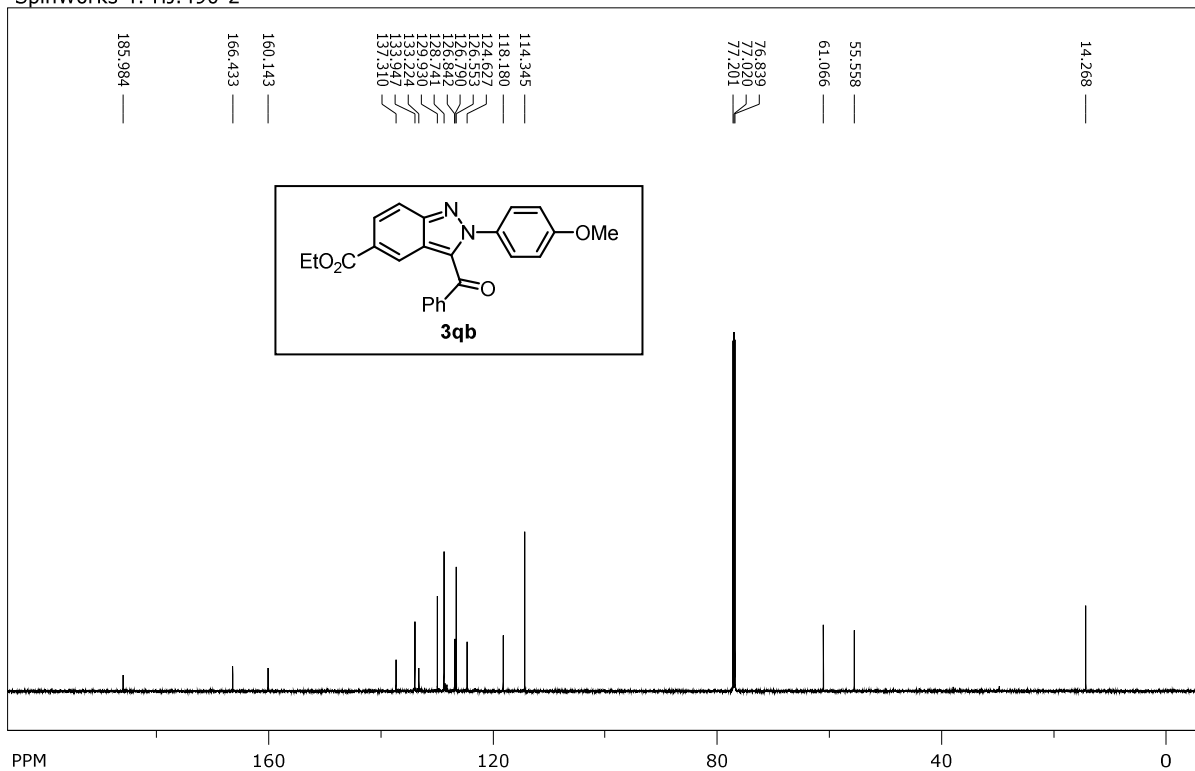
SpinWorks 4: HJ.490-1



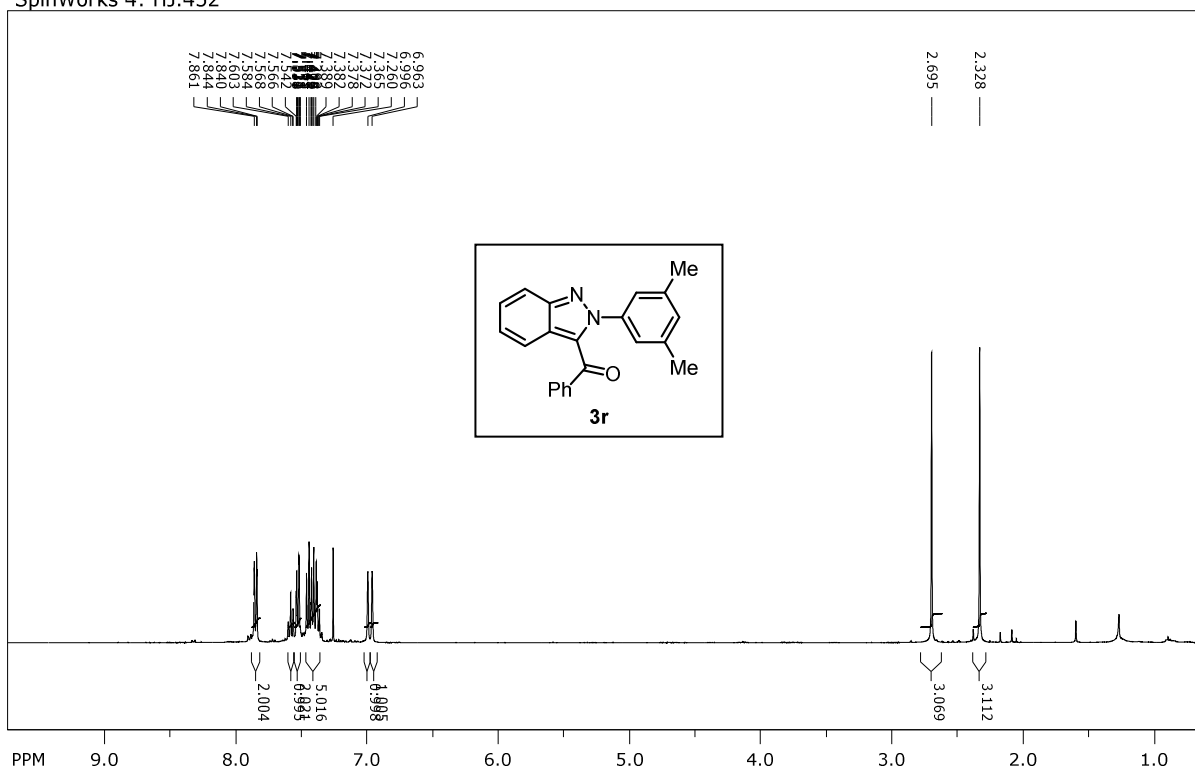
SpinWorks 4: HJ.490-2



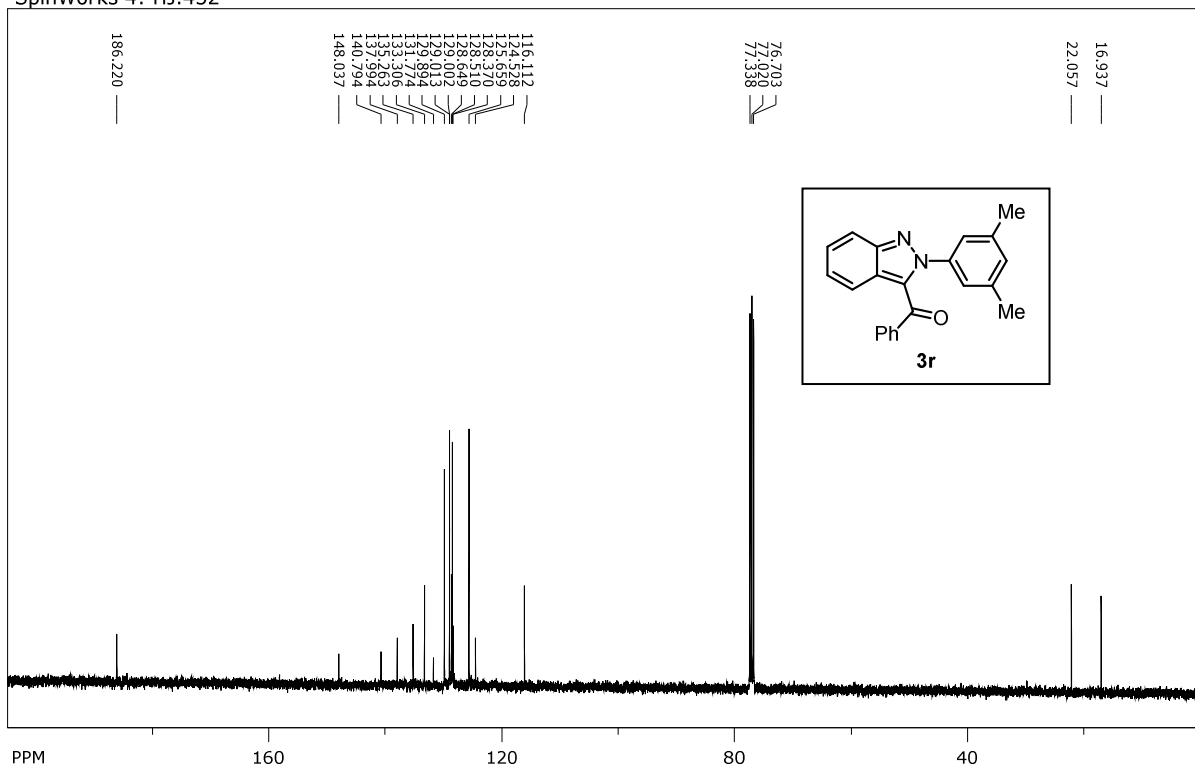
SpinWorks 4: HJ.490-2



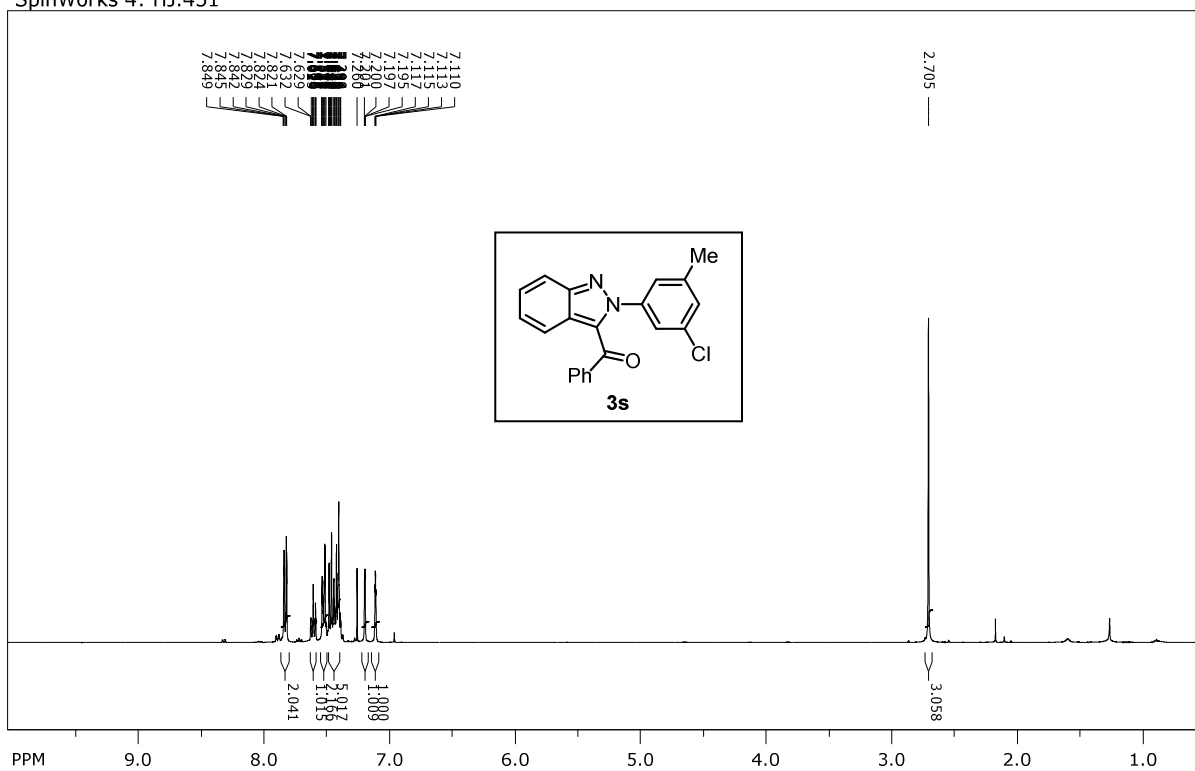
SpinWorks 4: HJ.452



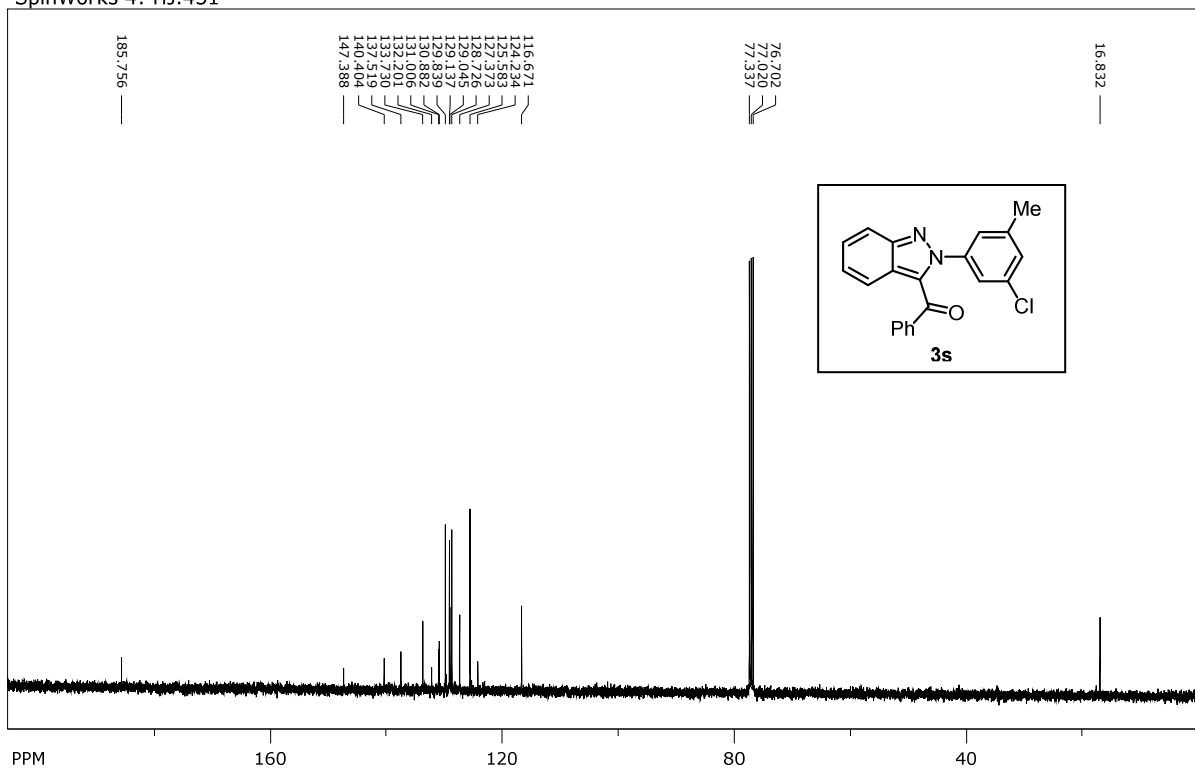
SpinWorks 4: HJ.452



SpinWorks 4: HJ.451



SpinWorks 4: HJ.451



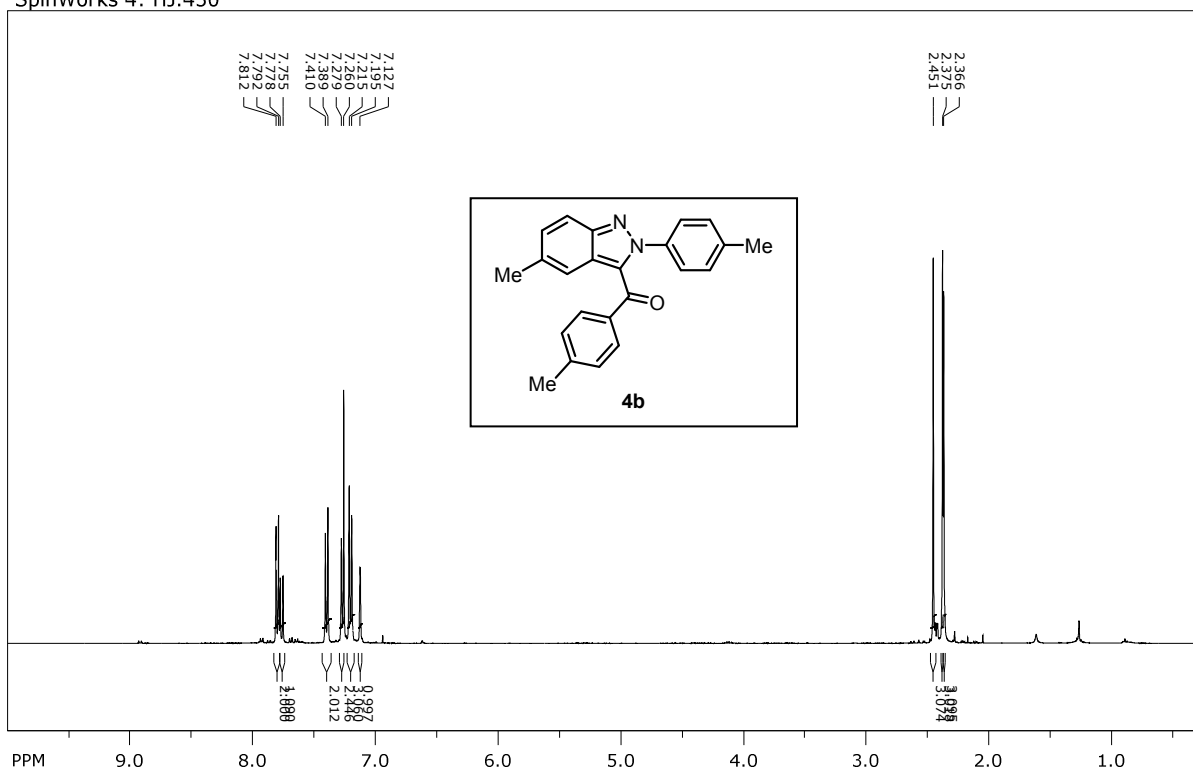
Chemical Structure of 3tb: 1-(4-methylphenyl)-2-phenyl-1H-benzotriazole

¹H NMR Spectrum (CDCl₃):

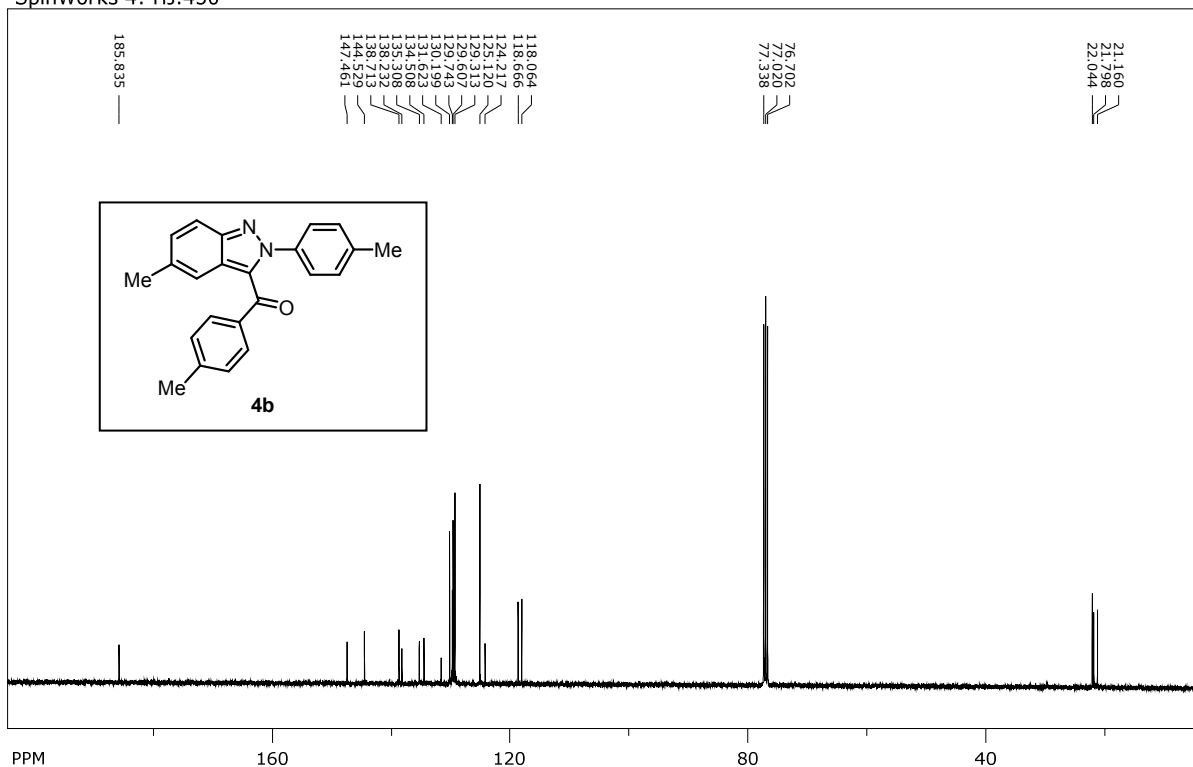
- Peak List (Top):** 7.917, 7.910, 7.903, 7.900, 7.894, 7.891, 7.887, 7.882, 7.875, 7.872, 7.869, 7.866, 7.863, 7.860, 7.857, 7.854, 7.851, 7.848, 7.845, 7.842, 7.839, 7.836, 7.833.
- Peak List (Bottom):** 7.995, 7.992, 7.989, 7.986, 7.983, 7.980, 7.977, 7.974, 7.971, 7.968, 7.965, 7.962, 7.959, 7.956, 7.953, 7.950, 7.947, 7.944, 7.941, 7.938, 7.935, 7.932, 7.929, 7.926, 7.923, 7.920, 7.917, 7.914, 7.911, 7.908, 7.905, 7.902, 7.899, 7.896, 7.893, 7.890, 7.887, 7.884, 7.881, 7.878, 7.875, 7.872, 7.869, 7.866, 7.863, 7.860, 7.857, 7.854, 7.851, 7.848, 7.845, 7.842, 7.839, 7.836, 7.833.
- Integration Values:** 1.048, 5.328, 4.981, 2.994.

[illegible]

