

Supporting Information: Syntheses of 3,3-disubstituted Dihydrobenzofurans, Indolines, Indolinones and Isochromanes by Palladium-Catalyzed Tandem Reaction Using $\text{Pd}(\text{PPh}_3)_2\text{Cl}_2/(\pm)\text{-BINAP}$ as a Catalytic System

Guizhou Yue ^{1,*},[†], Sicheng Li ^{1,†}, Dan Jiang ¹, Gang Ding ², Juhua Feng ¹, Huabao Chen ², Chunping Yang ², Zhongqiong Yin ³, Xu Song ³, Xiaoxia Liang ³, Li Zhang ¹, Xianxiang Wang ¹ and Cuifen Lu ^{4,*}

¹ College of Science, Sichuan Agricultural University, Ya'an, Sichuan, 625014, China; lisicheng@stu.sicau.edu.cn (S.L.); jiangdan18@stu.sicau.edu.cn (D.J.); jhfeng@sicau.edu.cn (J.F.); zhangli@sicau.edu.cn (L.Z.); wangxx@sicau.edu.cn (X.W.)

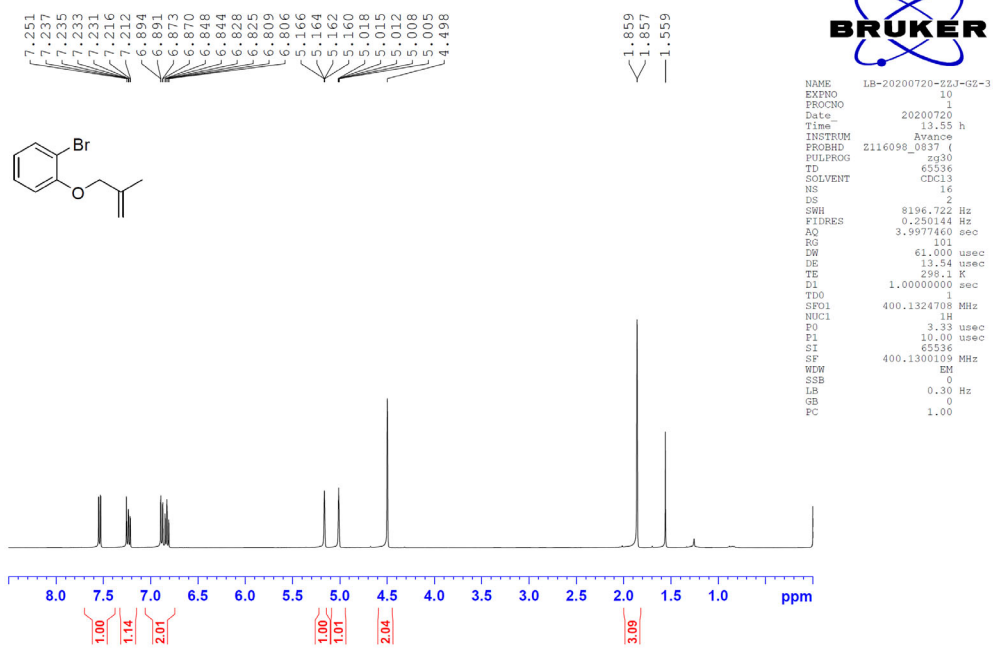
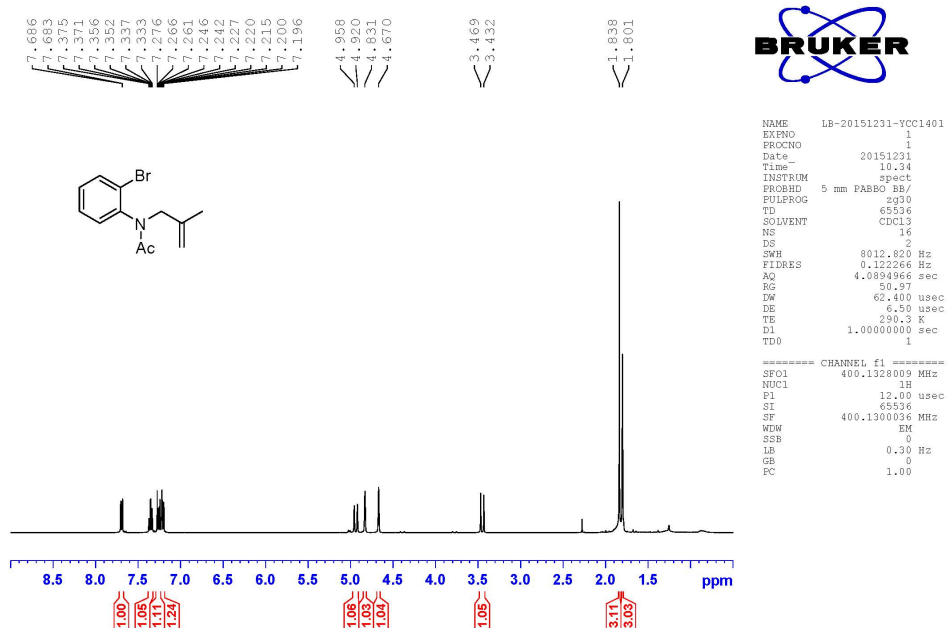
² College of Agricultural Sciences, Sichuan Agricultural University, Chengdu, Sichuan, 611130, China; cndyjhs@stu.sicau.edu.cn (G.D.); chenuabao@sicau.edu.cn (H.C.); yangchunping@sicau.edu.cn (C.Y.)

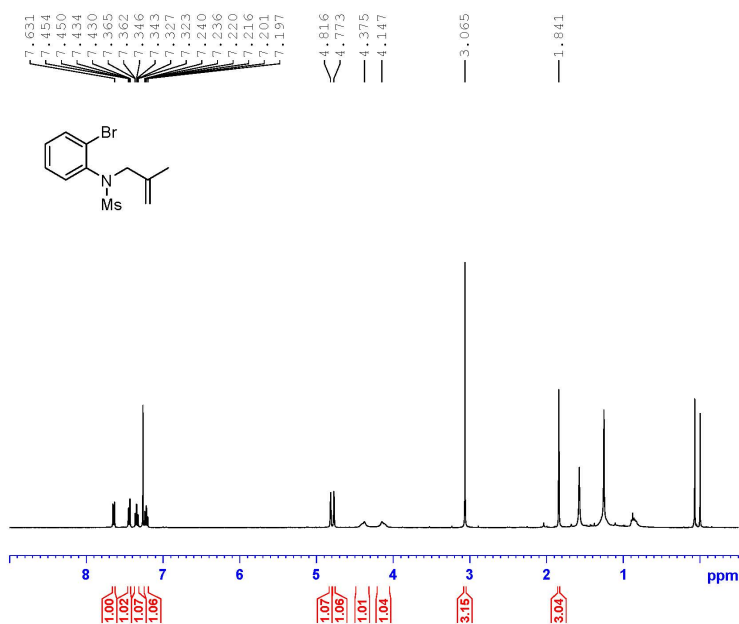
³ College of Veterinary Medicine, Sichuan Agricultural University, Chengdu, Sichuan, 611130, China; yinzhongq@sicau.edu.cn (Z.Y.); songx@sicau.edu.cn (X.S.); liangxiaoxia@sicau.edu.cn (X.L.)

⁴ Hubei Collaborative Innovation Center for Advanced Organochemical Materials & Ministry-of-Education Key Laboratory for the Synthesis and Application of Organic Functional Molecules, Hubei University, Wuhan 430062, China

* Correspondence: yueguizhou@sicau.edu.cn (G.Y.); lucf@hubu.edu.cn (C.L.); Tel.: +86-0835-2886189 (G.Y.); +86-027-50865322 (C.L.)

[†] The authors contributed equally to this work

¹H NMR Spectra for Compound 1a¹H NMR Spectra for Compound 1b

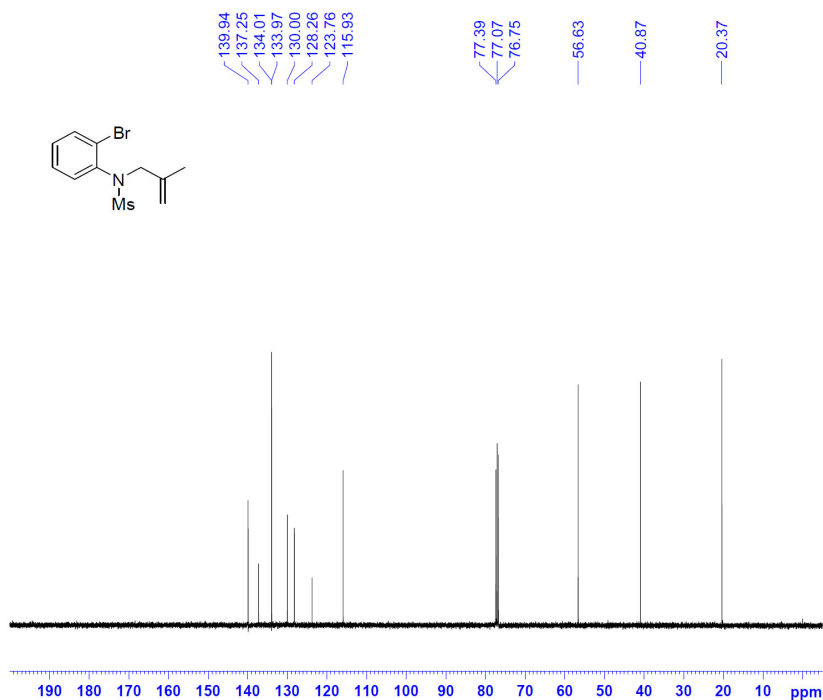
^1H and ^{13}C NMR Spectra for Compound 1c

```

NAME      Ms-yuanliao
EXPNO     1
PROCNO    1
Date_     20151202
Time      23.09
INSTRUM   spect
PROBHD    5 mm PABBO BB/
PULPROG   zgpg30
TD        65536
SOLVENT   CDCl3
NS         16
DS         0
SWH        8223.685 Hz
FIDRES     0.125483 Hz
AQ         3.9846397 sec
RG         121.24
DW         60.800 usec
DE         6.90 usec
TE         0.0 K
D1         1.00000000 sec
TD0        1
  
```

```

===== CHANNEL f1 =====
NUC1      1H
P1        9.93 usec
SI        65536
SF        400.1300095 MHz
WDW        EM
SSB        0
LB        0.30 Hz
GB         0
PC         1.00
  
```

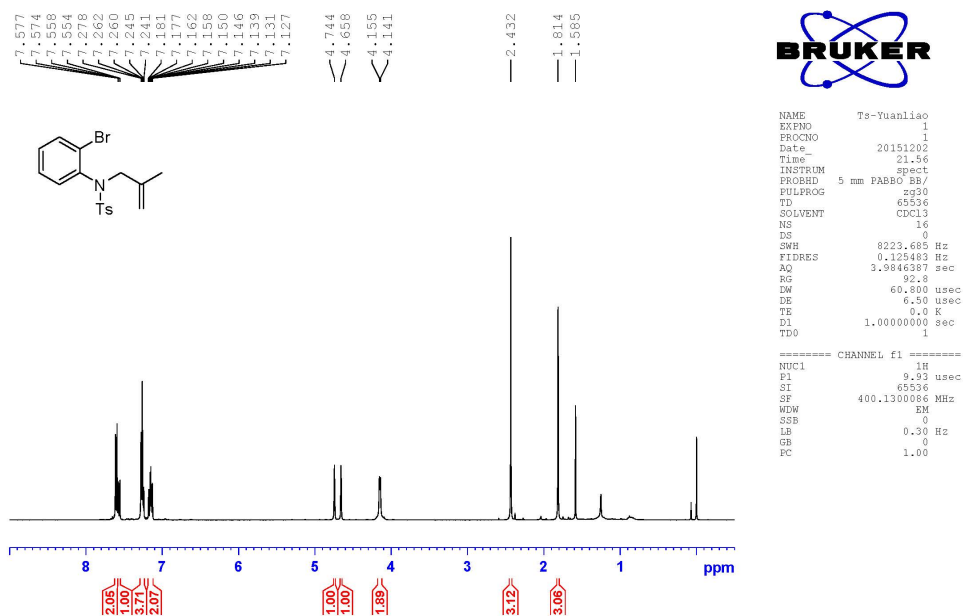
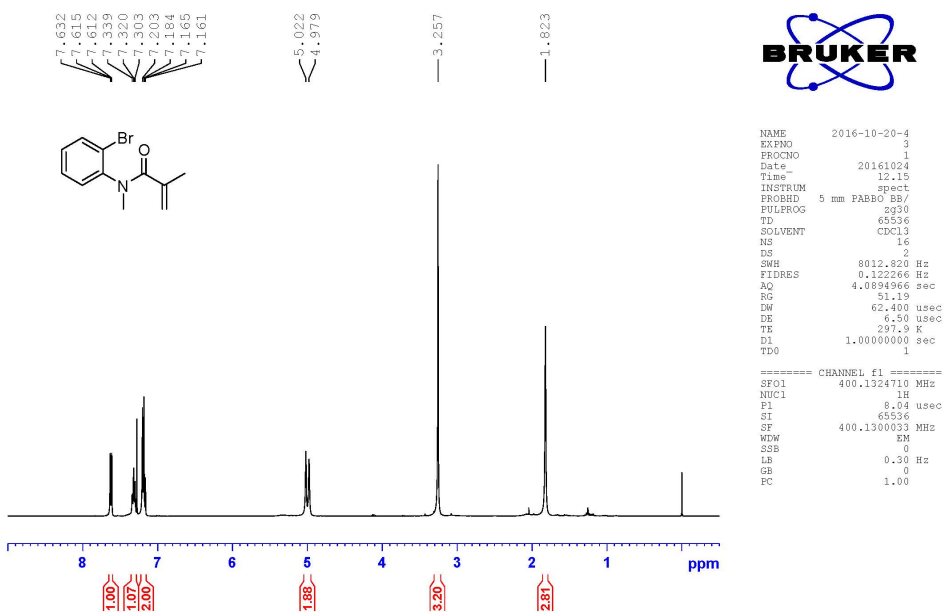