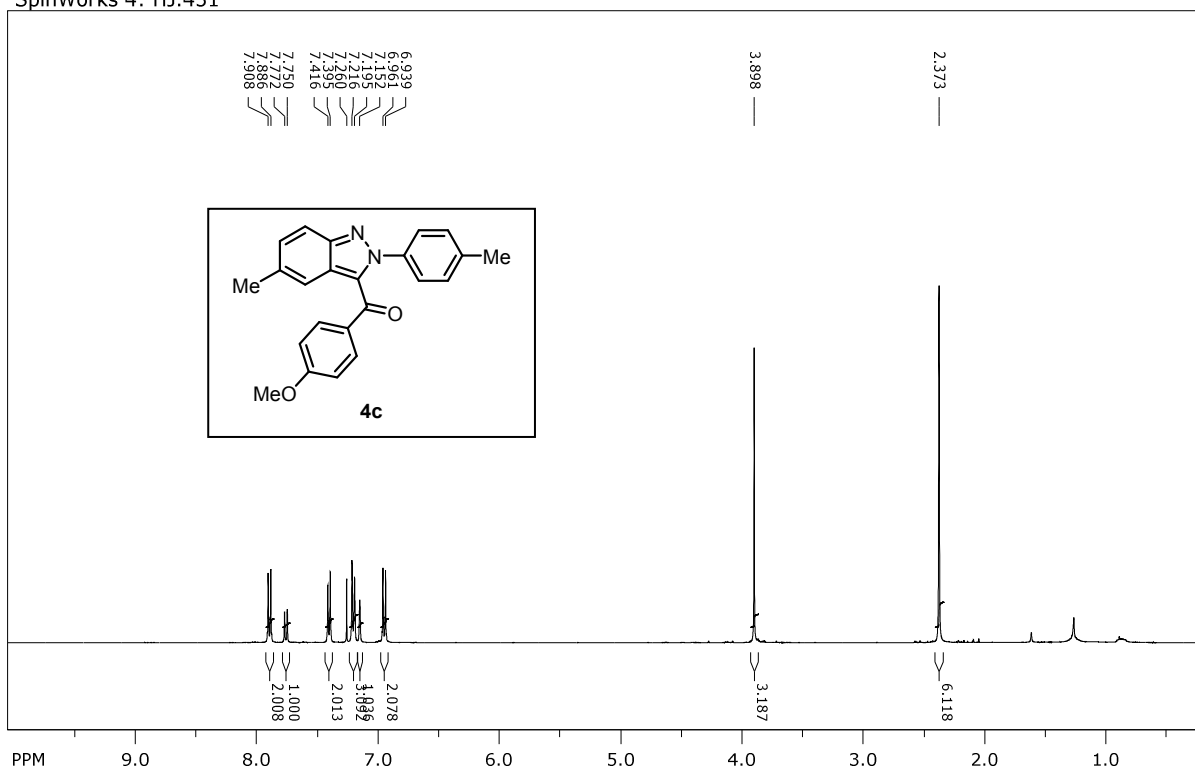
SpinWorks 4: HJ.430



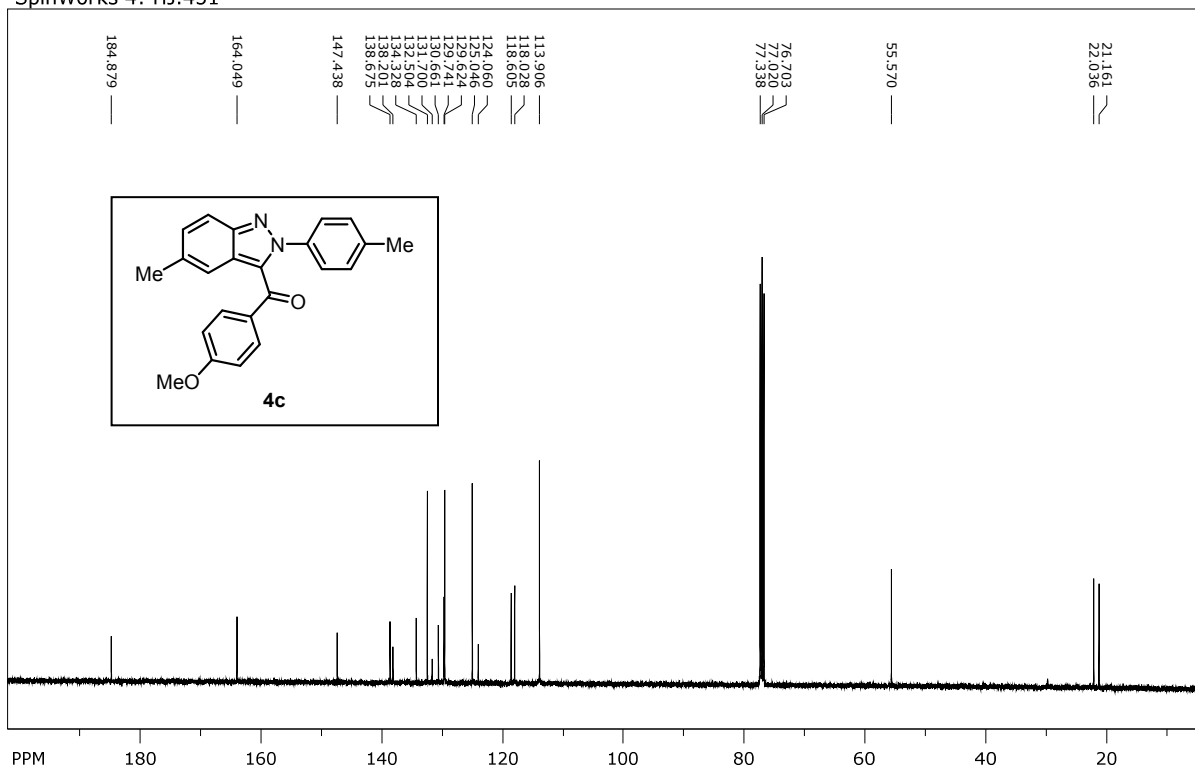
SpinWorks 4: HJ.430



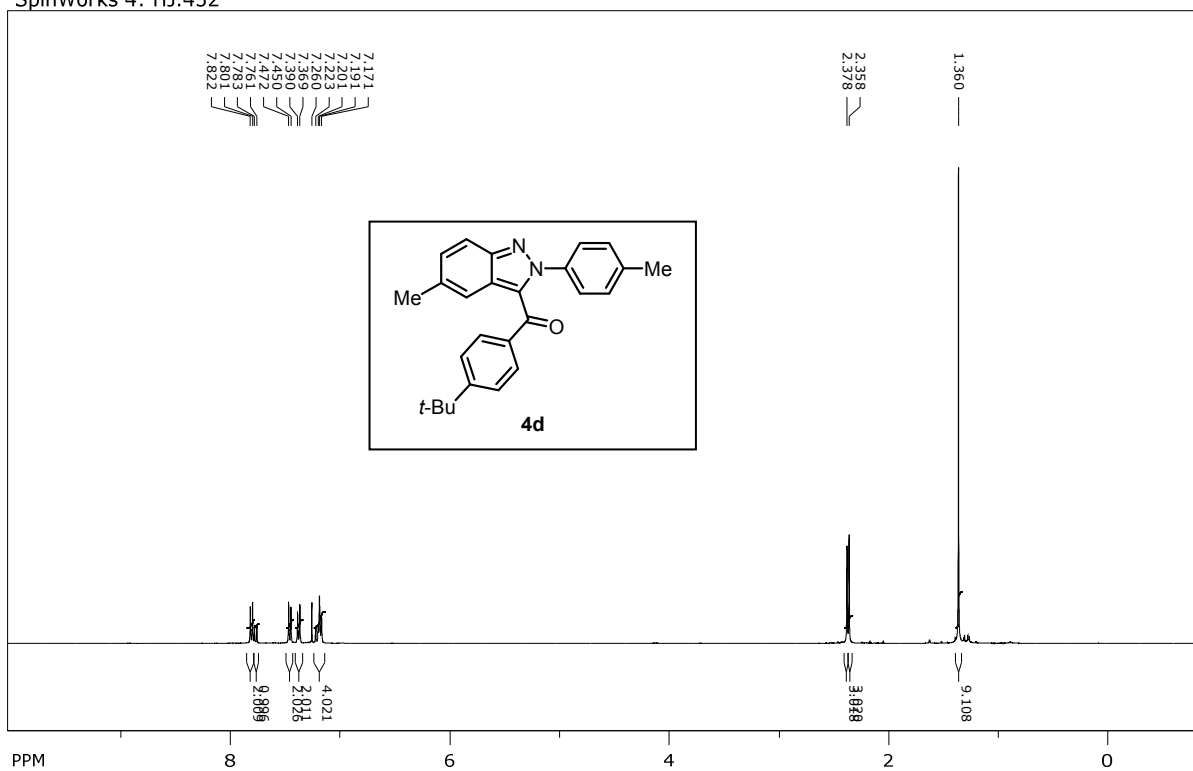
SpinWorks 4: HJ.431



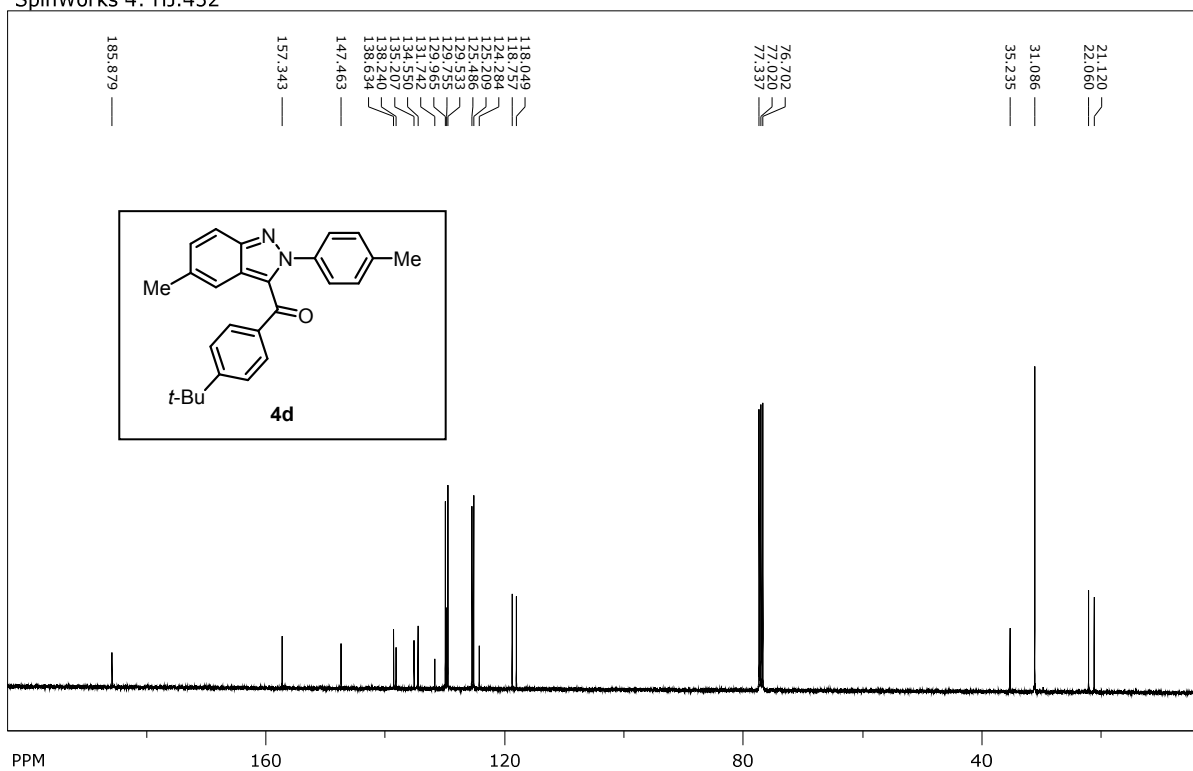
SpinWorks 4: HJ.431



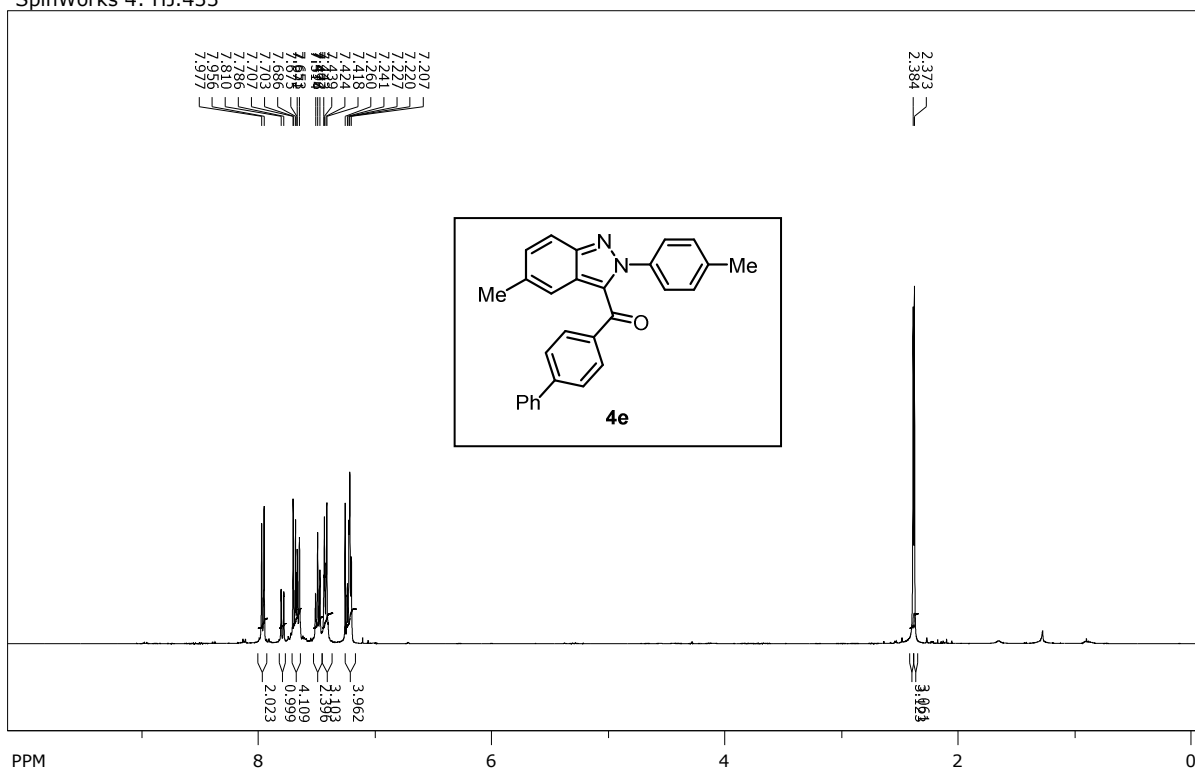
SpinWorks 4: HJ.432



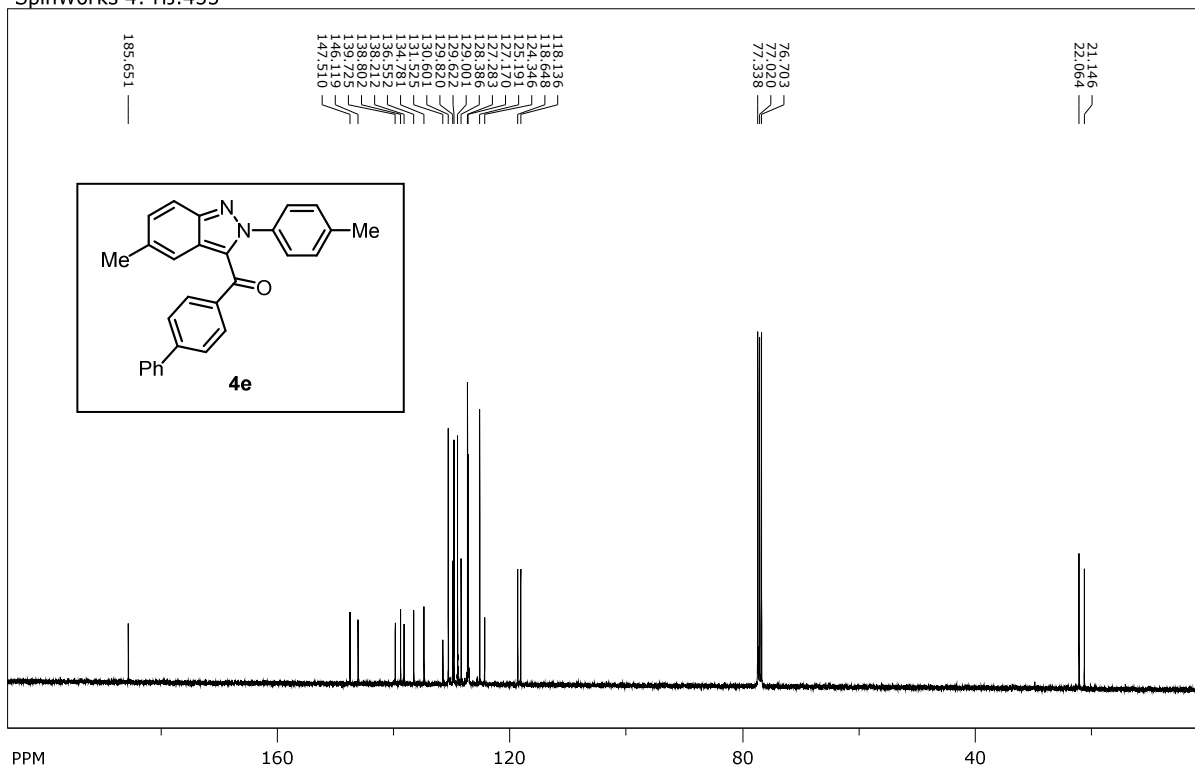
SpinWorks 4: HJ.432



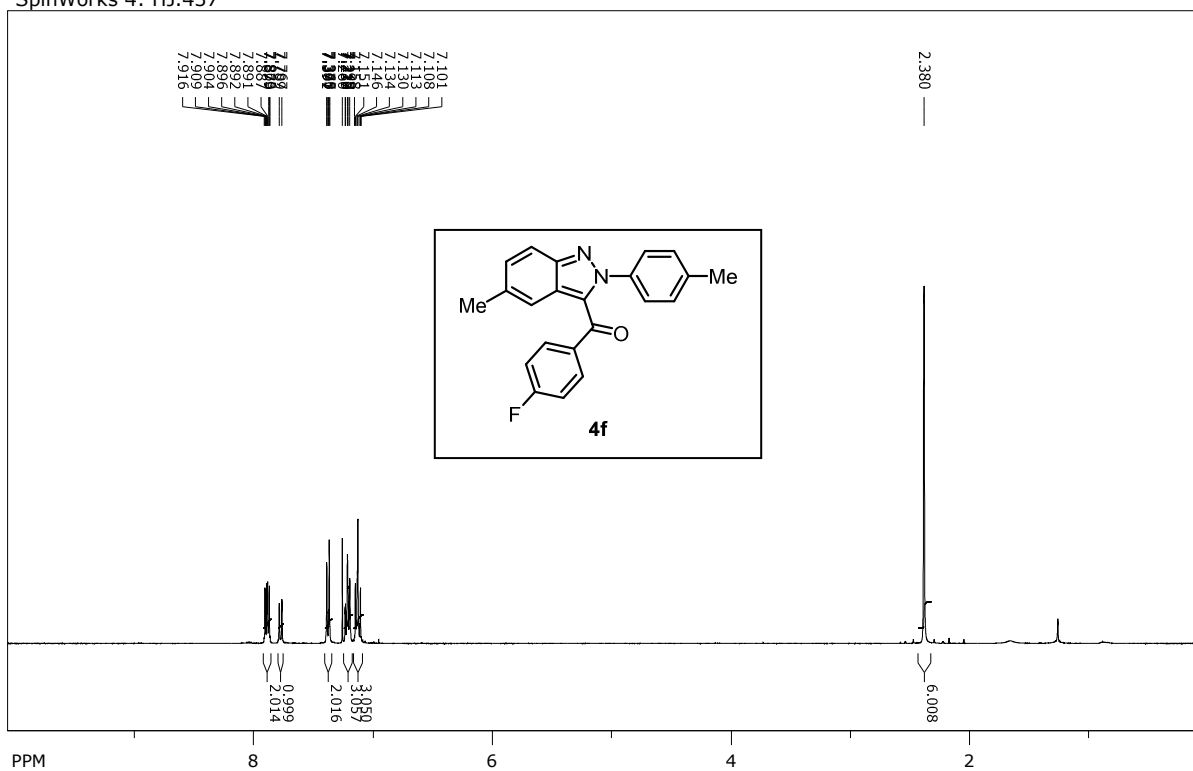