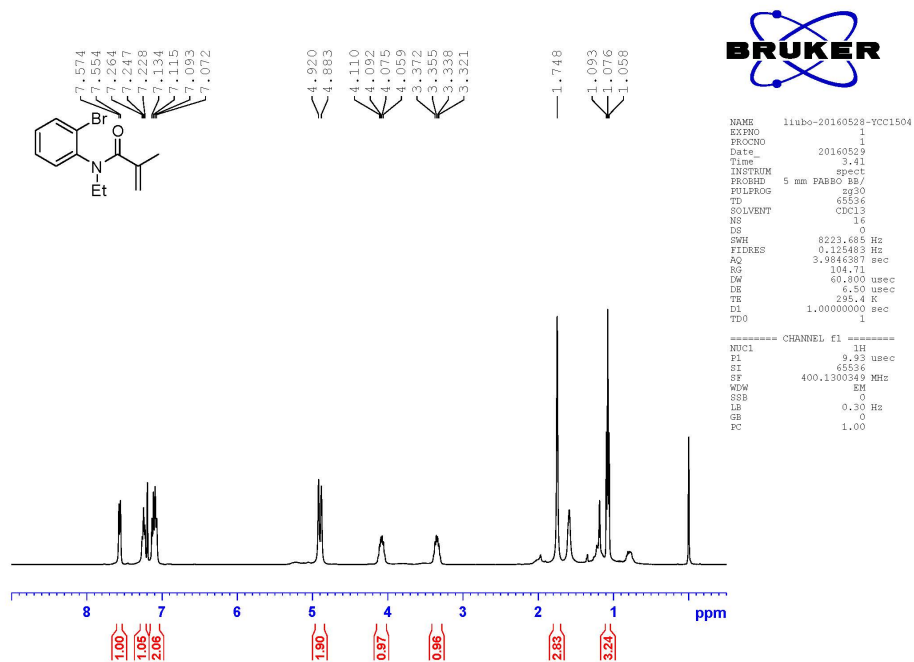
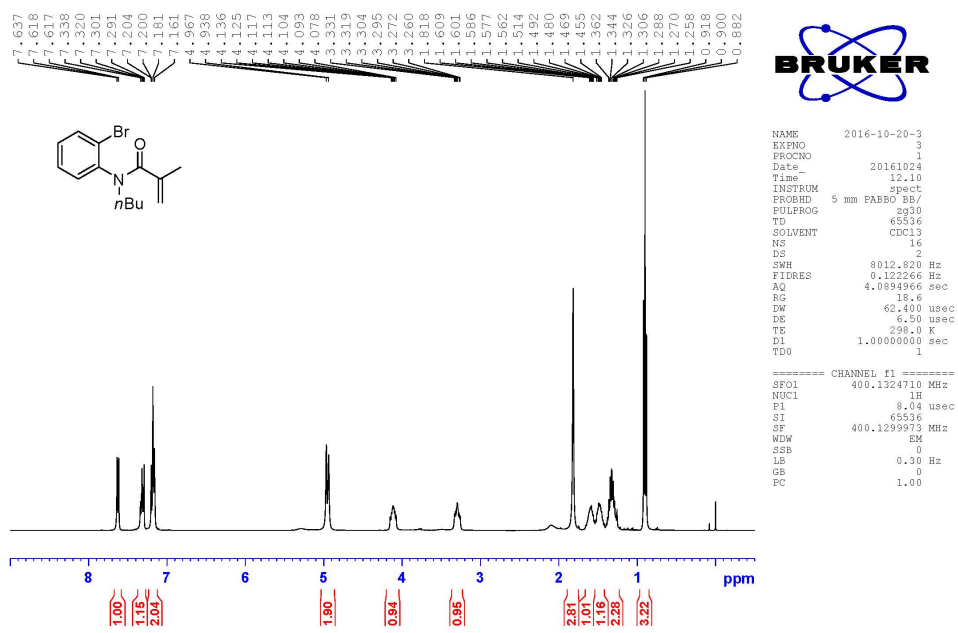
```

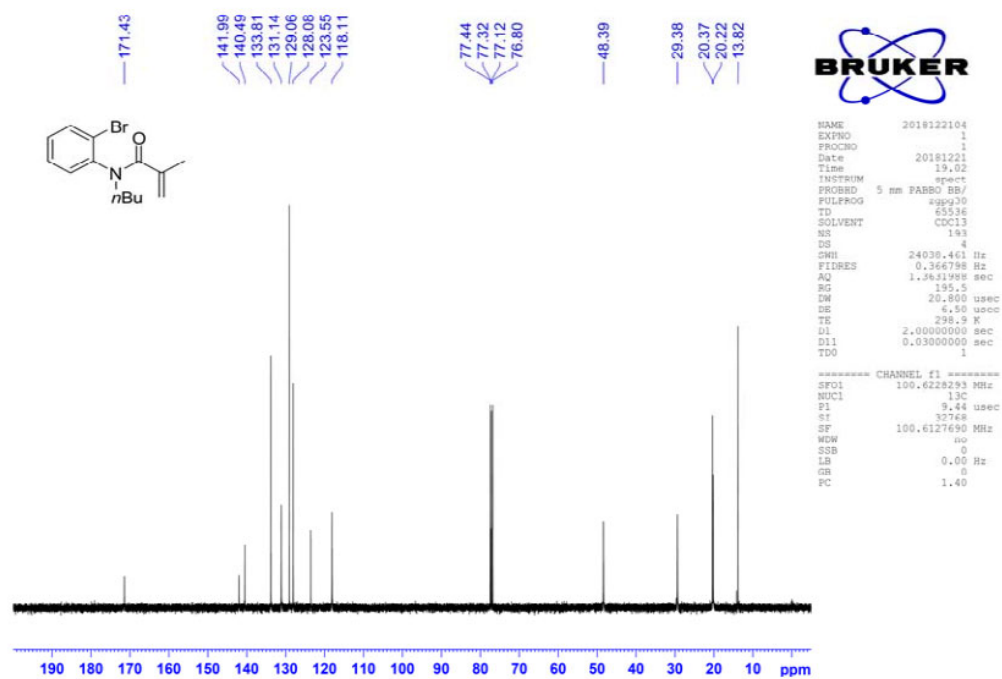
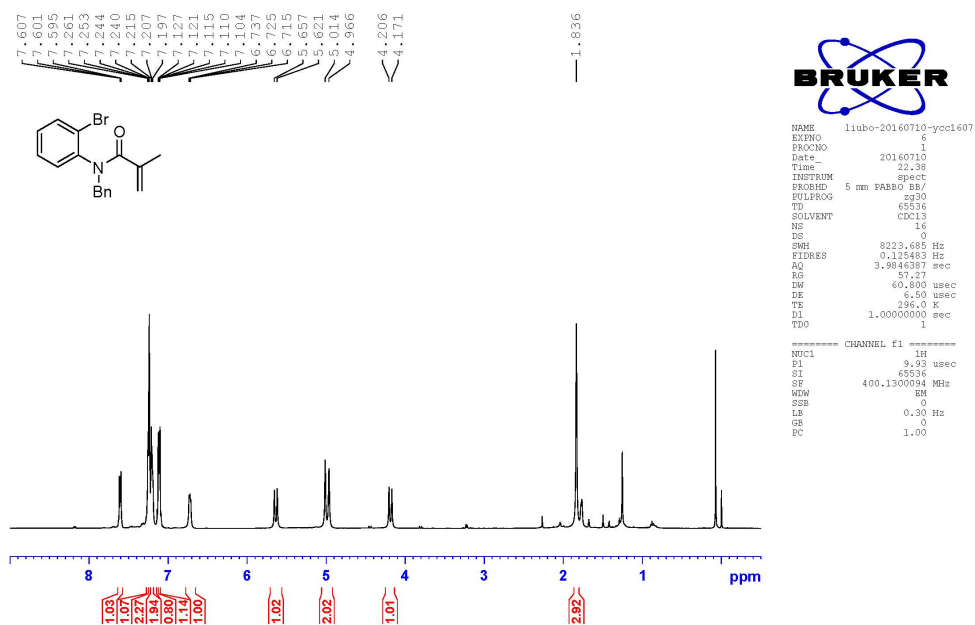
NAME      2018122103
EXPNO     1
PROCNO    1
Date_     20181221
Time      18.45
INSTRUM   spect
PROBHD    5 mm PABBO BB/
PULPROG   zgpg30
TD        65536
SOLVENT   CDCl3
NS         261
DS         4
SWH        24038.461 Hz
FIDRES     0.366798 Hz
AQ         1.3631988 sec
RG         185.5
DW         20.800 usec
DE         6.50 usec
TE         298.9 K
D1         2.00000000 sec
D11        0.03000000 sec
TD0        1
  
```

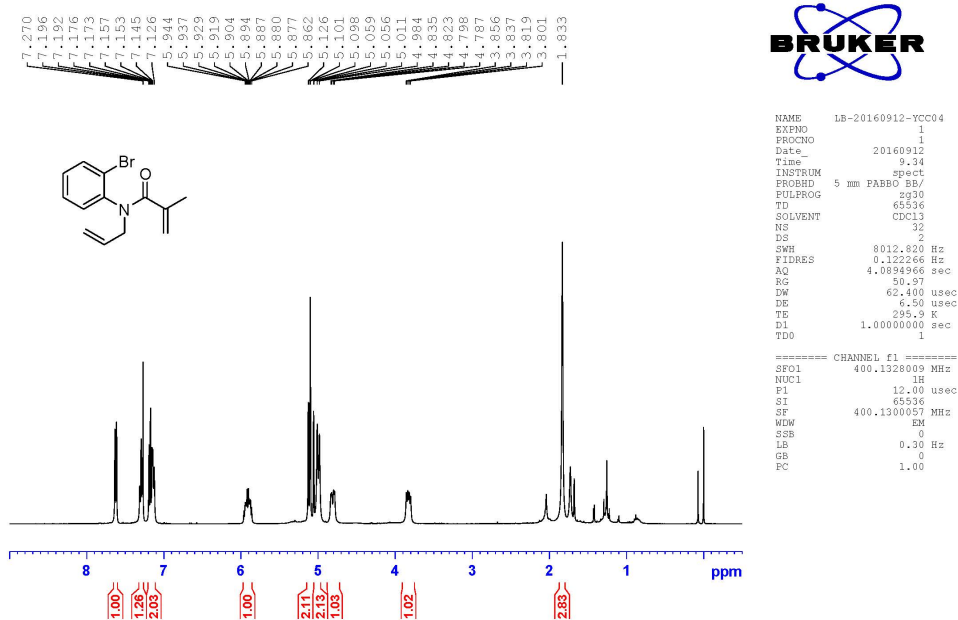
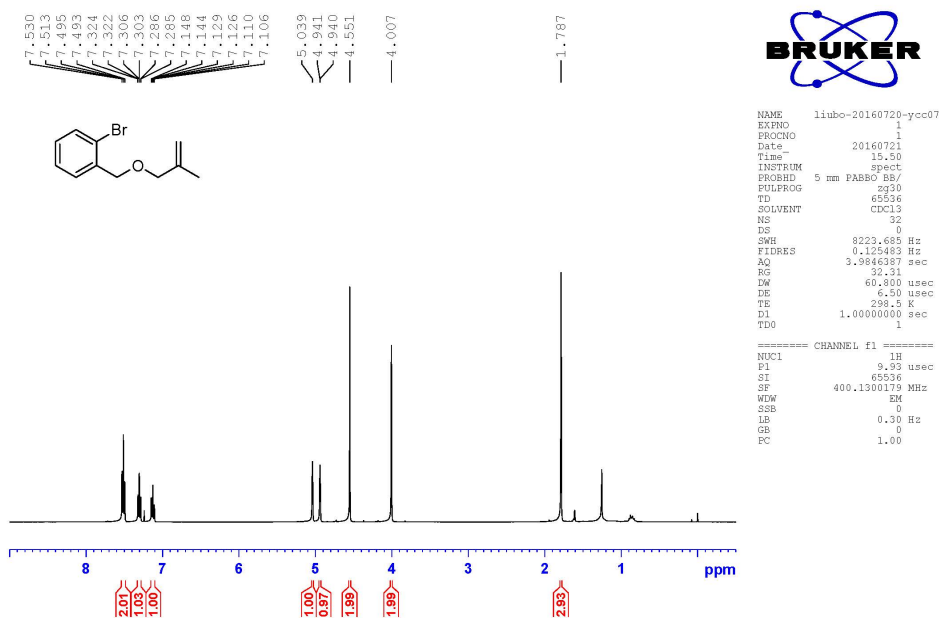
```

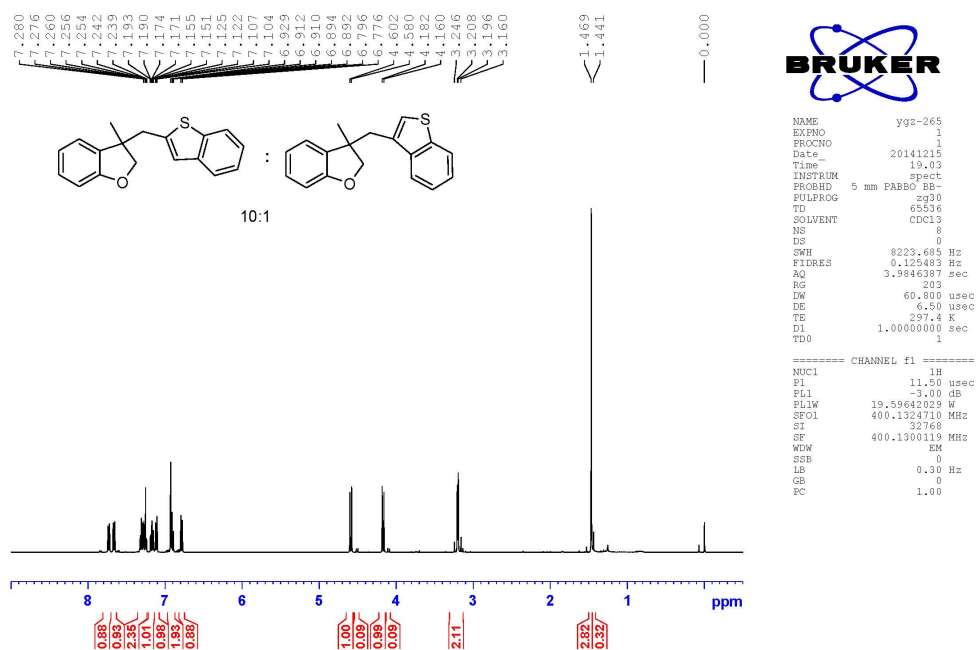
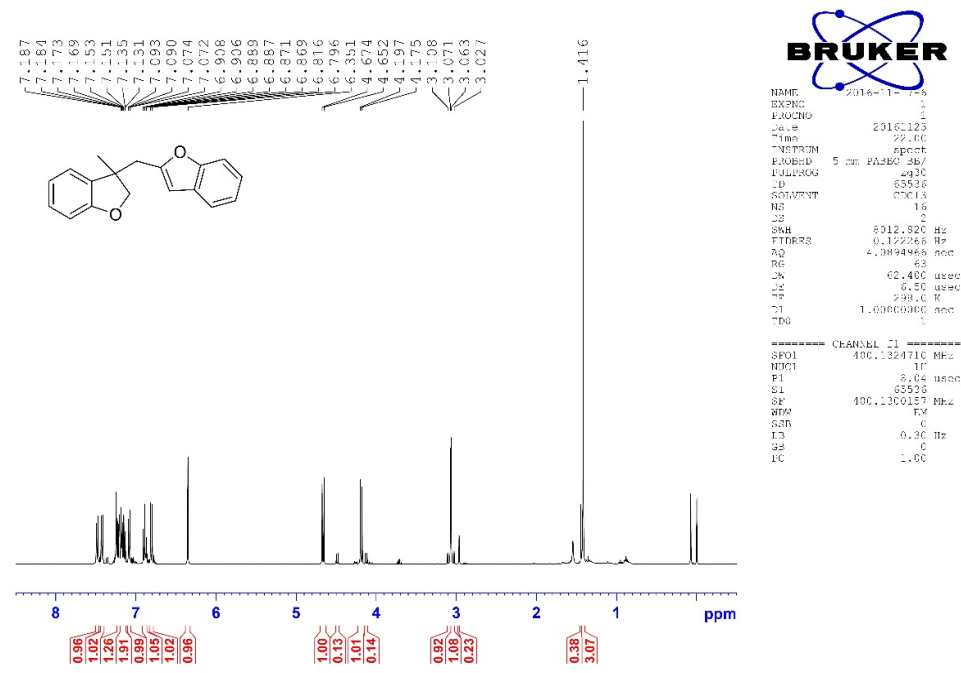
===== CHANNEL f1 =====
SFO1      100.6228293 MHz
NUC1      13C
P1        9.44 usec
SI        32768
SF        100.6127690 MHz
WDW        no
SSB        0
LB        0.00 Hz
GB         0
PC         1.40
  
```

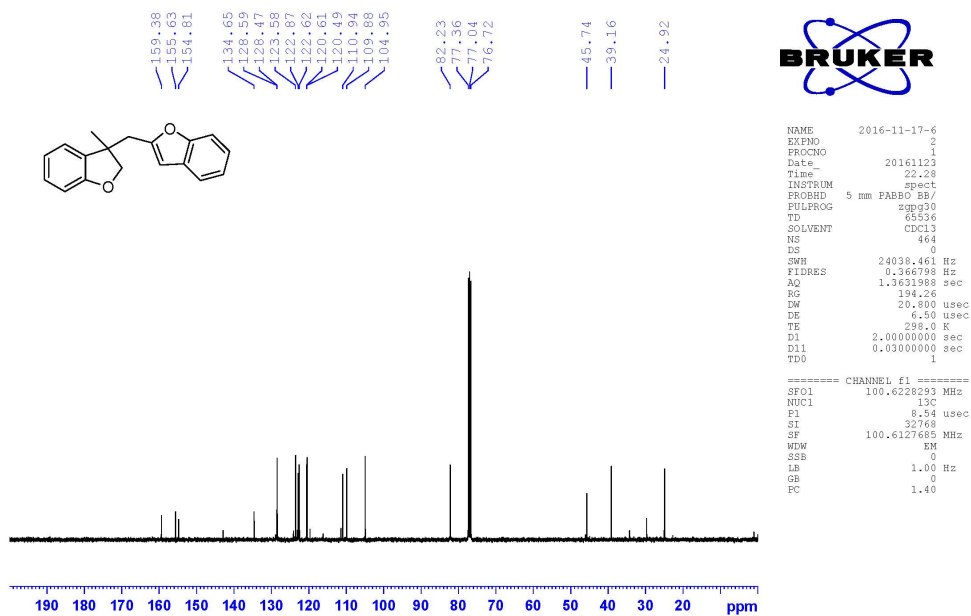
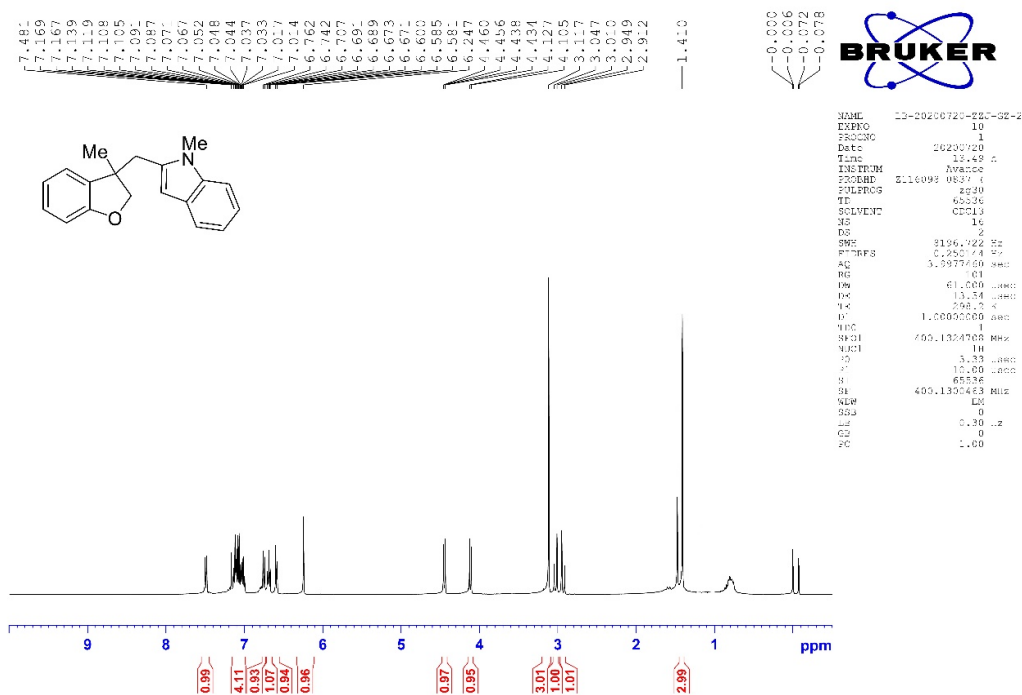
¹H NMR Spectra for Compound 1d¹H NMR Spectra for Compound 1e

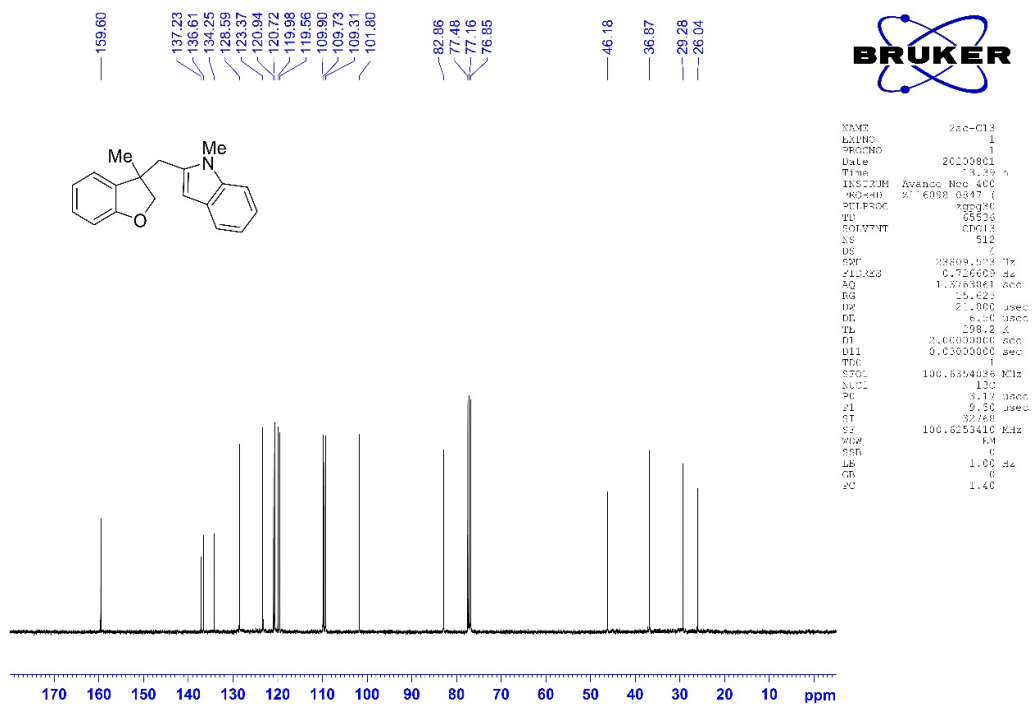
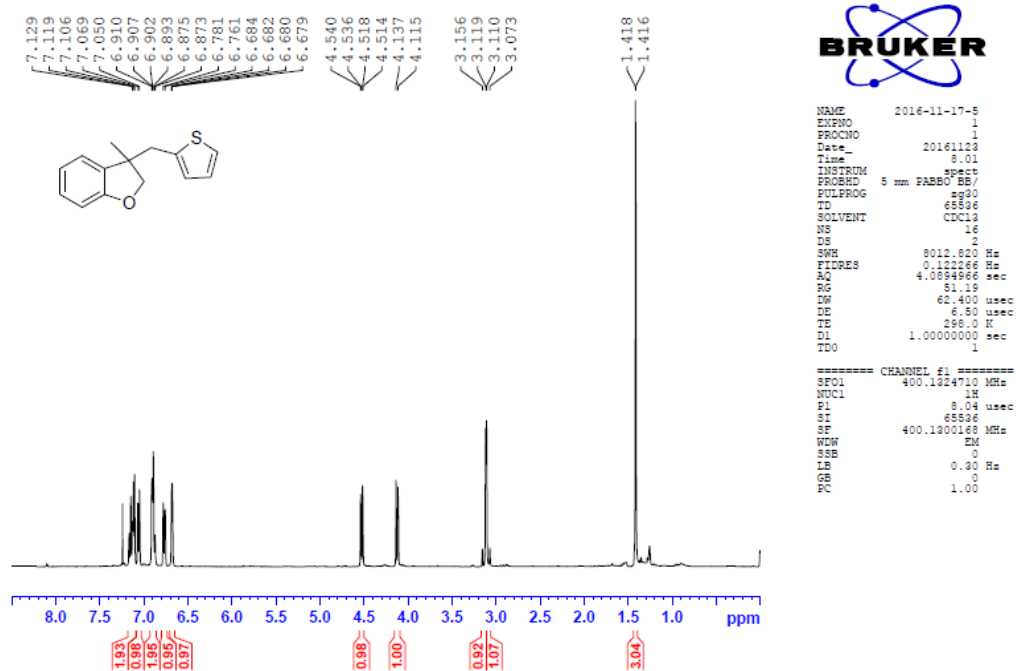
¹H NMR Spectra for Compound 1f¹H NMR Spectra for Compound 1g

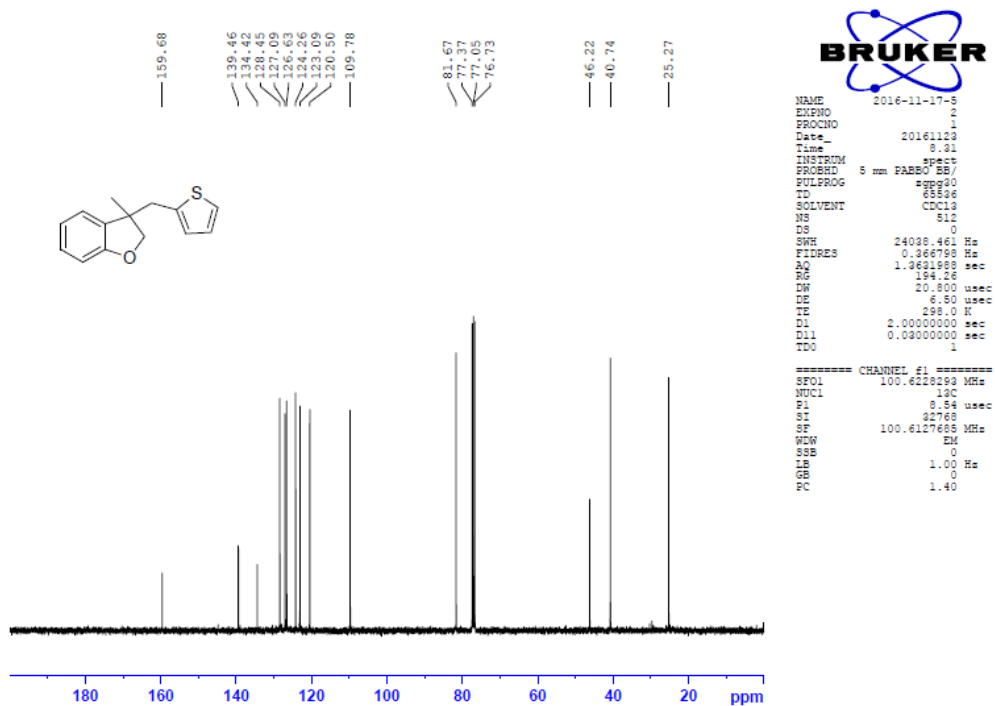
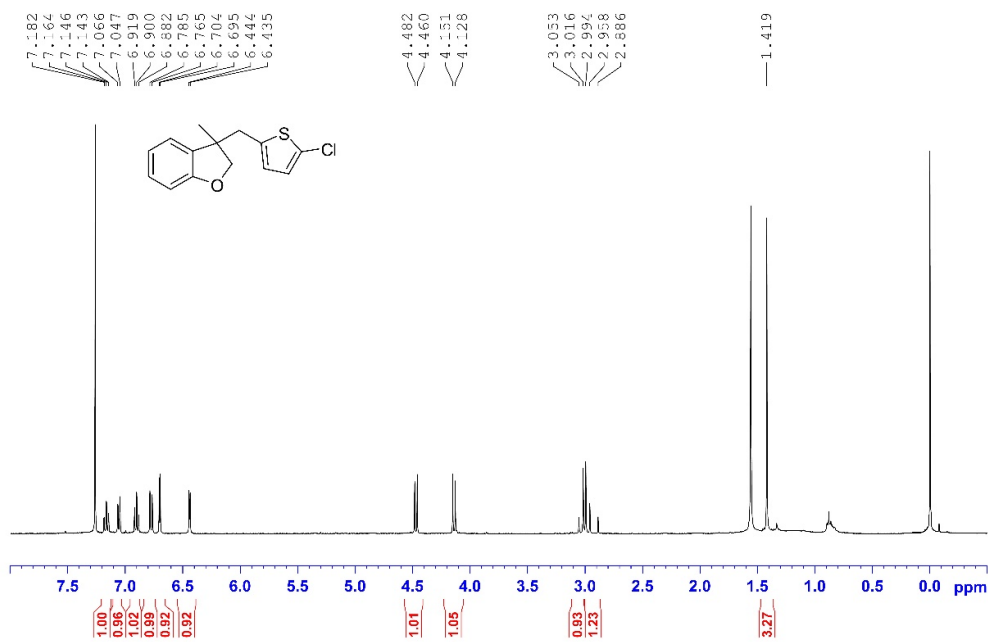
¹H NMR Spectra for Compound 1h

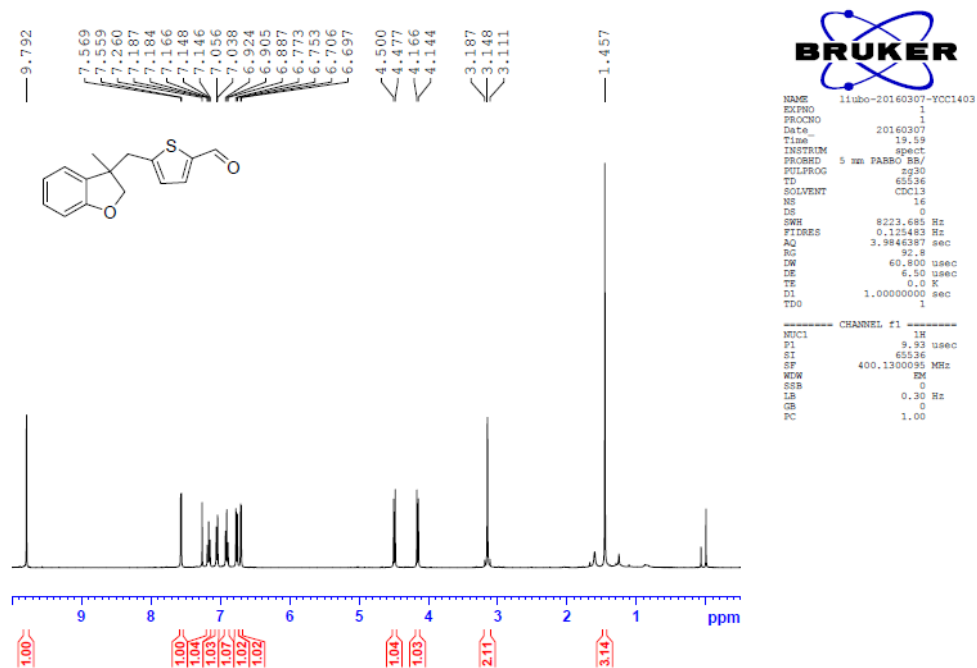
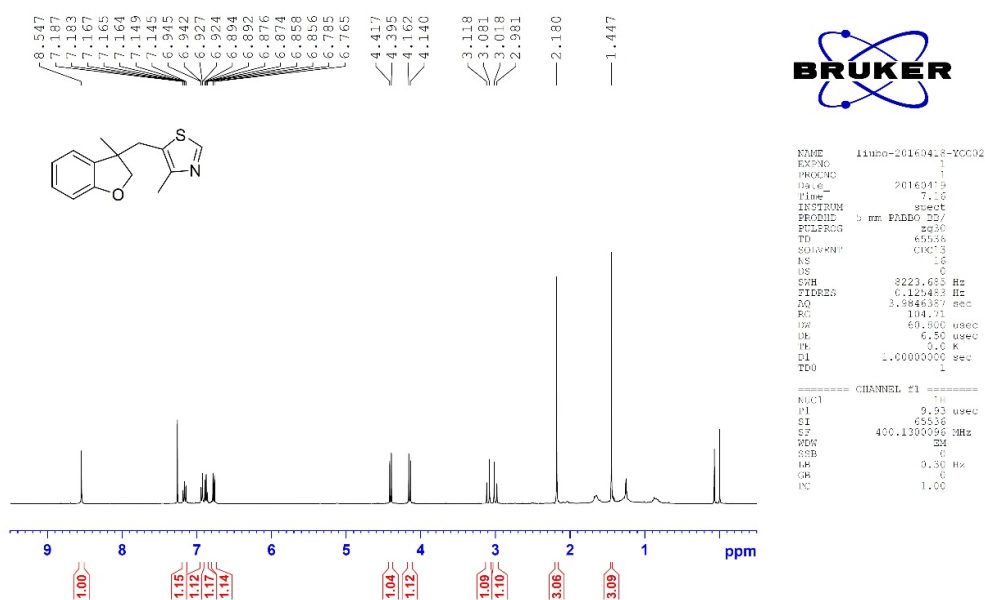
¹H NMR Spectra for Compound 1i¹H NMR Spectra for Compound 1j