SpinWorks 4: HJ.433



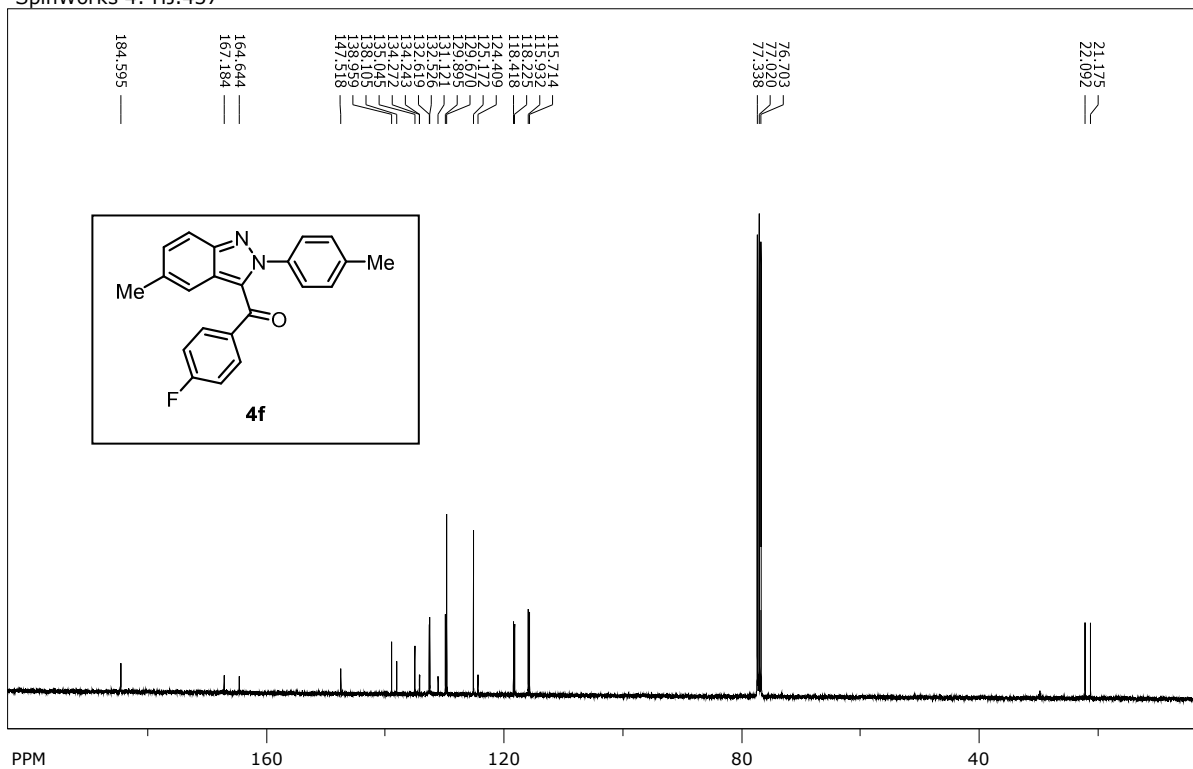
SpinWorks 4: HJ.433



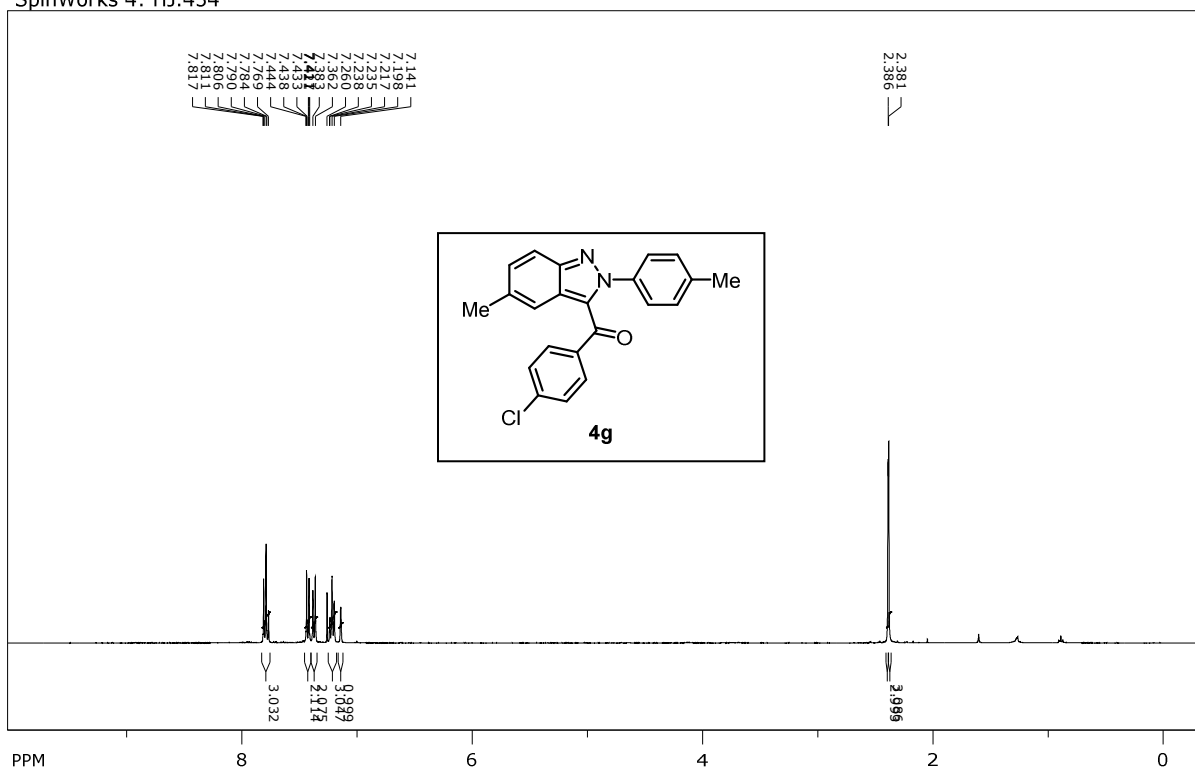
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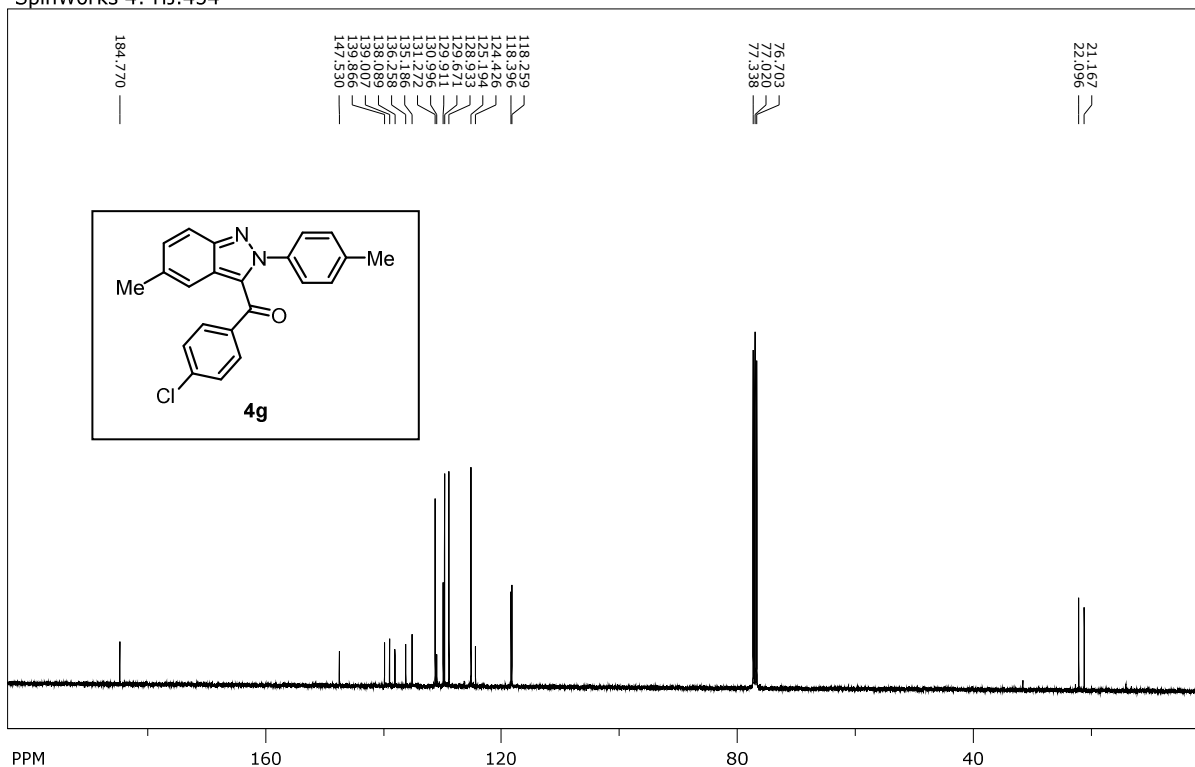
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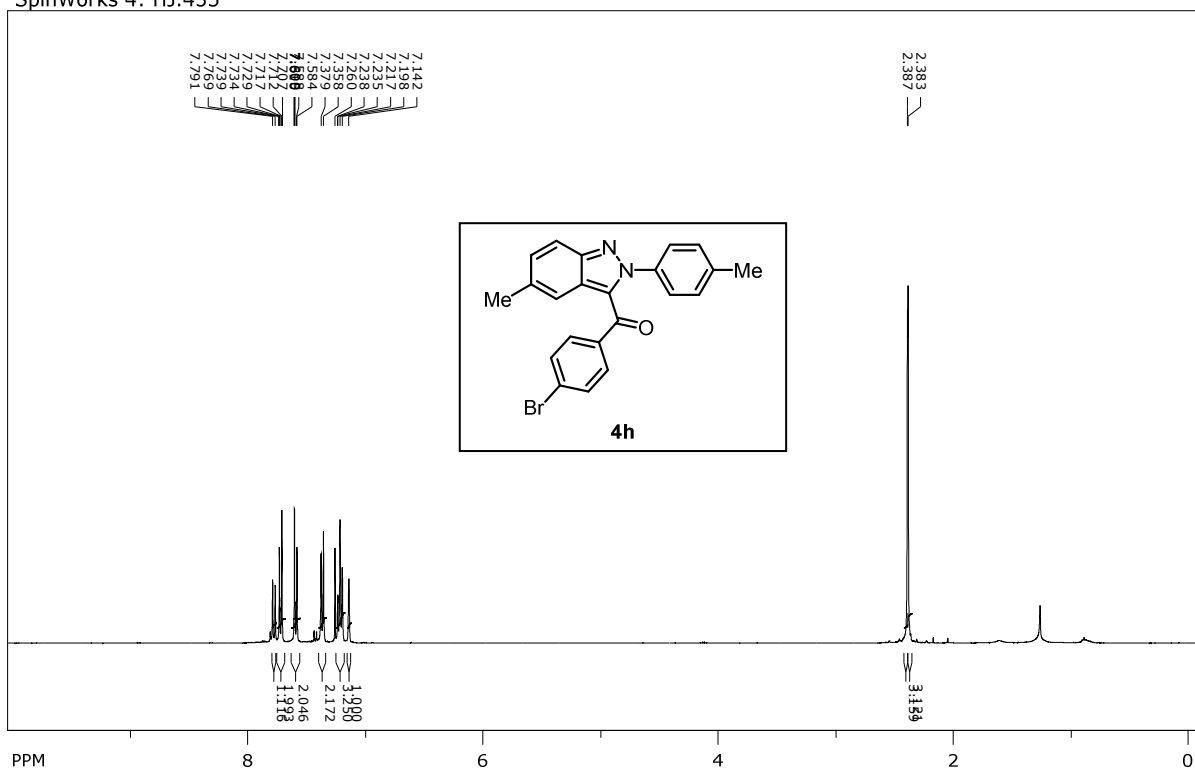
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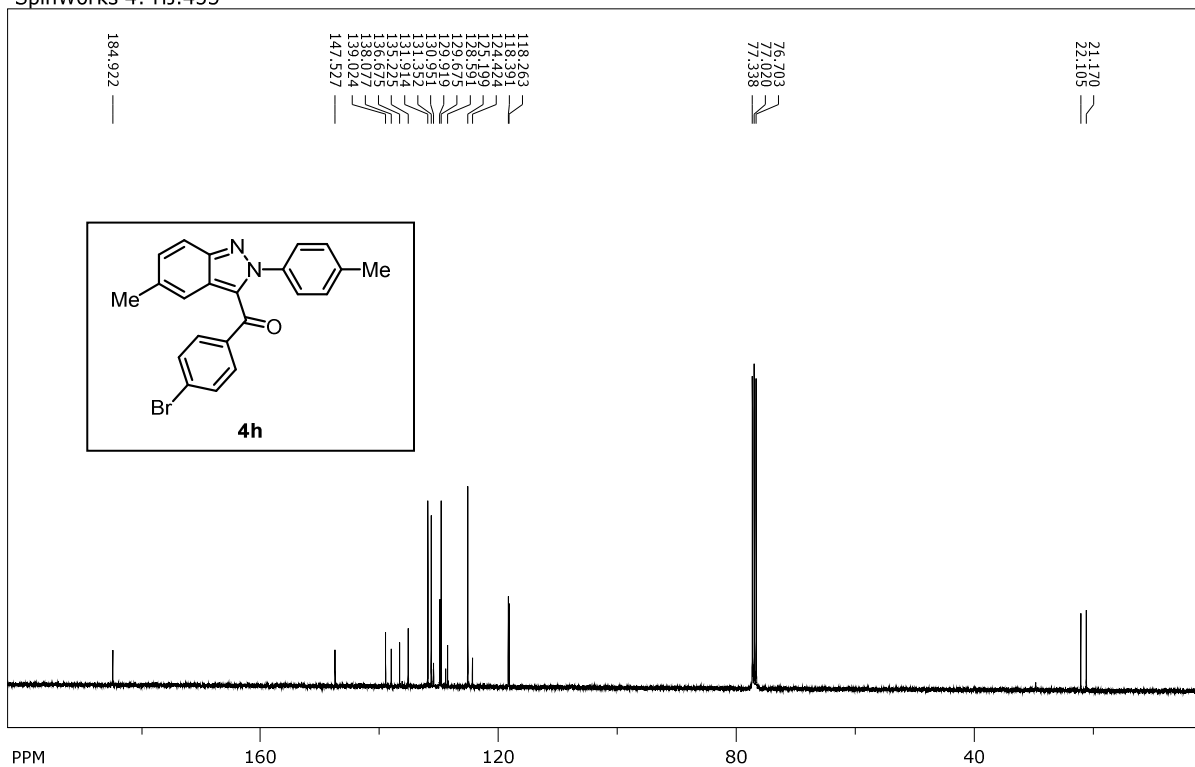