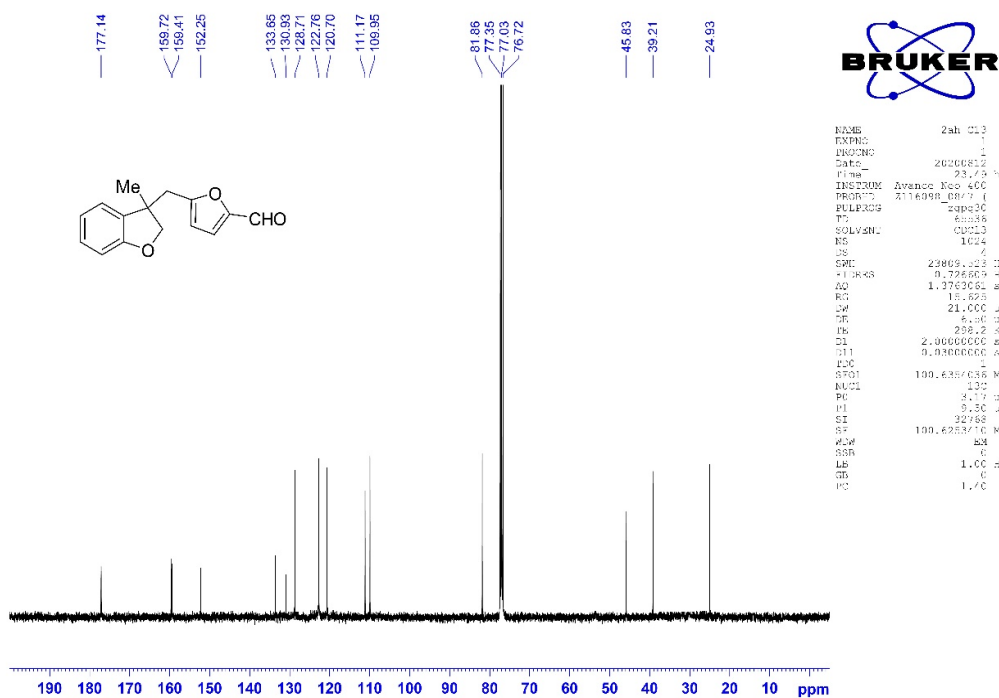
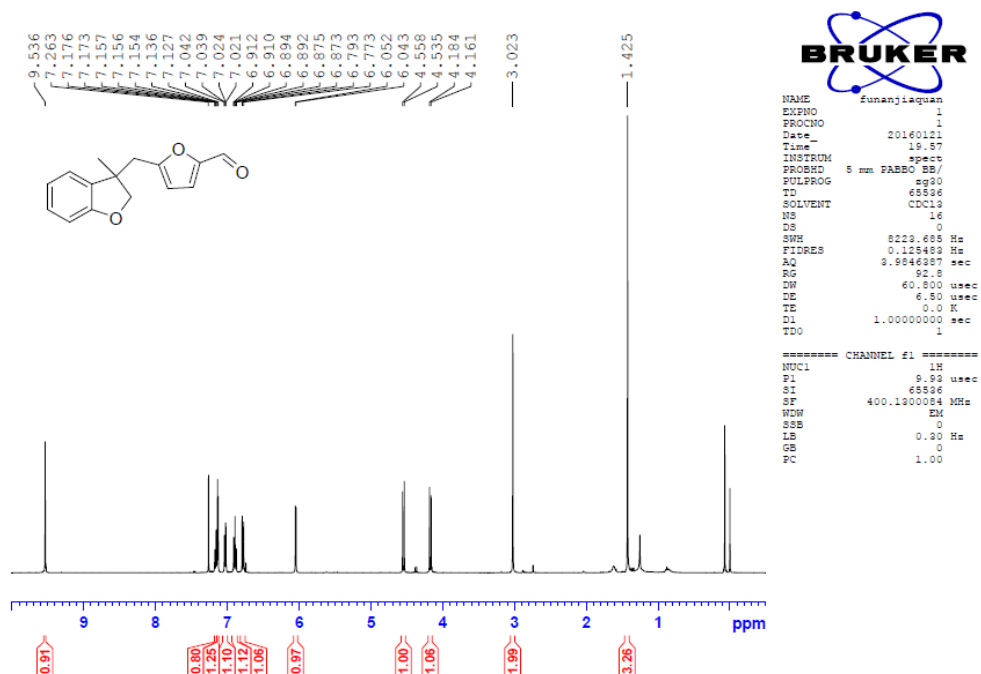
¹H NMR Spectra for Compound 2aa¹H and ¹³C NMR Spectra for Compound 2ab

¹H and ¹³C NMR Spectra for Compound 2ac

¹H and ¹³C NMR Spectra for Compound 2ad

¹H NMR Spectra for Compound 2ae

¹H NMR Spectra for Compound 2af¹H NMR Spectra for Compound 2ag

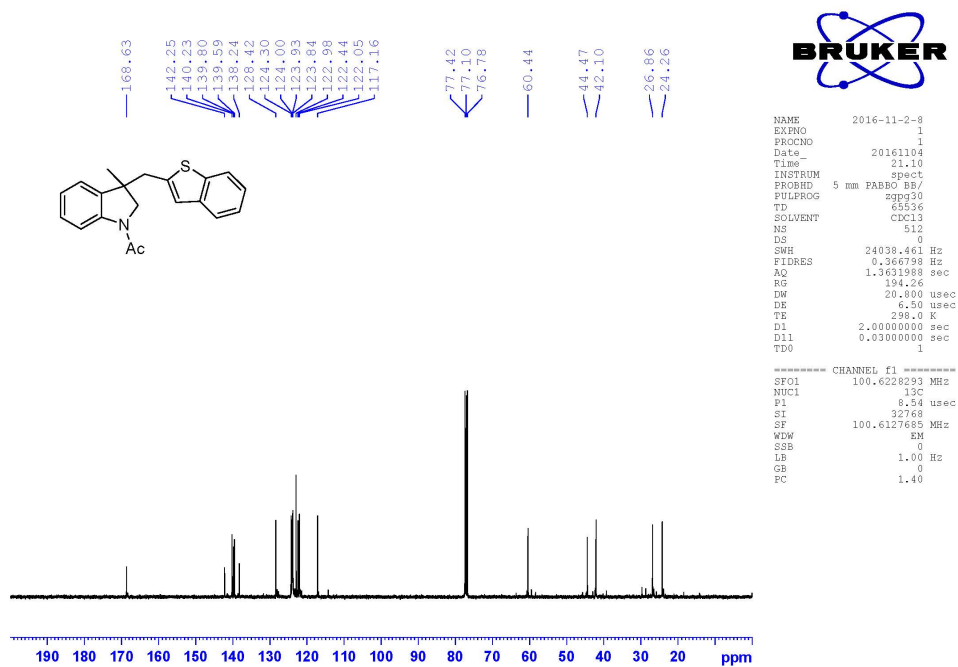
^1H and ^{13}C NMR Spectra for Compound 2ah

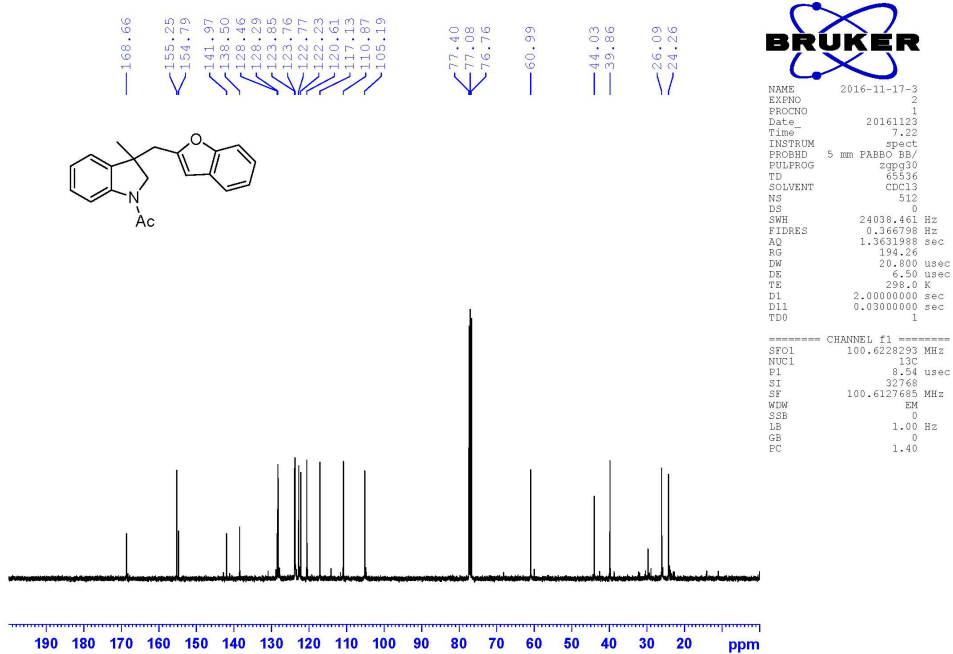
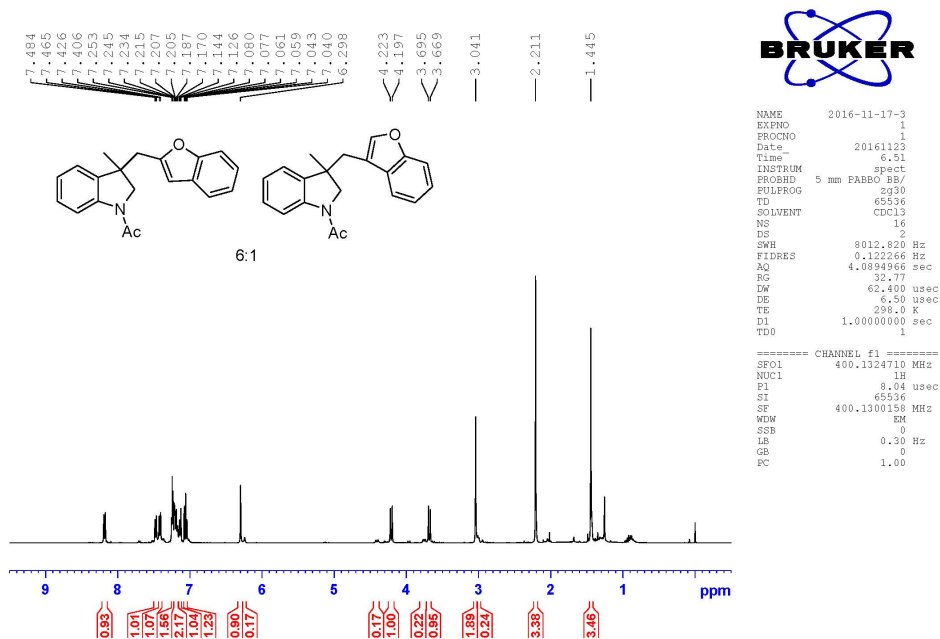
Chemical structures of the 5:1 ratio of diastereomers:

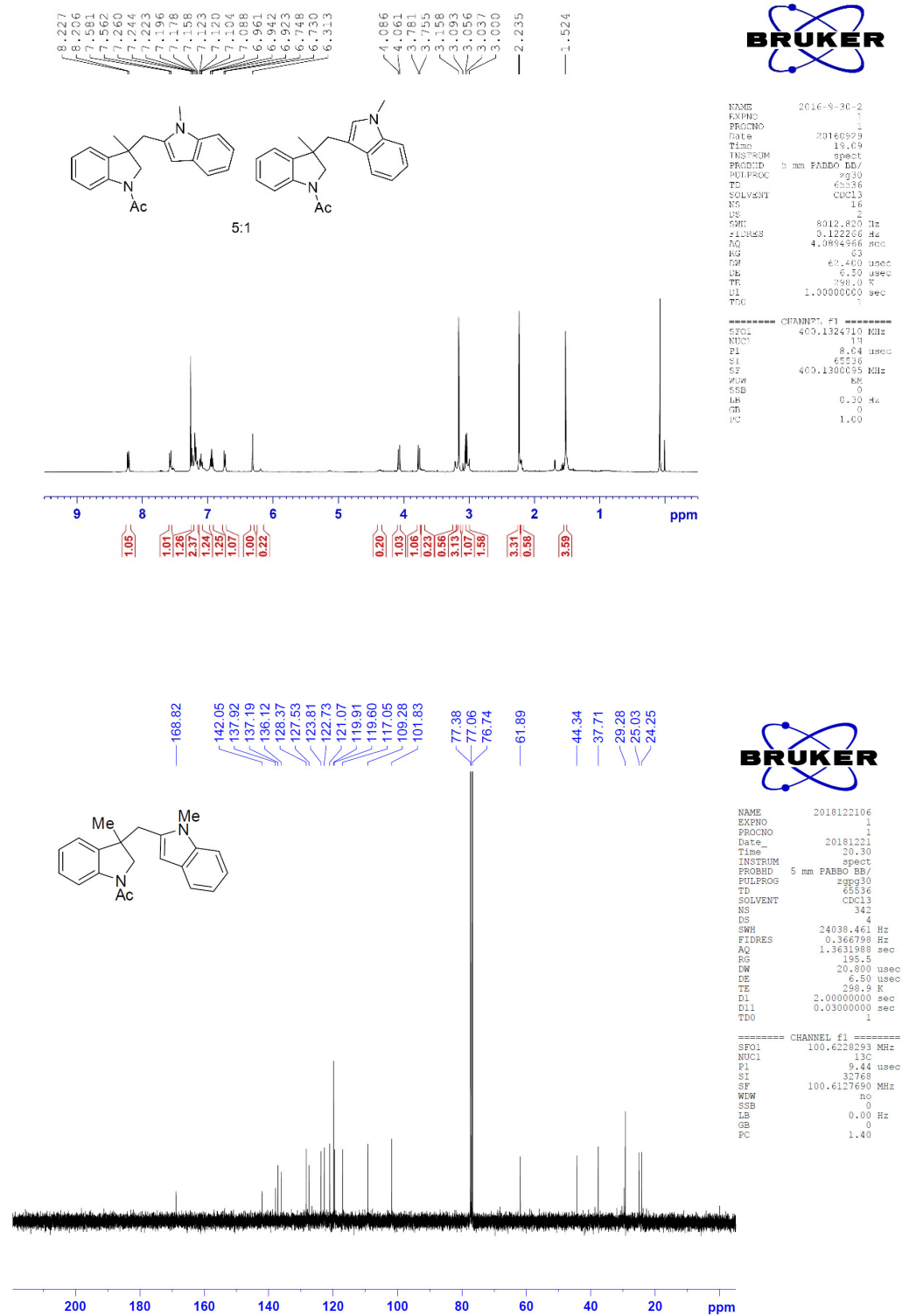
CC(=O)N1CC[C@H](C1)Cc2sc3ccccc32 (Major isomer, 5 parts)
CC(=O)N1CC[C@H](C1)Cc2cc3ccccc3s2 (Minor isomer, 1 part)

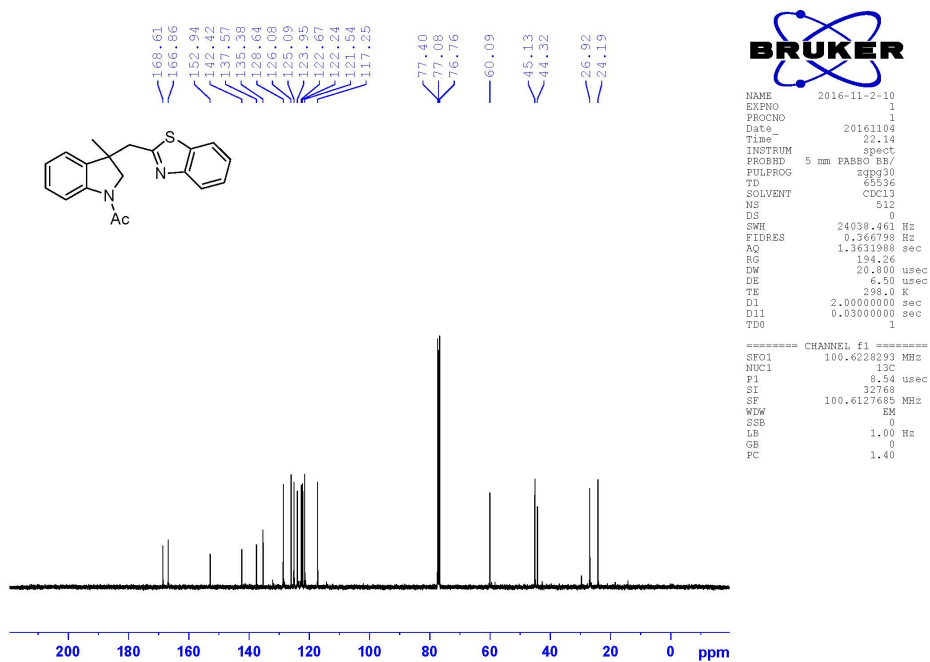
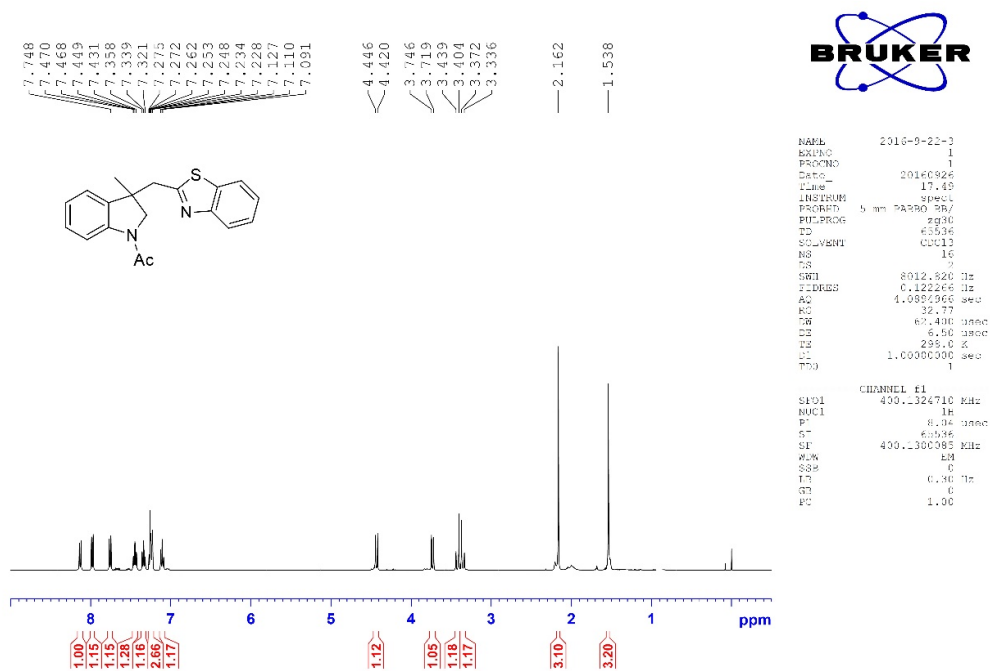
¹H NMR spectrum (CDCl₃) showing peaks from 0 to 8 ppm. The spectrum includes a list of peak chemical shifts (delta) and integrations.

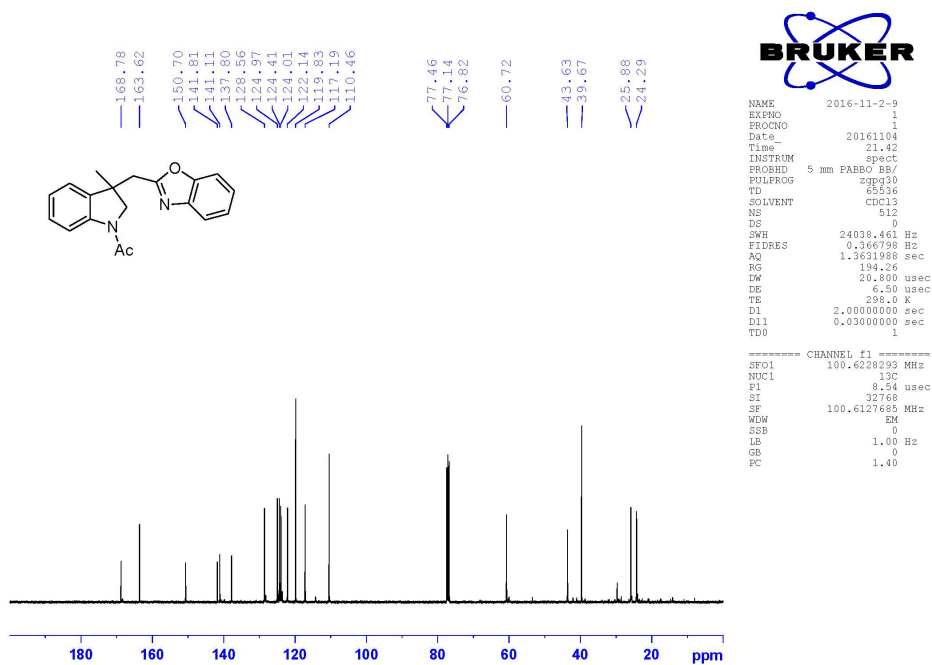
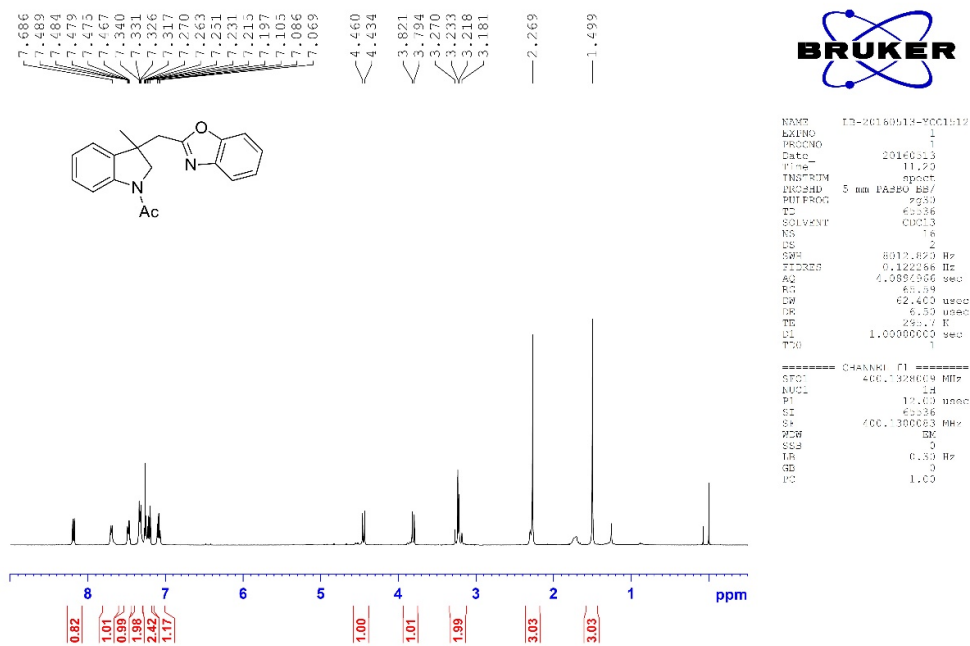
Chemical Shift (ppm)	Integration
7.731	1.00
7.712	0.10
7.667	0.90
7.649	1.26
7.325	4.37
7.310	1.19
7.307	1.33
7.291	0.86
7.284	0.29
7.261	
7.262	
7.252	
7.245	
7.243	
7.167	
7.170	
7.111	
7.108	
7.092	
7.090	
7.074	
6.504	
4.125	
4.098	
3.766	
3.676	
3.650	
3.197	
3.190	
2.166	
1.491	

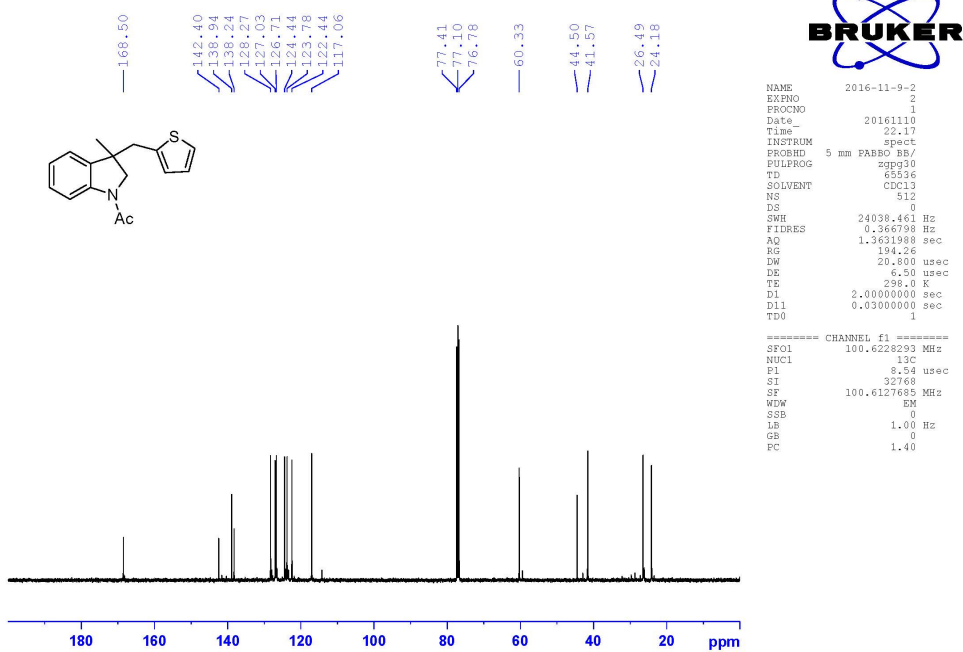
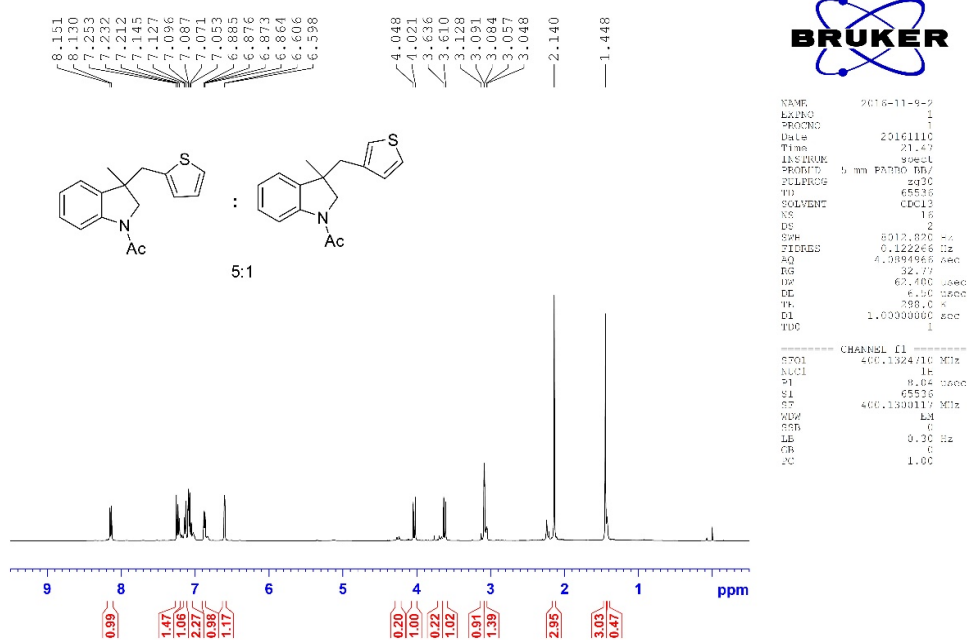


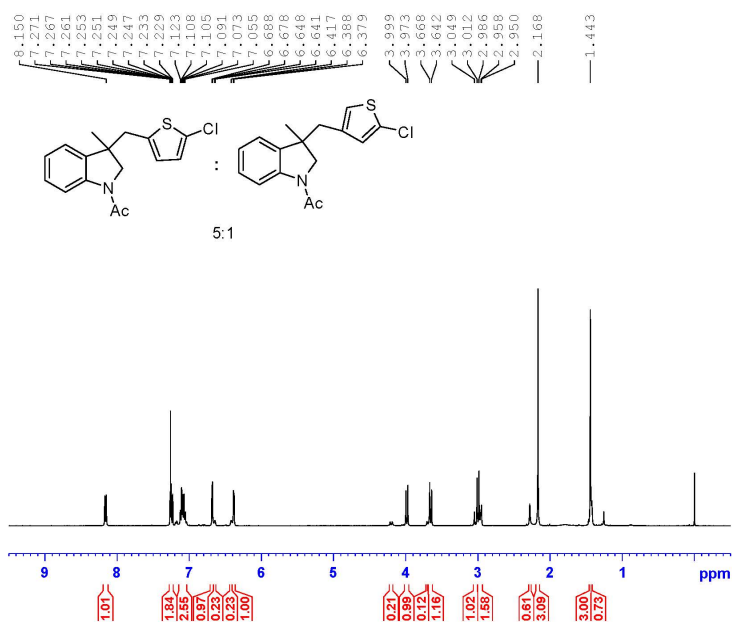
¹H and ¹³C NMR Spectra for Compound **2bb**

^1H and ^{13}C NMR Spectra for Compound **2b**

^1H and ^{13}C NMR Spectra for Compound **2bd**

¹H and ¹³C NMR Spectra for Compound 2b

¹H and ¹³C NMR Spectra for Compound 2bf

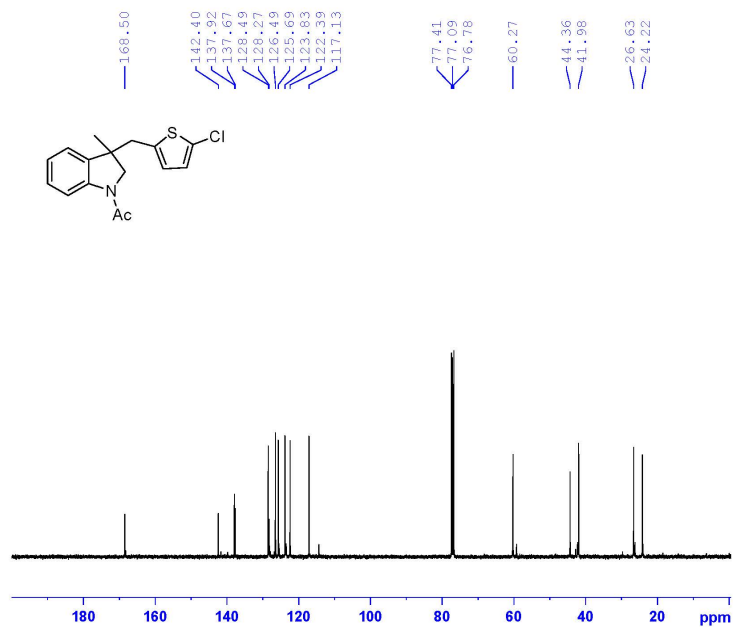
¹H and ¹³C NMR Spectra for Compound 2bg

```

NAME      2016-10-19-3
EXPNO     2
PROCNO    1
Date_     20161021
Time      19.15
INSTRUM   spect
PROBHD    5 mm PABBO BB/
PULPROG   zg30
TD         65536
SOLVENT   CDCl3
NS         16
DS         2
SWH        8012.820 Hz
FIDRES     0.122266 Hz
AQ         4.0894966 sec
RG         63
DM         62.400 usec
DE         6.50 usec
TE         298.0 K
D1         1.00000000 sec
TD0        1

===== CHANNEL f1 =====
SF01      400.1324710 MHz
NUC1       1H
P1         8.04 usec
SI         65536
SF         400.1300096 MHz
WDW        EM
SSB         0
LB         0.30 Hz
GB         0
PC         1.00