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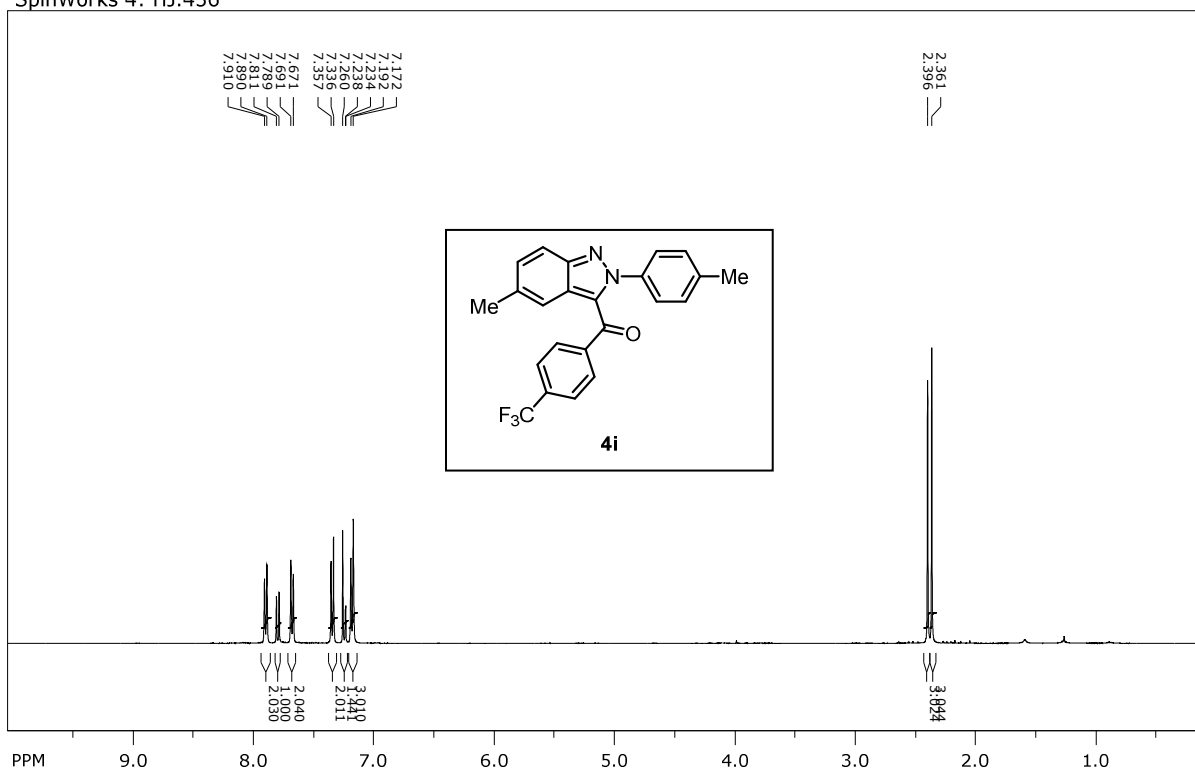
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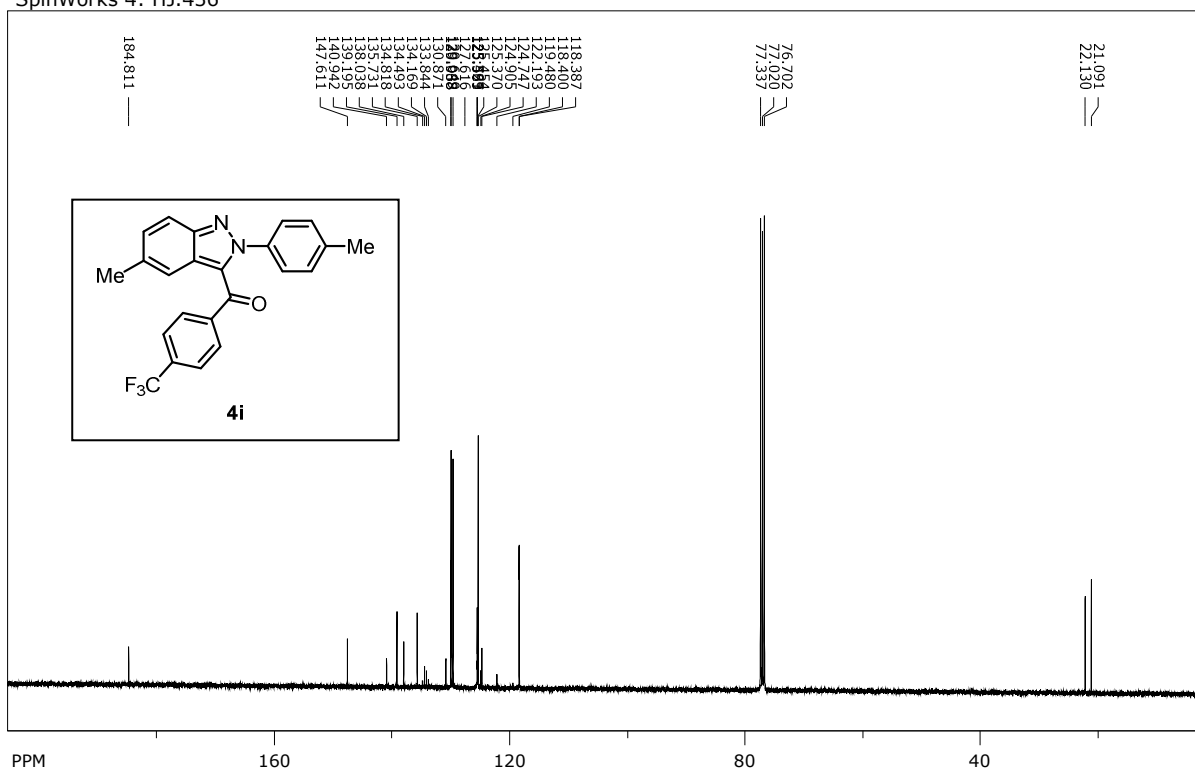
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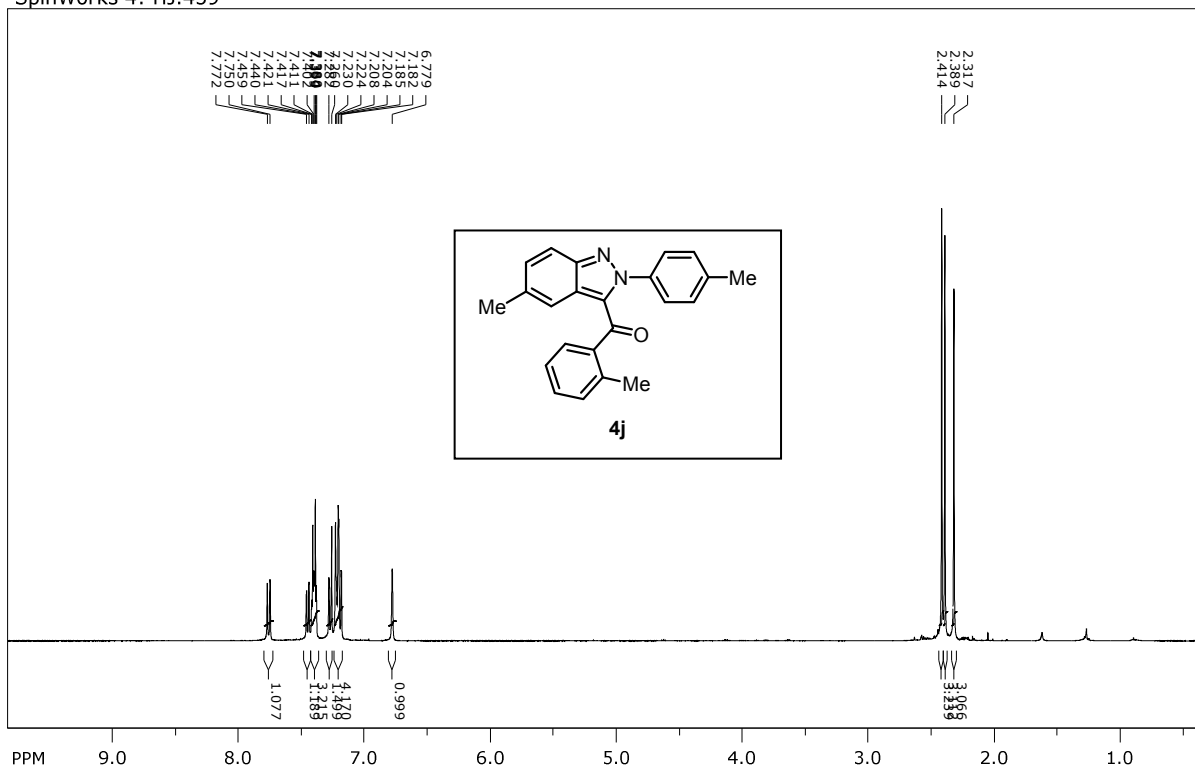
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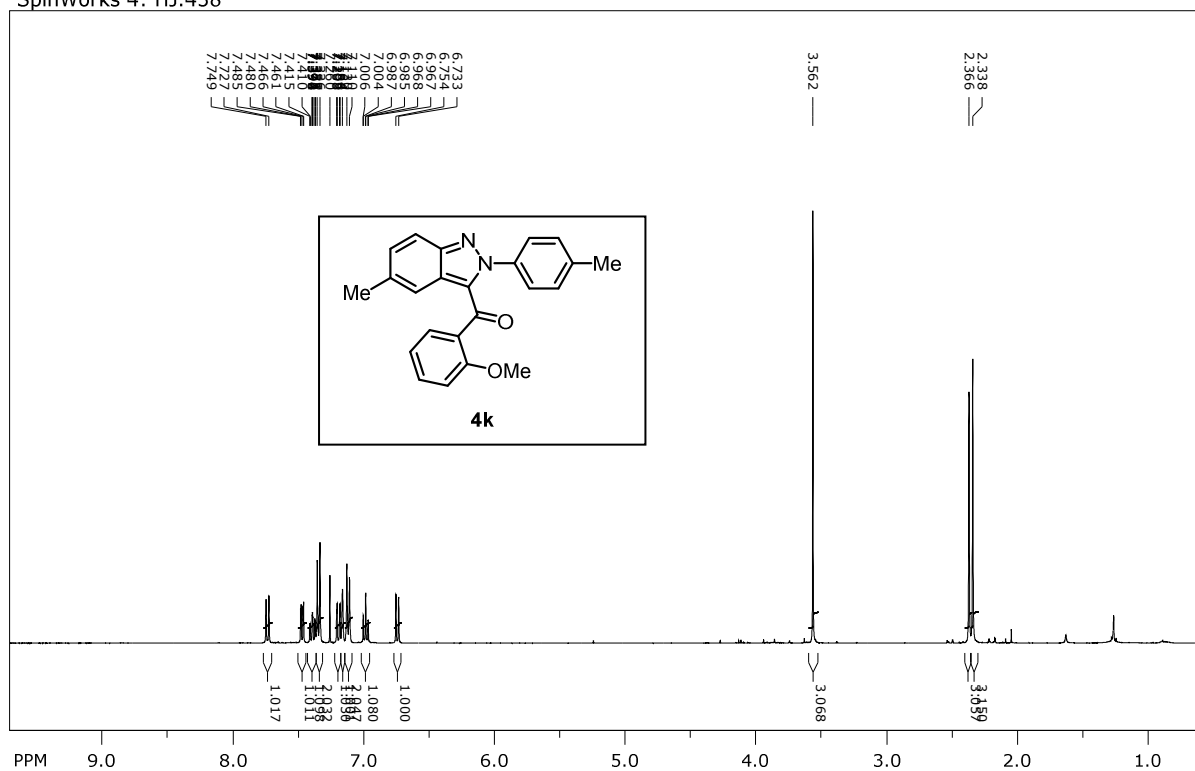
SpinWorks 4: HJ.436