```

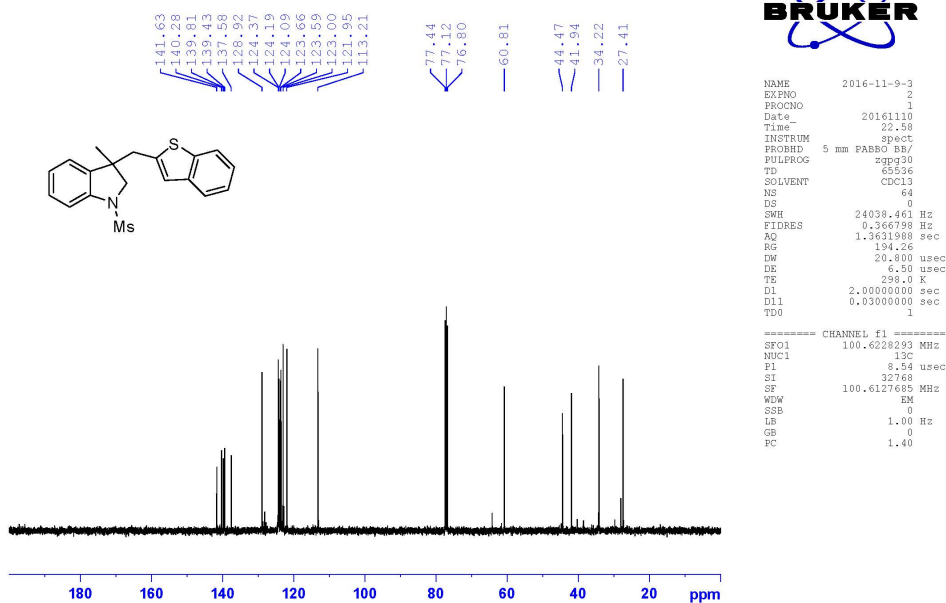
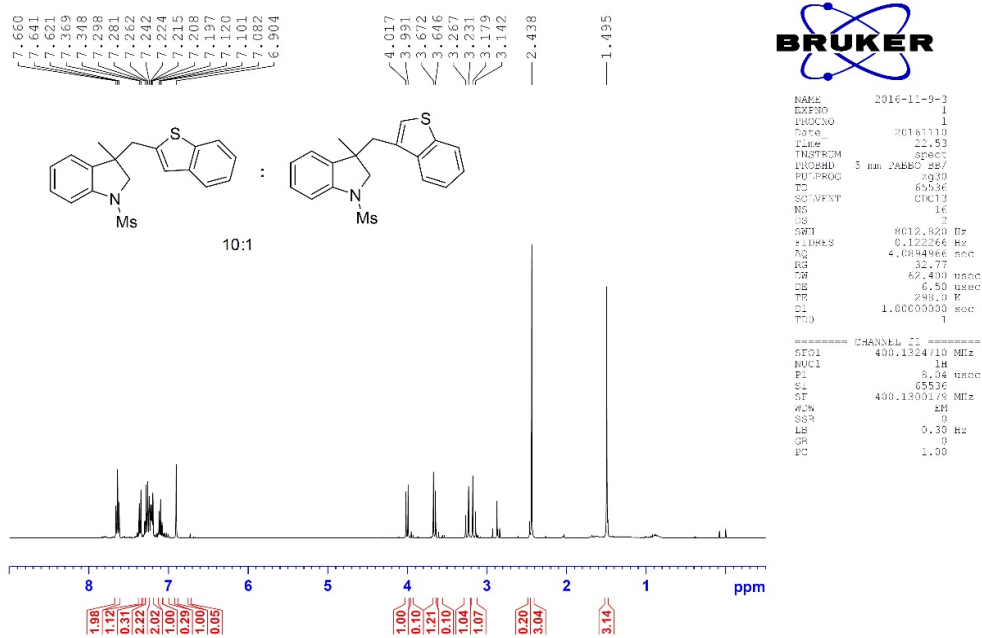


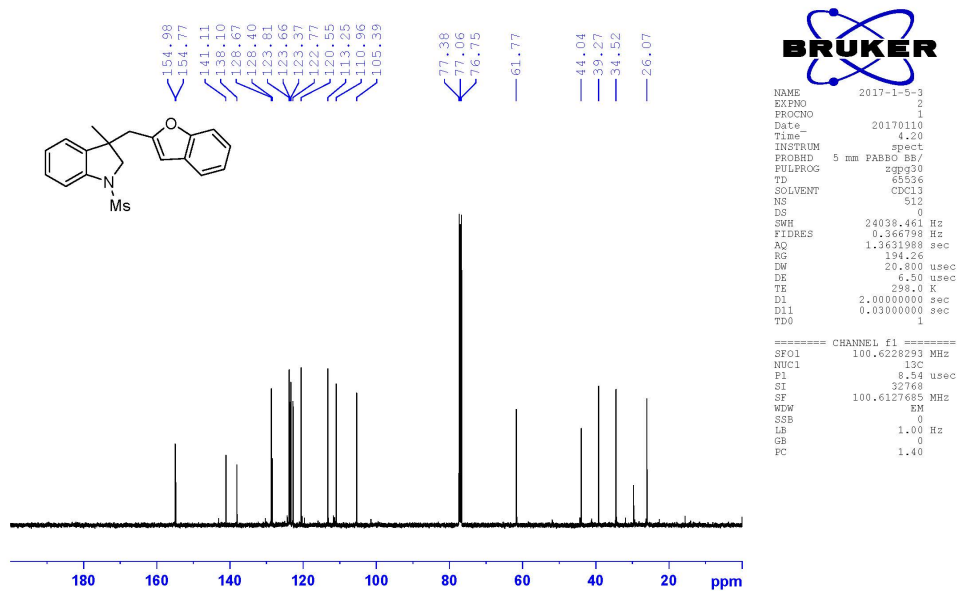
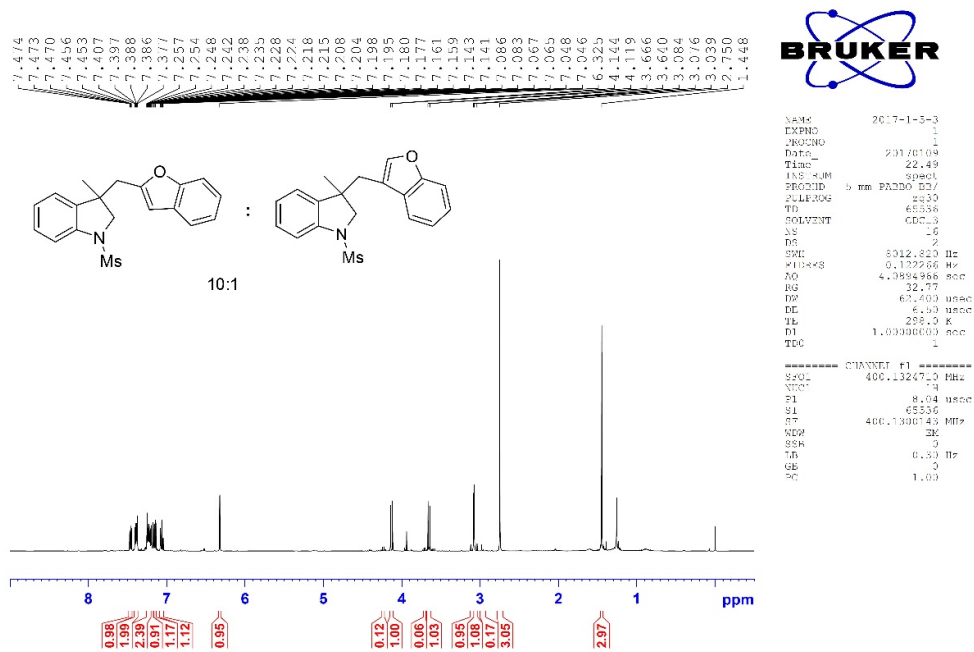
```

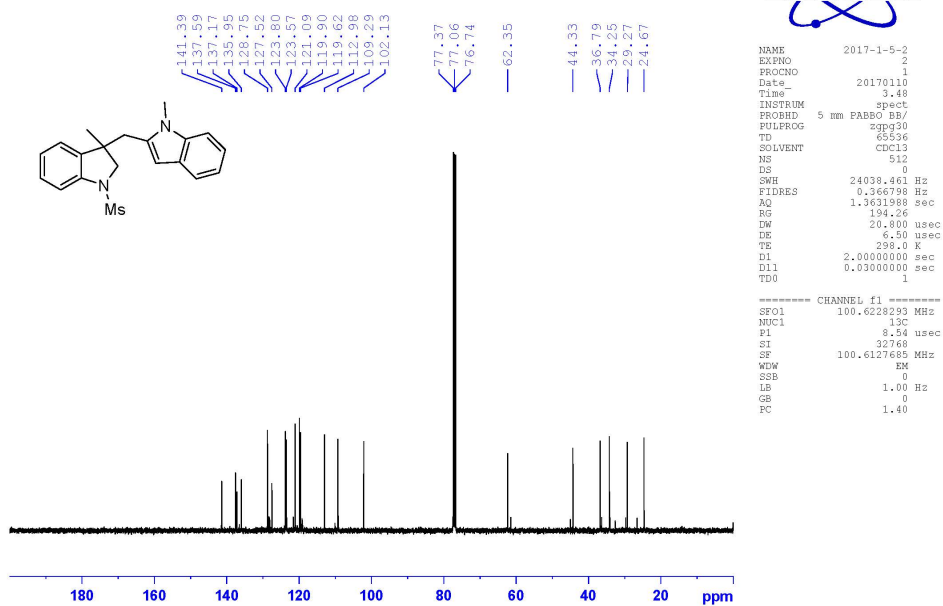
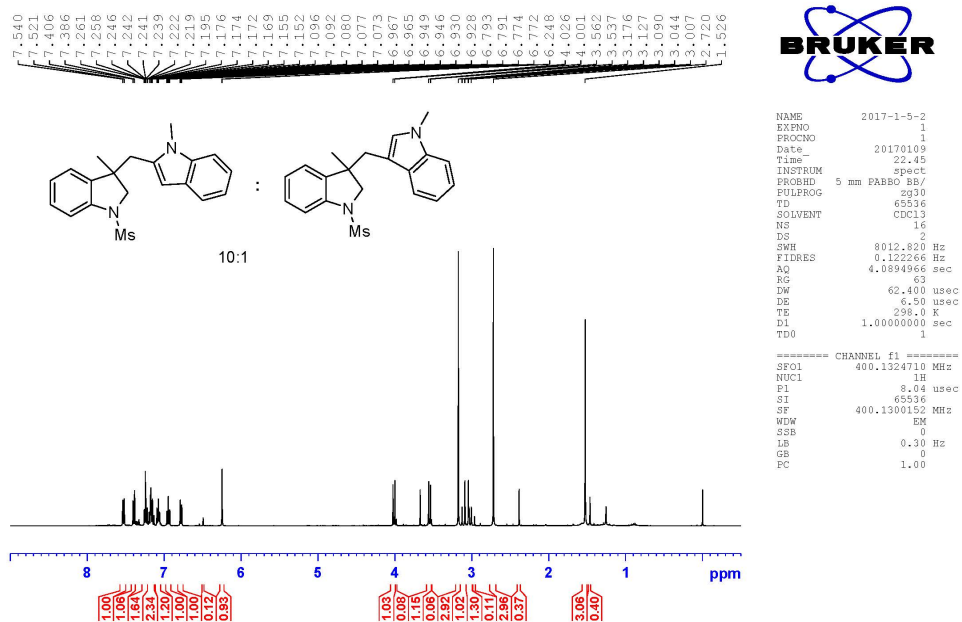
NAME      2016-11-2-7
EXPNO     1
PROCNO    1
Date_     20161104
Time      20.38
INSTRUM   spect
PROBHD    5 mm PABBO BB/
PULPROG   zgpg30
TD         65536
SOLVENT   CDCl3
NS         512
DS         0
SWH        24038.461 Hz
FIDRES     0.366798 Hz
AQ         1.3631988 sec
RG         194.25
DM         20.800 usec
DE         6.50 usec
TE         298.0 K
D1         2.00000000 sec
D11        0.03000000 sec
TD0        1

===== CHANNEL f1 =====
SF01      100.6228293 MHz
NUC1       13C
P1         8.54 usec
SI         32768
SF         100.6127685 MHz
WDW        EM
SSB         0
LB         1.00 Hz
GB         0
PC         1.40

```


¹H and ¹³C NMR Spectra for Compound 2ca

¹H and ¹³C NMR Spectra for Compound 2cb

¹H and ¹³C NMR Spectra for Compound 2cc

Chemical structure: 1-methyl-2-(thiophen-2-ylmethyl)pyrrolidine

¹H NMR spectrum (400 MHz, CDCl₃) showing peaks from 2.4 to 7.2 ppm. The x-axis is labeled 'ppm'.

Integration values (from left to right): 1.00, 1.06, 1.06, 1.12, 1.04, 0.99, 0.99, 1.02, 1.02, 1.04, 1.04, 3.13, 3.09.

Chemical structure inset:

CN1CCCC1Cc2ccsc2

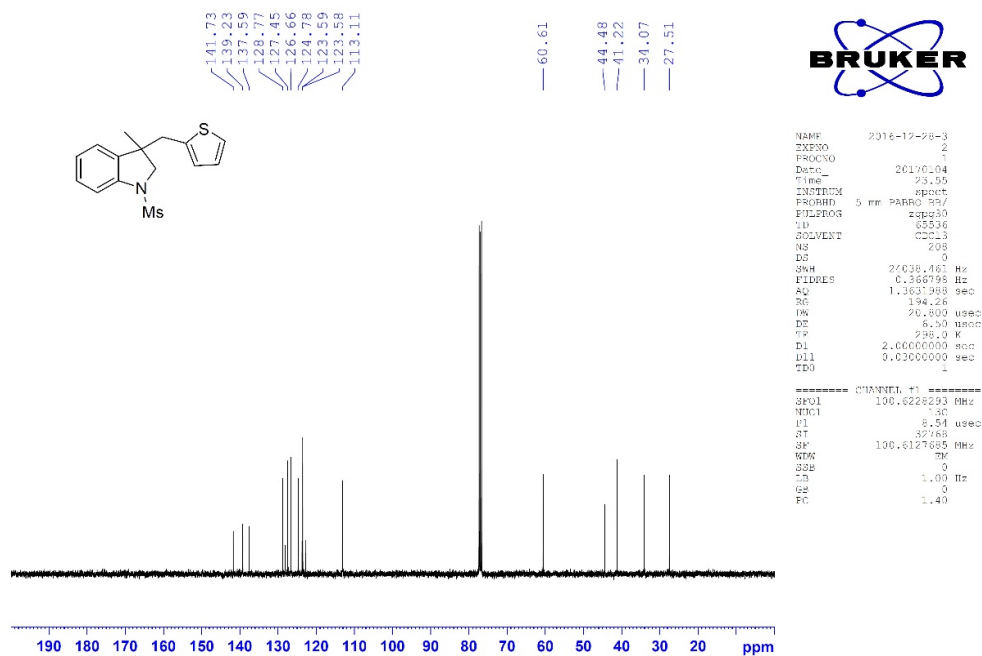
Peak list (ppm): 7.2419, 7.2232, 7.2229, 7.2000, 7.1977, 7.1817, 7.1719, 7.1119, 7.1117, 7.1017, 7.0987, 7.0800, 7.0663, 7.0663, 7.0500, 7.0500, 6.8777, 6.8668, 6.8684, 6.8566, 6.6577, 6.6448, 3.9448, 3.9222, 3.6866, 3.6666, 3.6330, 3.6330, 3.2117, 3.2117, 3.1755, 3.1117, 3.0800, 2.4933, 1.478.

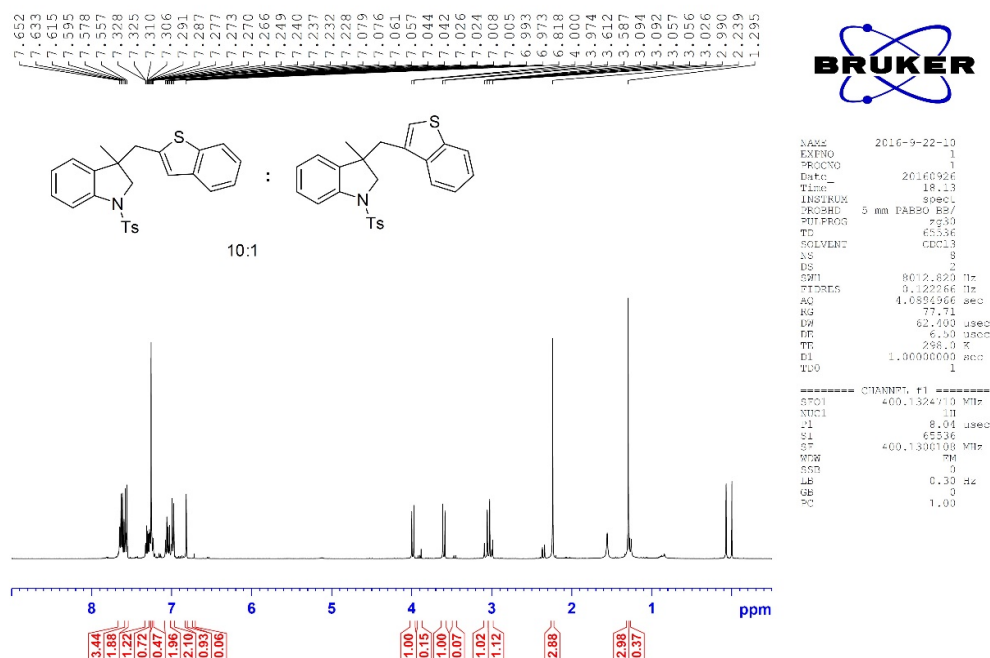
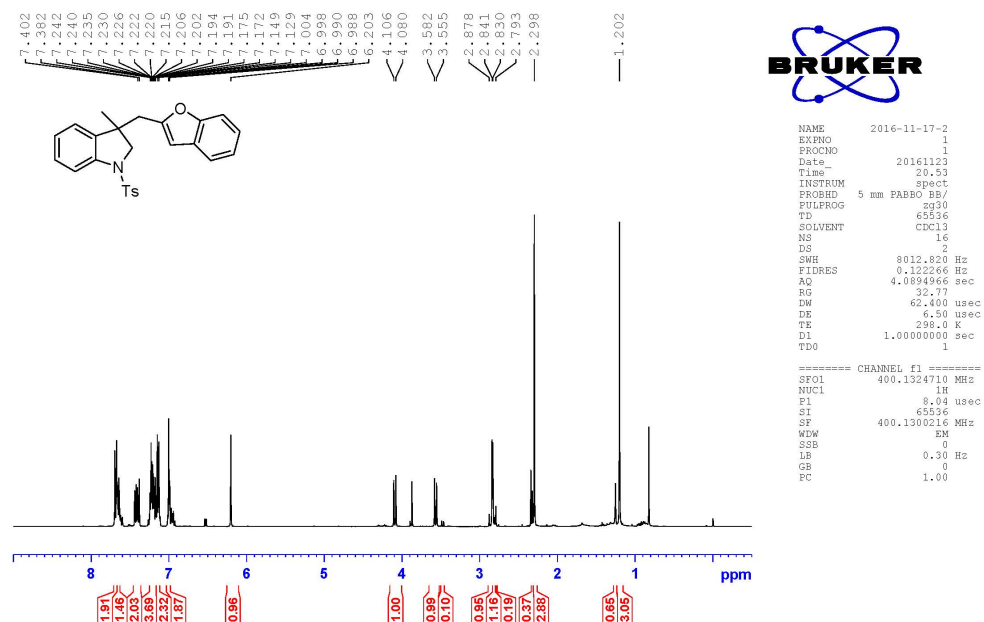
Integration values (from left to right): 1.00, 1.06, 1.06, 1.12, 1.04, 0.99, 0.99, 1.02, 1.02, 1.04, 1.04, 3.13, 3.09.

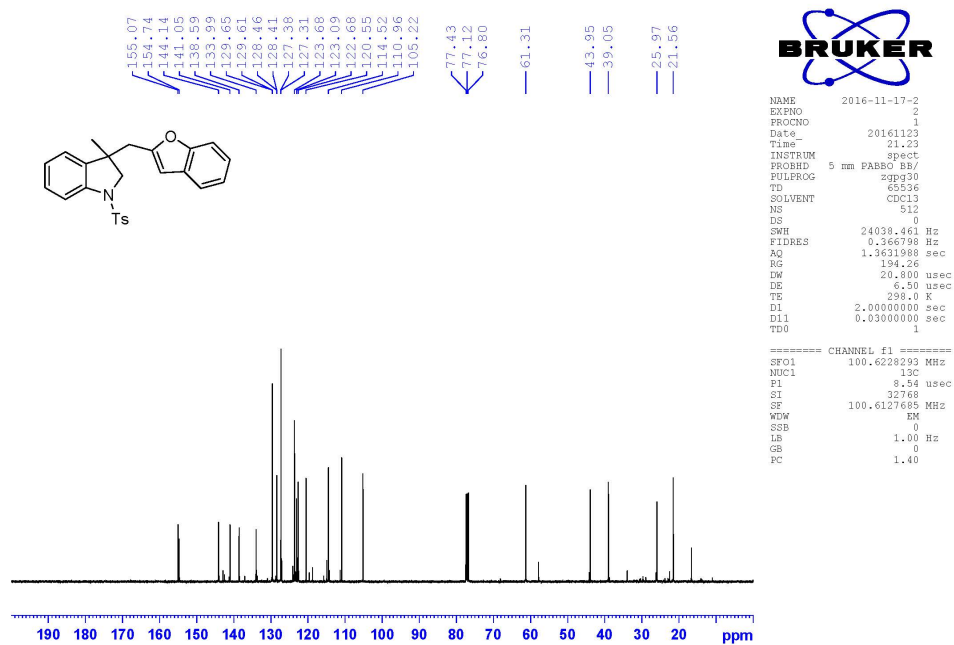
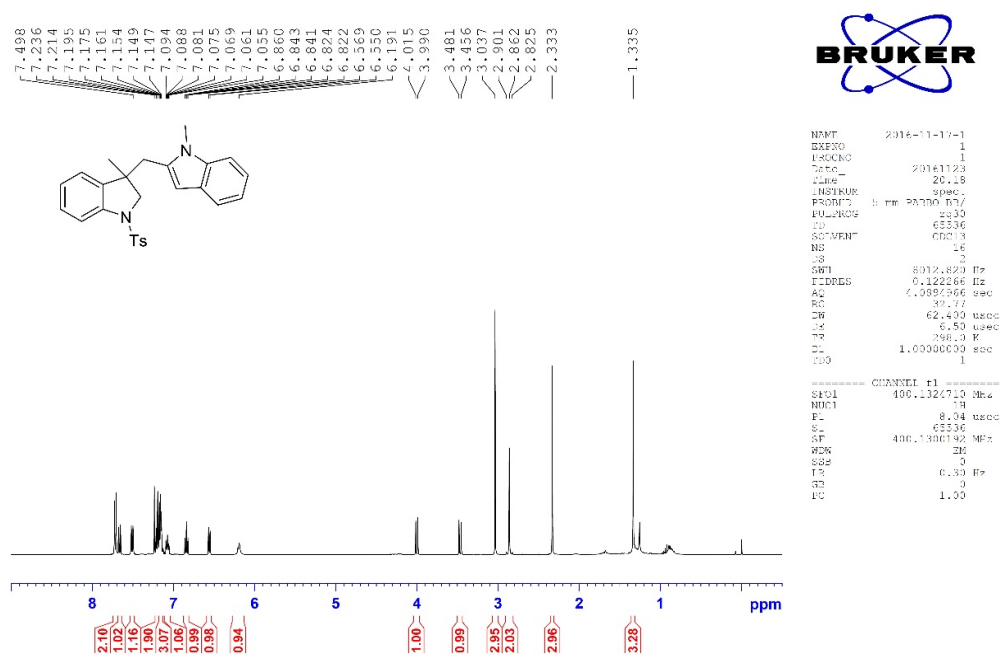
Chemical structure inset:

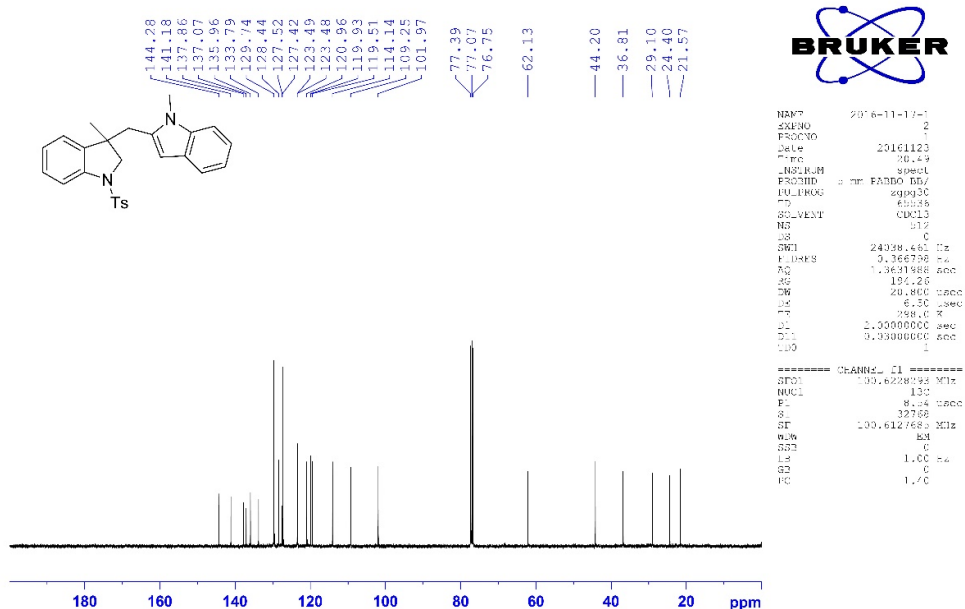
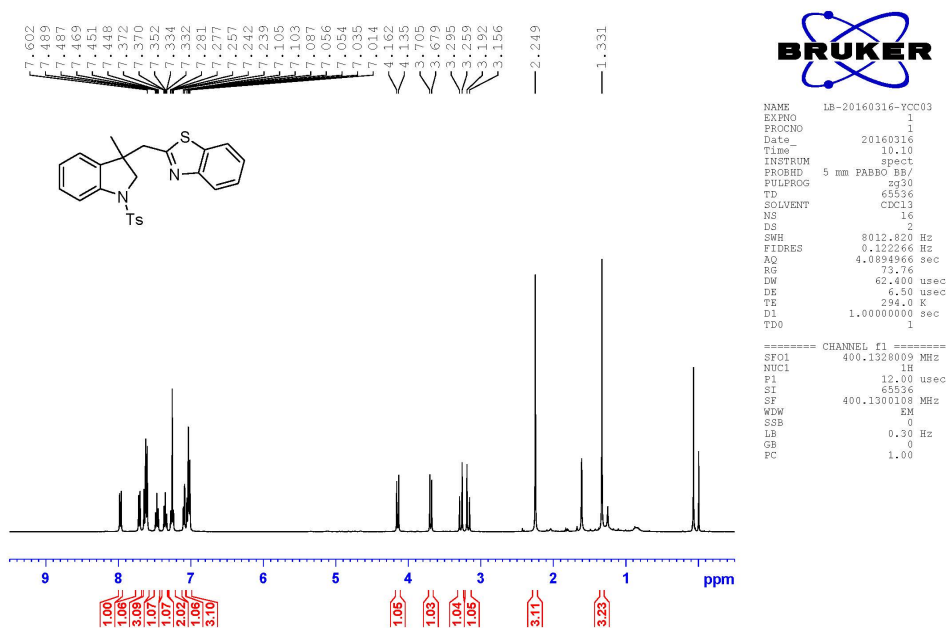
CN1CCCC1Cc2ccsc2

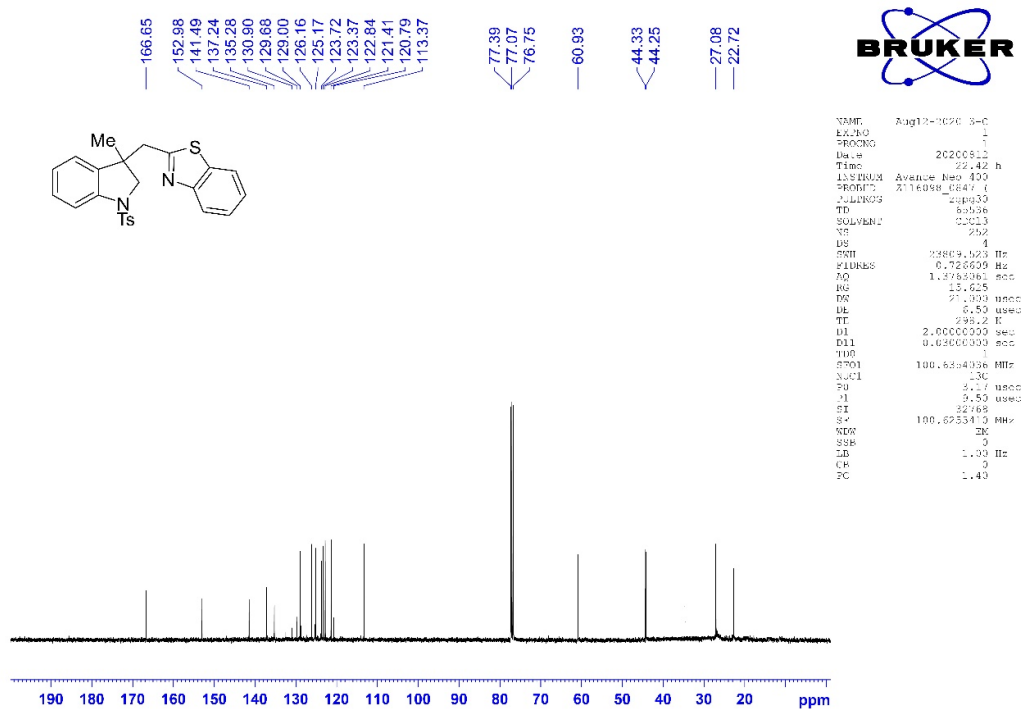
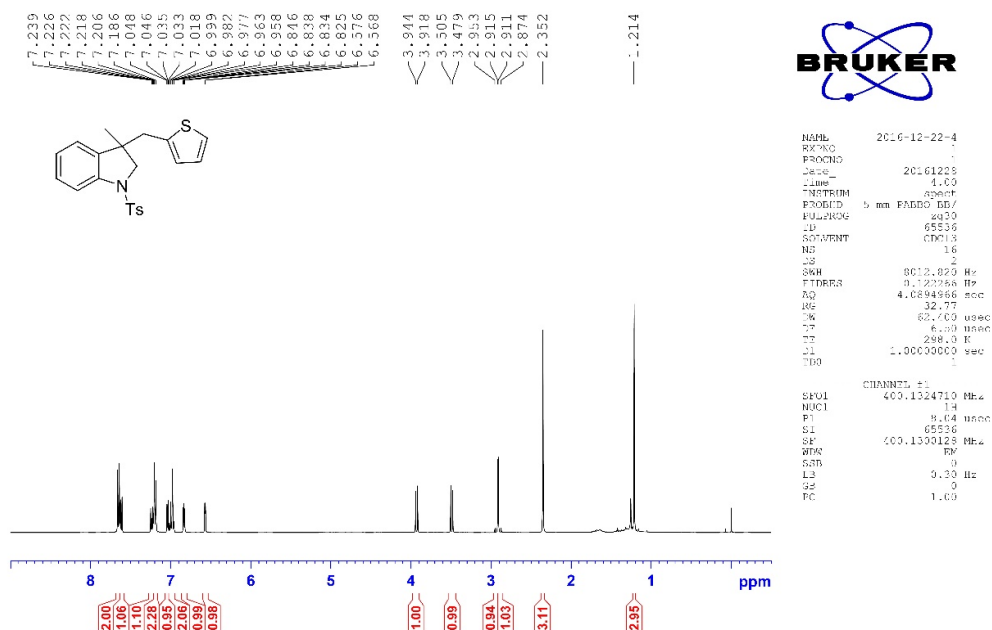
Peak list (ppm): 7.2419, 7.2232, 7.2229, 7.2000, 7.1977, 7.1817, 7.1719, 7.1119, 7.1117, 7.1017, 7.0987, 7.0800, 7.0663, 7.0663, 7.0500, 7.0500, 6.8777, 6.8668, 6.8684, 6.8566, 6.6577, 6.6448, 3.9448, 3.9222, 3.6866, 3.6666, 3.6330, 3.6330, 3.2117, 3.2117, 3.1755, 3.1117, 3.0800, 2.4933, 1.478.