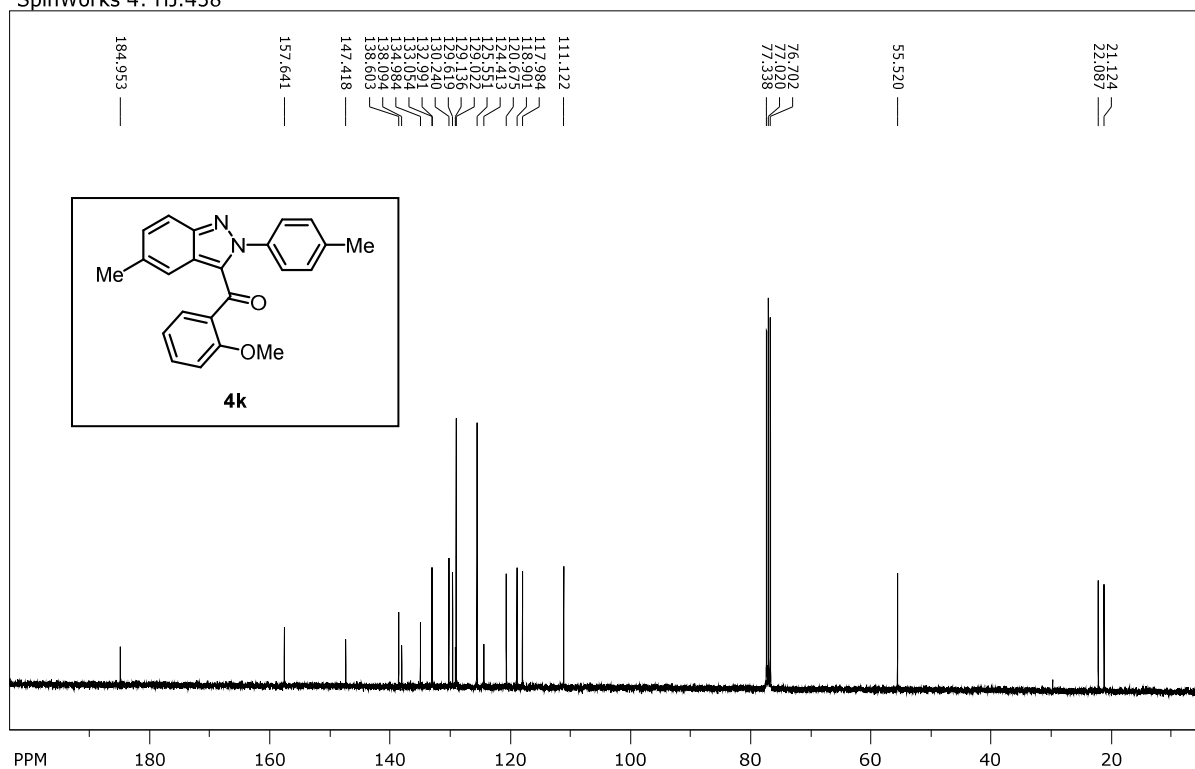
SpinWorks 4: HJ.439



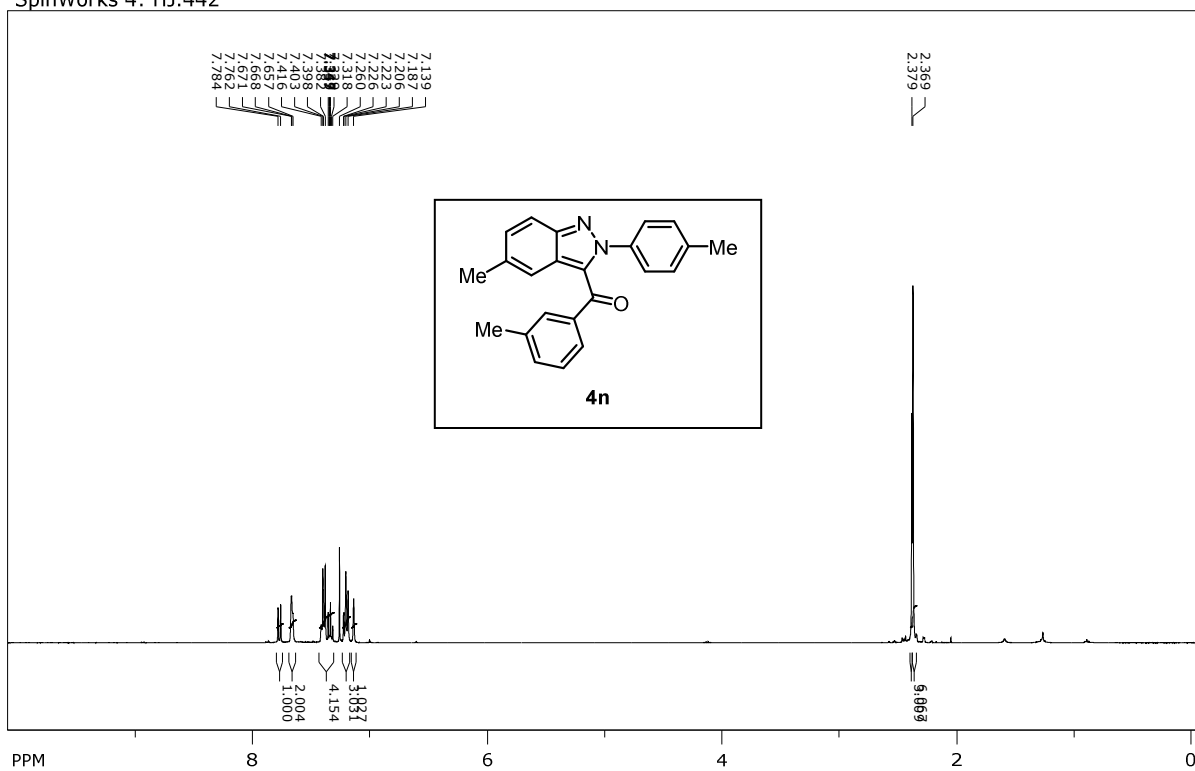
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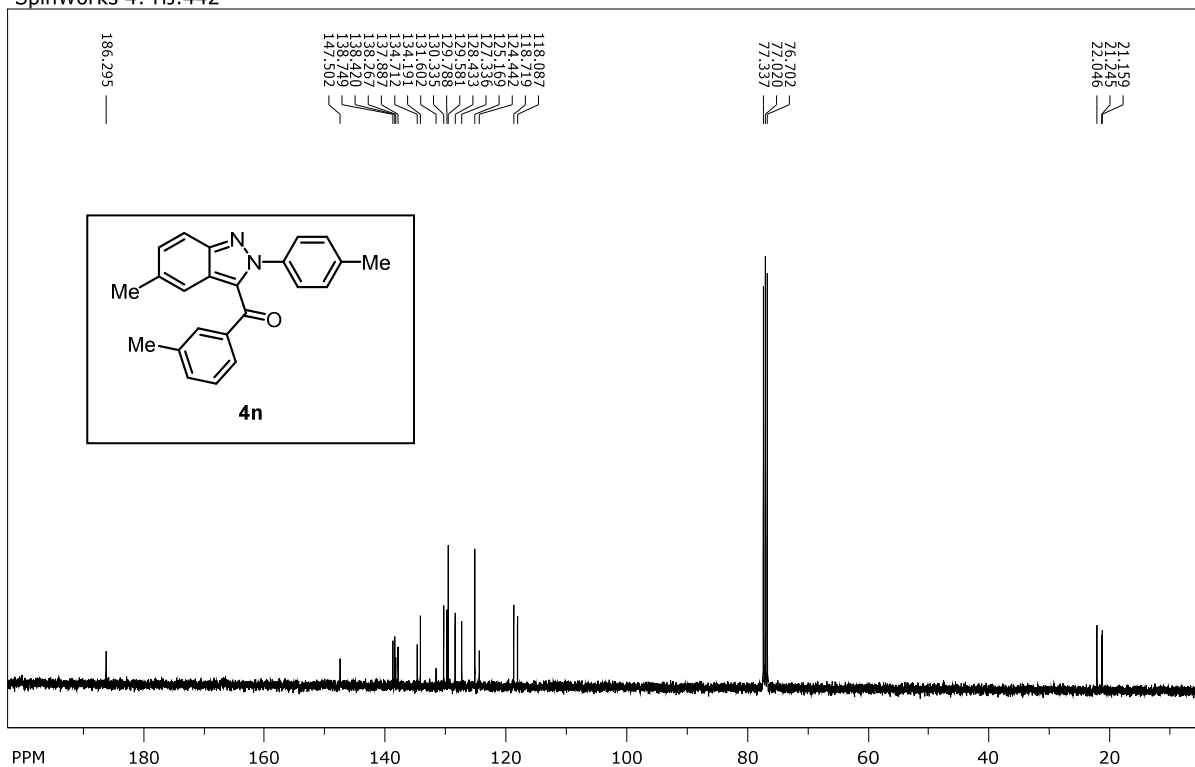
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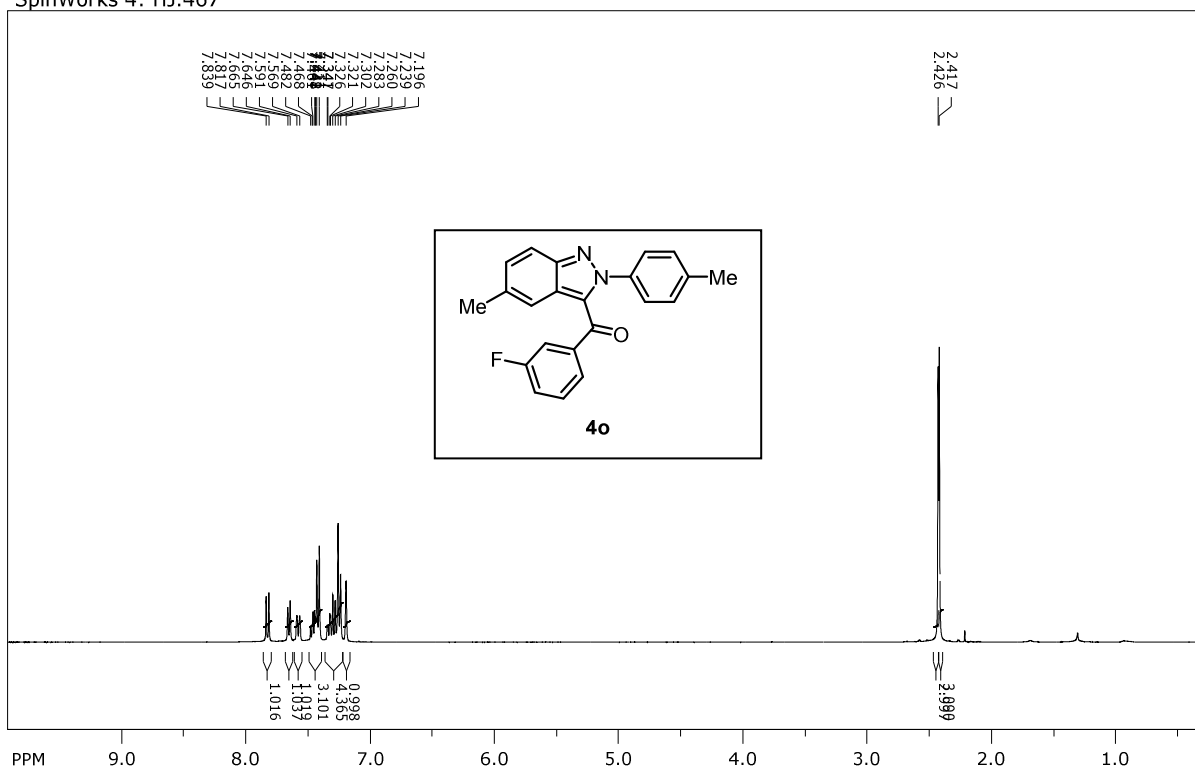
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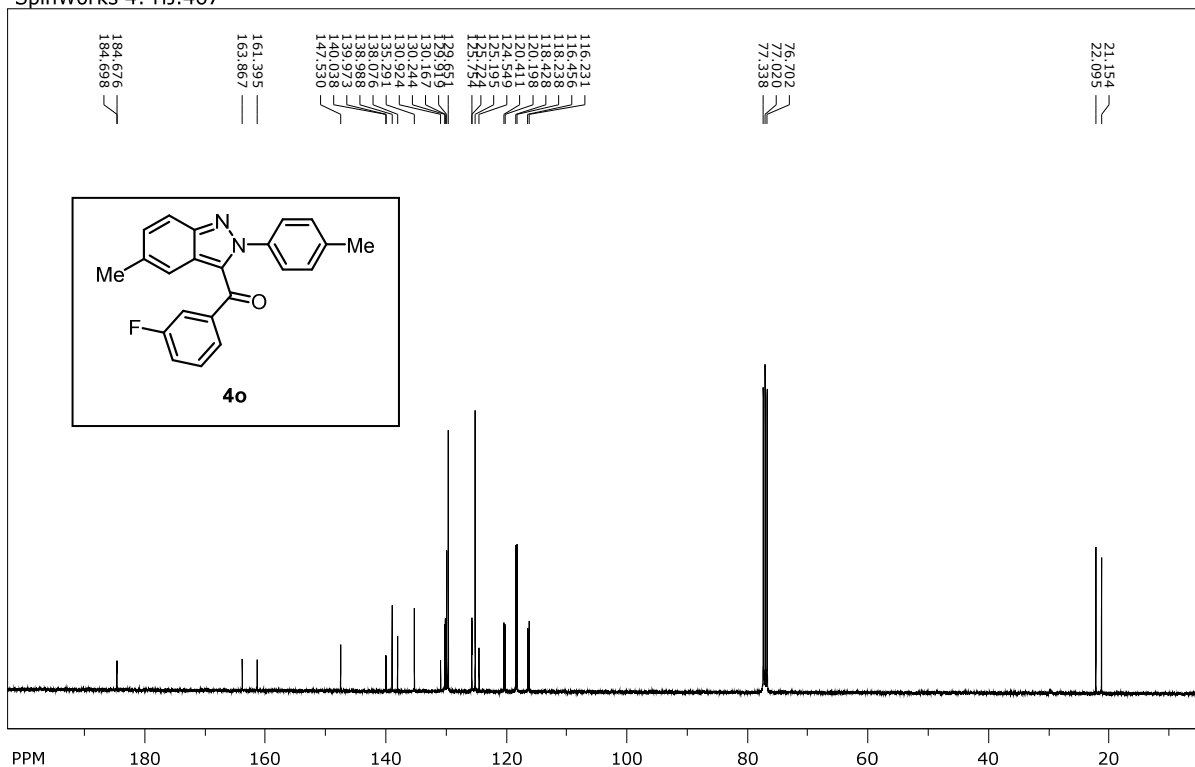
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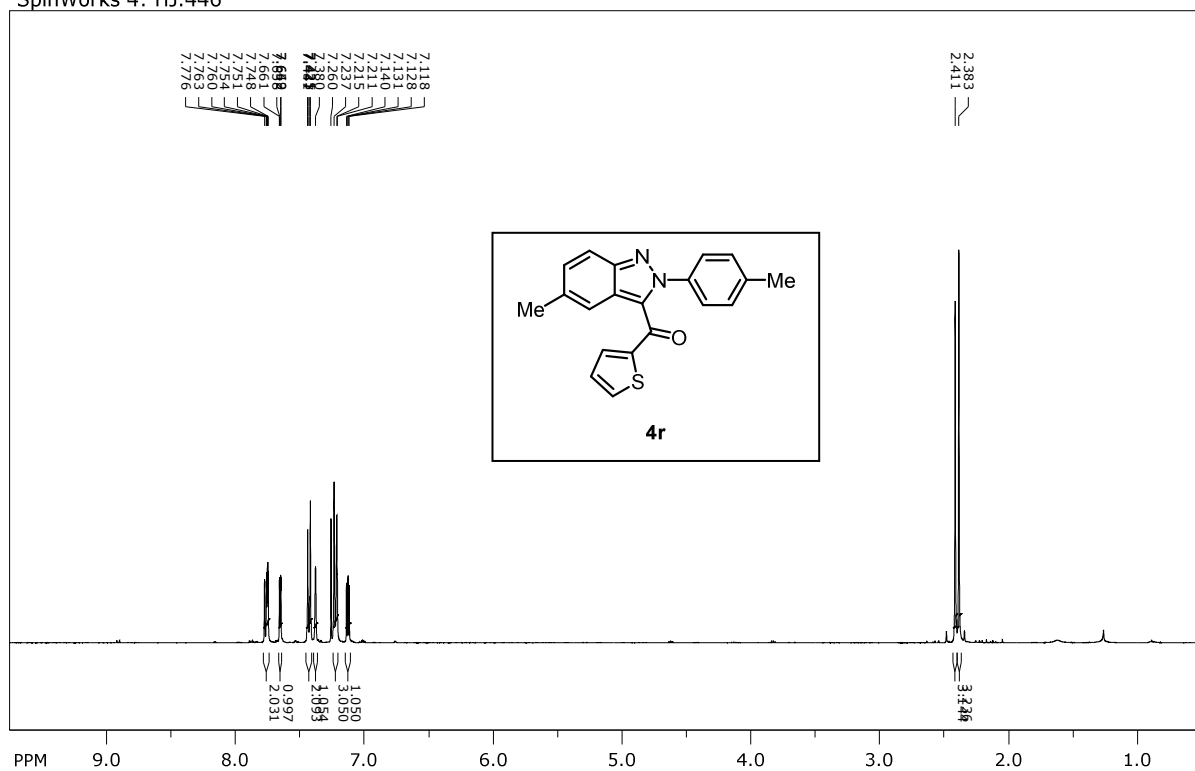
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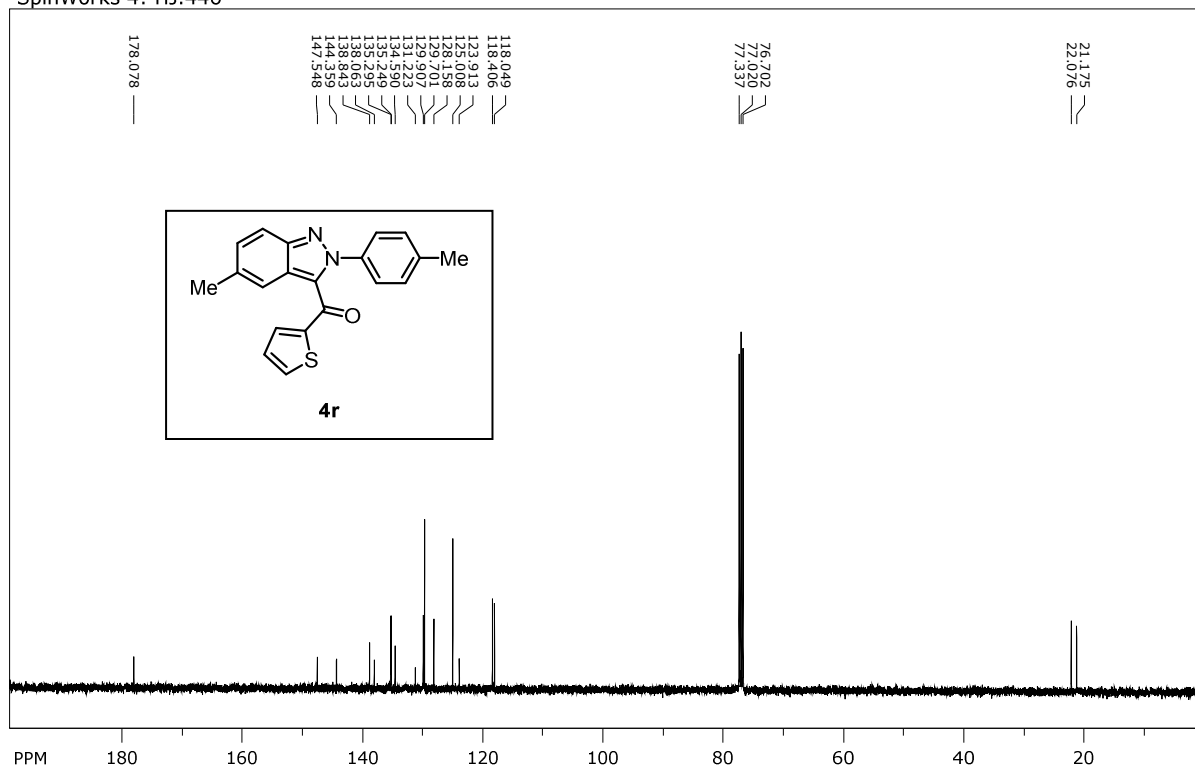
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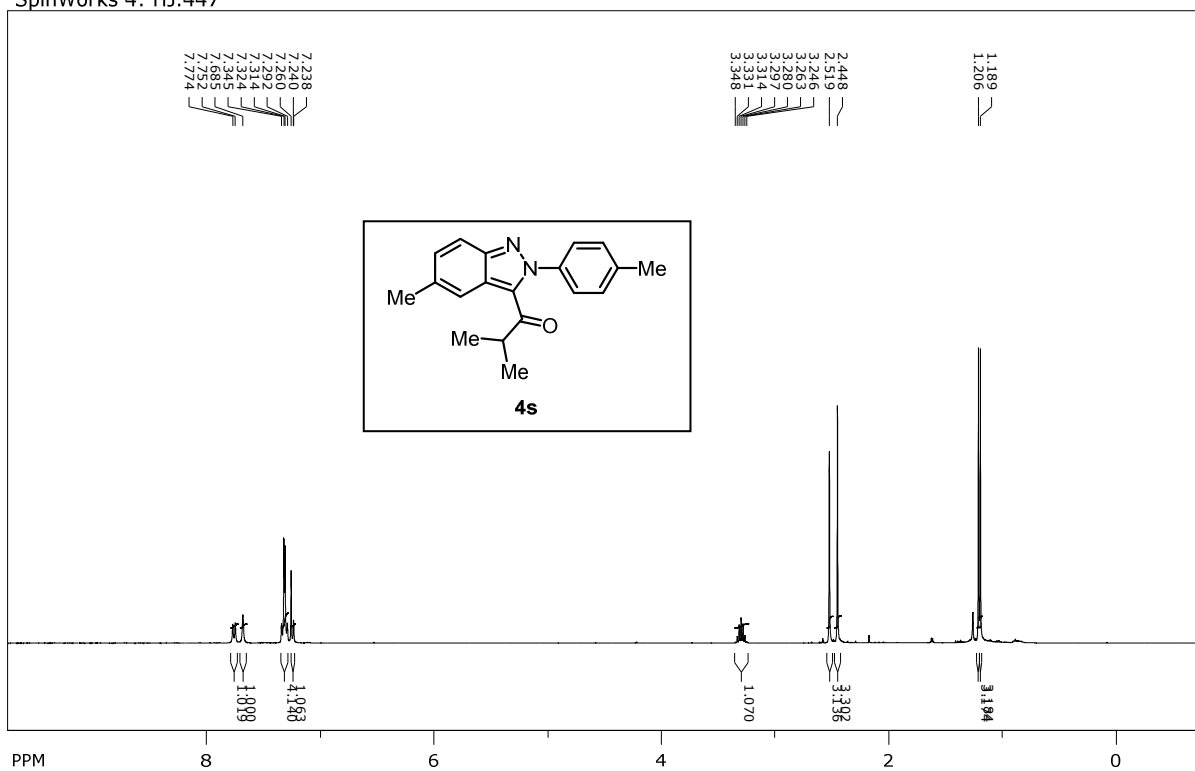
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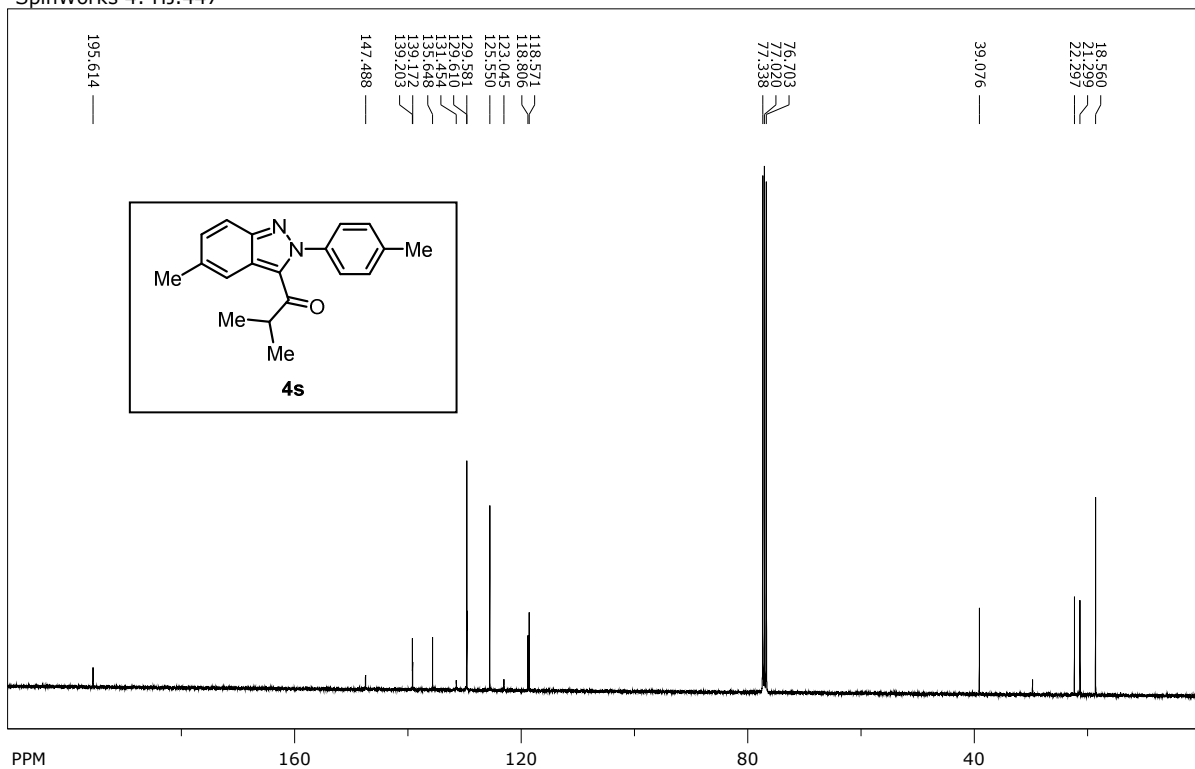
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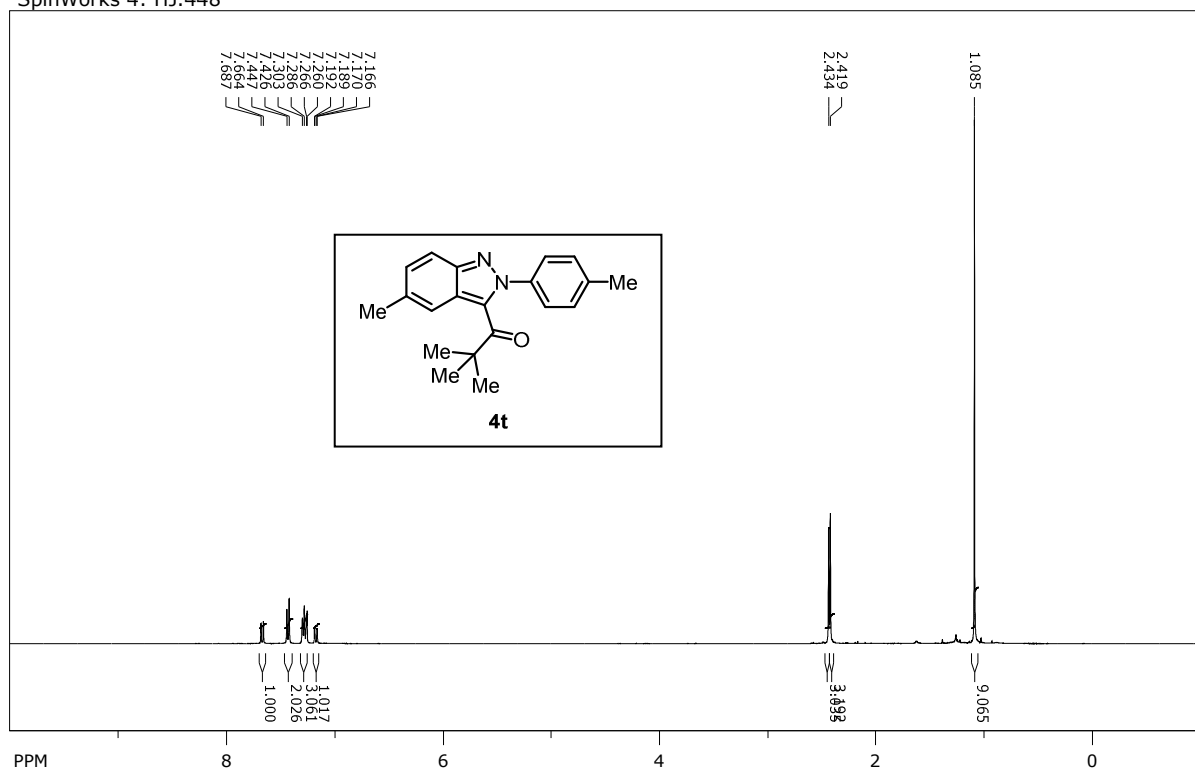
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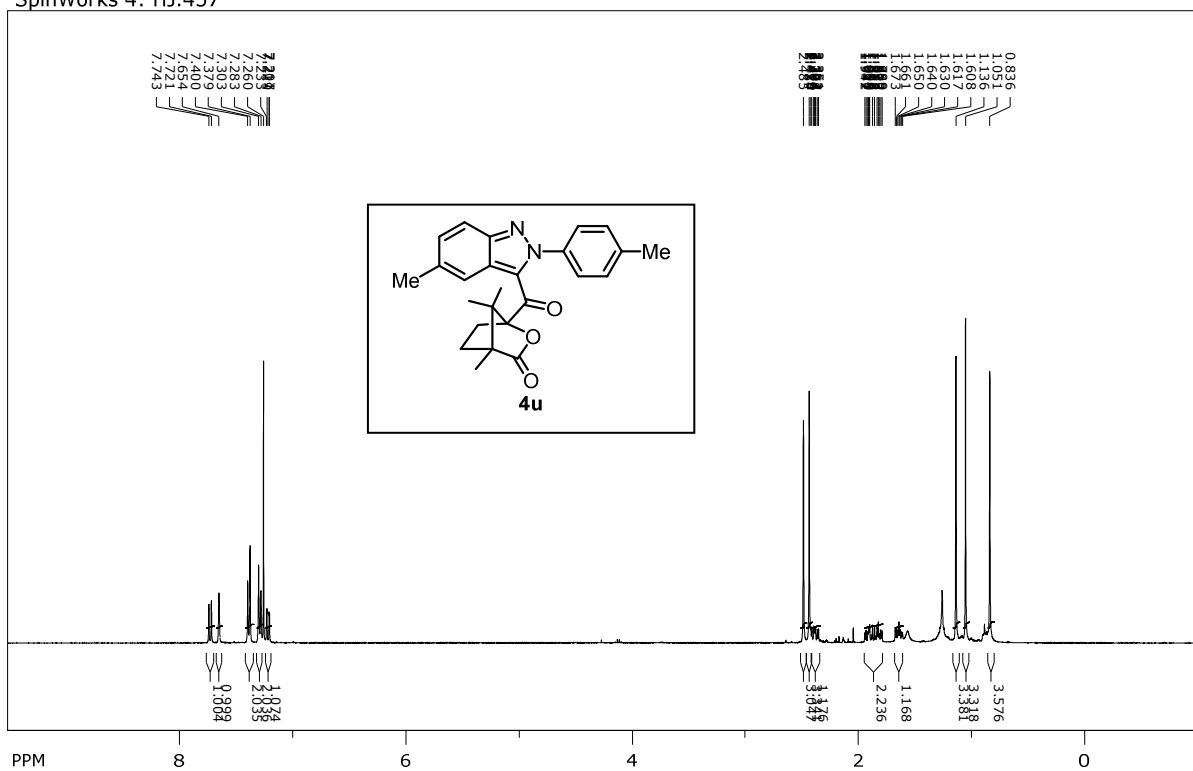
SpinWorks 4: HJ.447



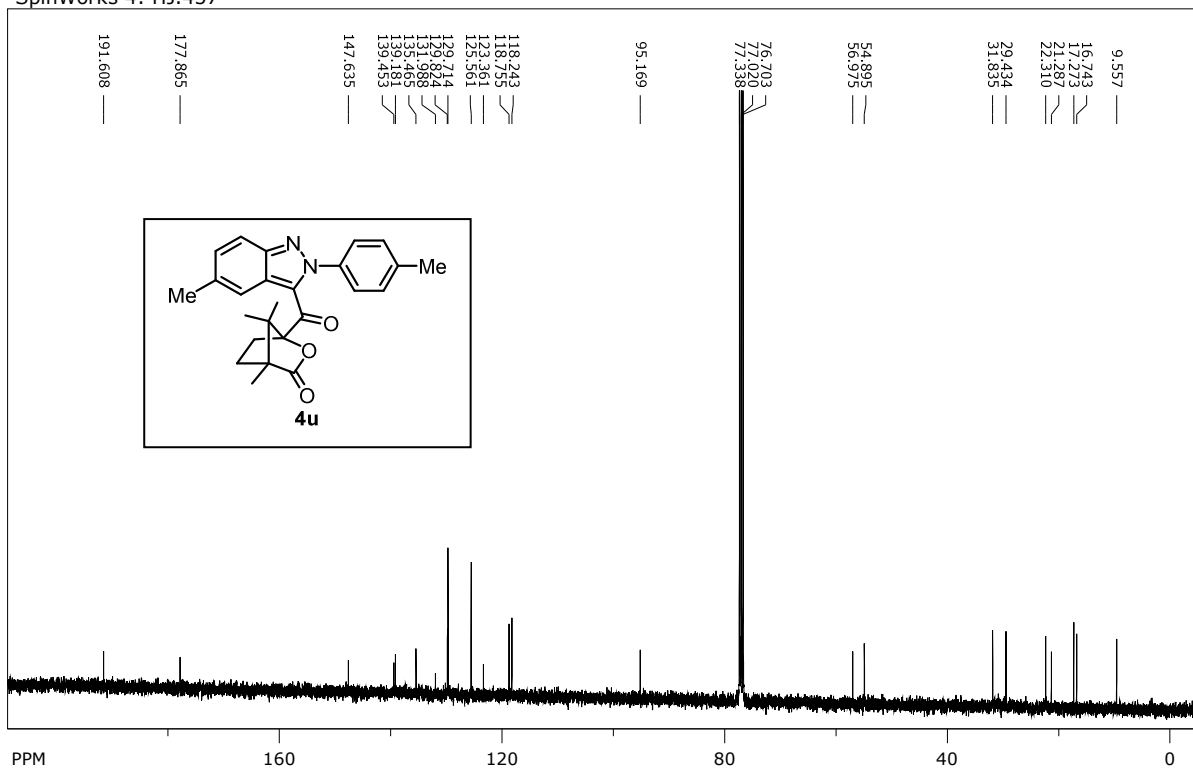
SpinWorks 4: HJ.448



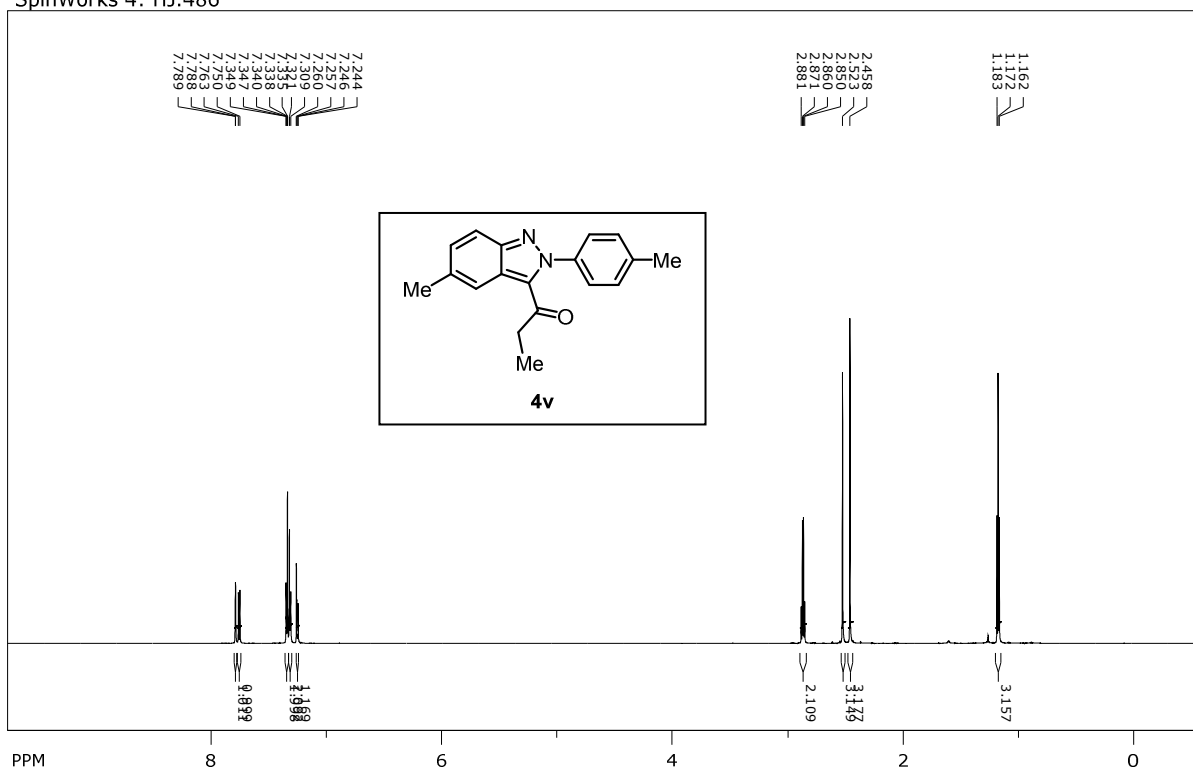
SpinWorks 4: HJ.457



SpinWorks 4: HJ.457



SpinWorks 4: HJ.486



SpinWorks 4: HJ.486