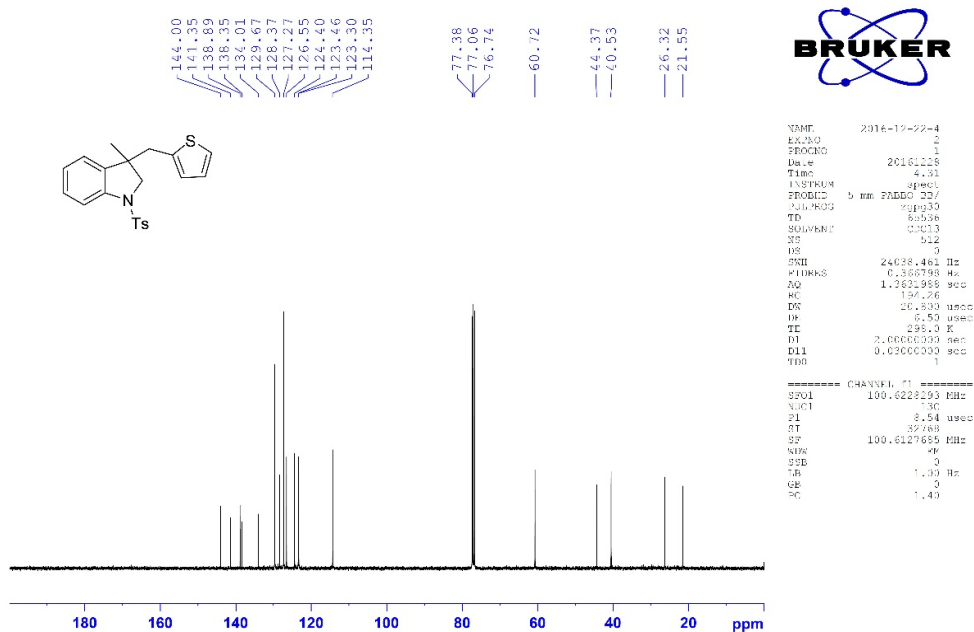


¹H Spectra for Compound 2da¹H and ¹³C NMR Spectra for Compound 2db

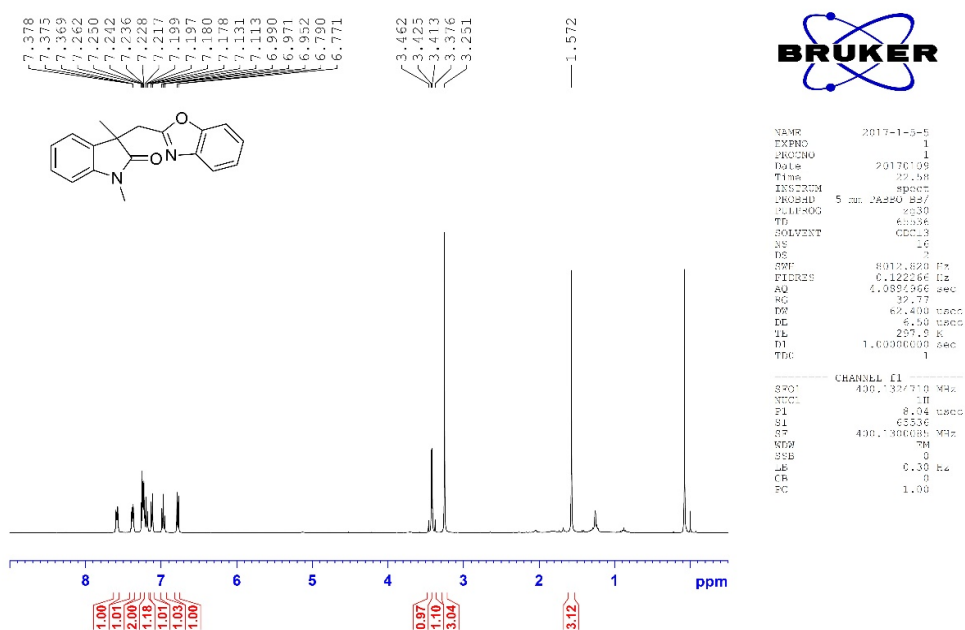
¹H and ¹³C NMR Spectra for Compound 2dc

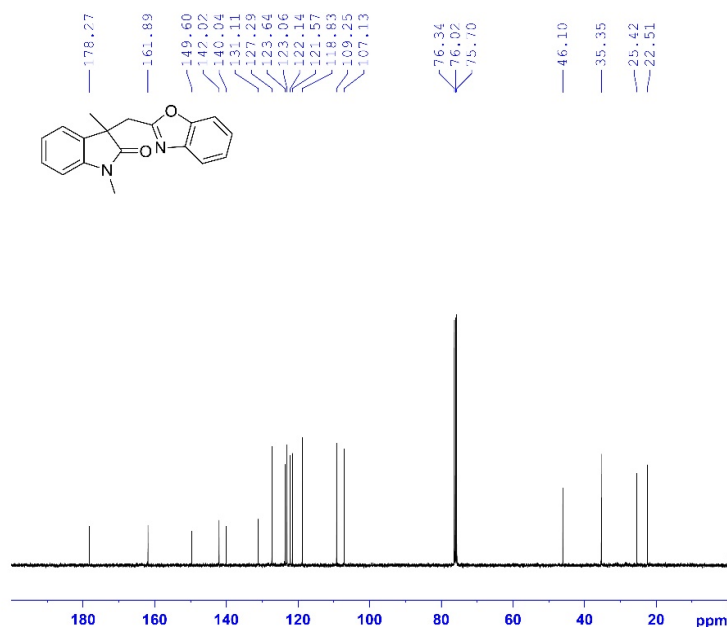
¹H and ¹³C NMR Spectra for Compound 2dd

¹H and ¹³C NMR Spectra for Compound 2de



¹H and ¹³C NMR Spectra for Compound 2ea





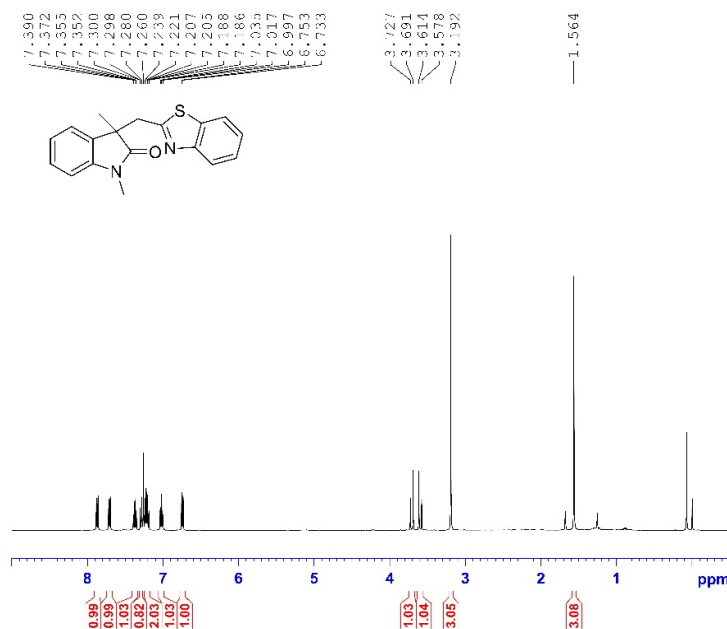
```

NAME      2017-1-5-3
EXPNO     1
PROCNO    1
Date_     2017/01/06
Time      9.24
INSTRUM   spect
PROBHD    5 mm PABBO BB/
PULPROG   zgpg30
TD         65536
SOLVENT   CDCl3
NS         512
DS         4
SWH        24038.441 Hz
FIDRES     0.366788 Hz
AQ         1.3631988 sec
RG         194.25
FR         20.806 sec
DE         6.50 usec
TE         298.0 K
D1         2.0000000 sec
D11        0.0200000 sec
TDC        1
  
```

```

===== CHANNEL f1 =====
NUC1       13C, 6228234 MHz
P1         130
PL         0.57 usec
PT         32.788
SF         100.6129730 MHz
WDW        EM
SSB         0
GB         1.00 Hz
CB         0
PC         1.40
  
```

¹H and ¹³C NMR Spectra for Compound 2eb

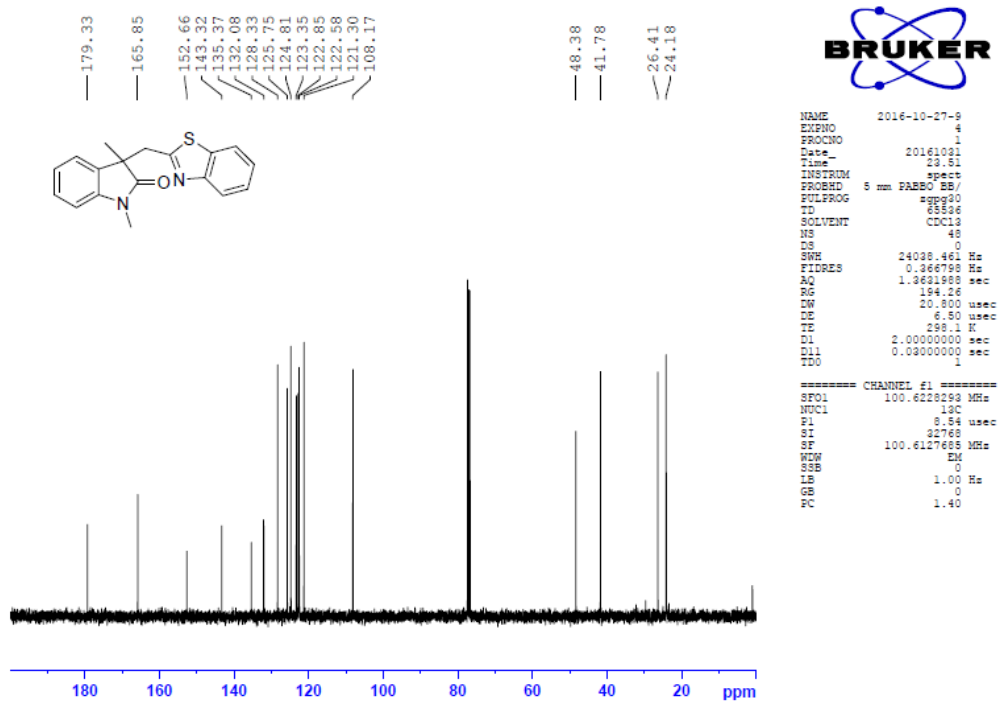
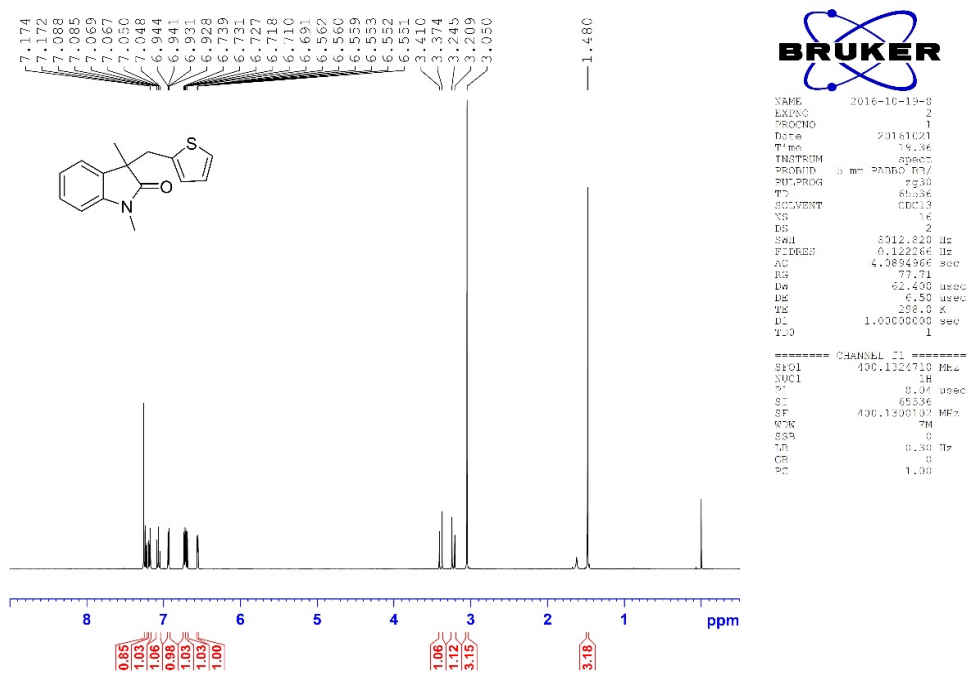


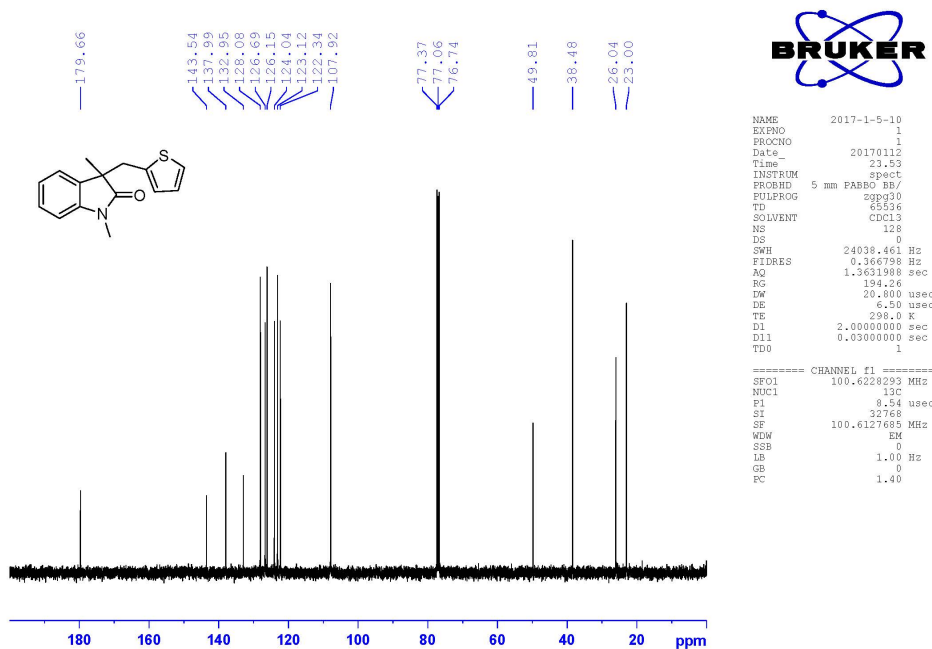
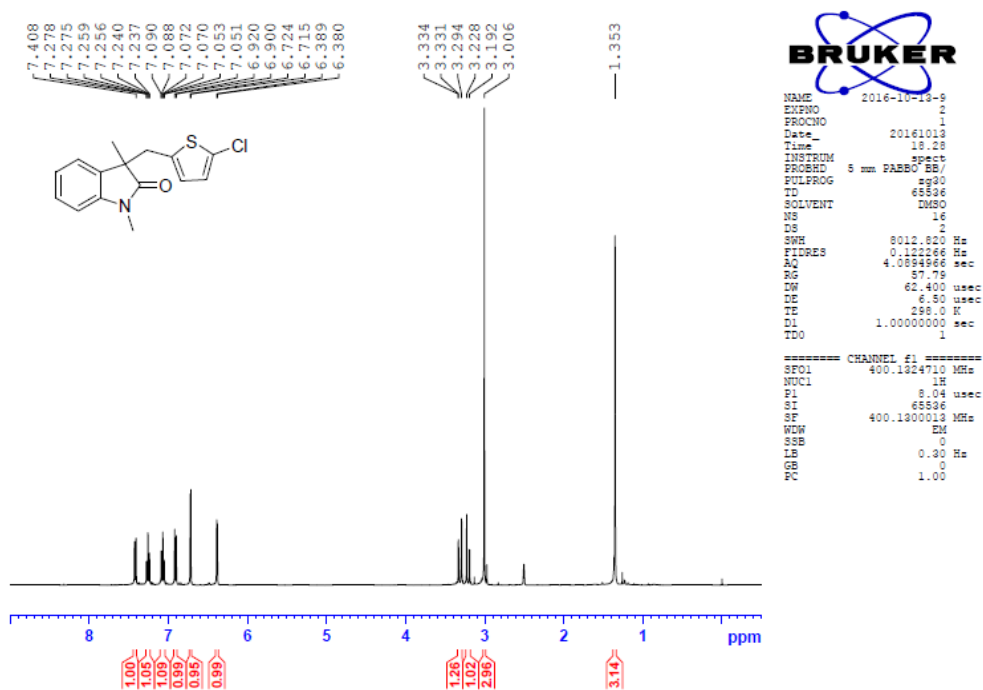
```

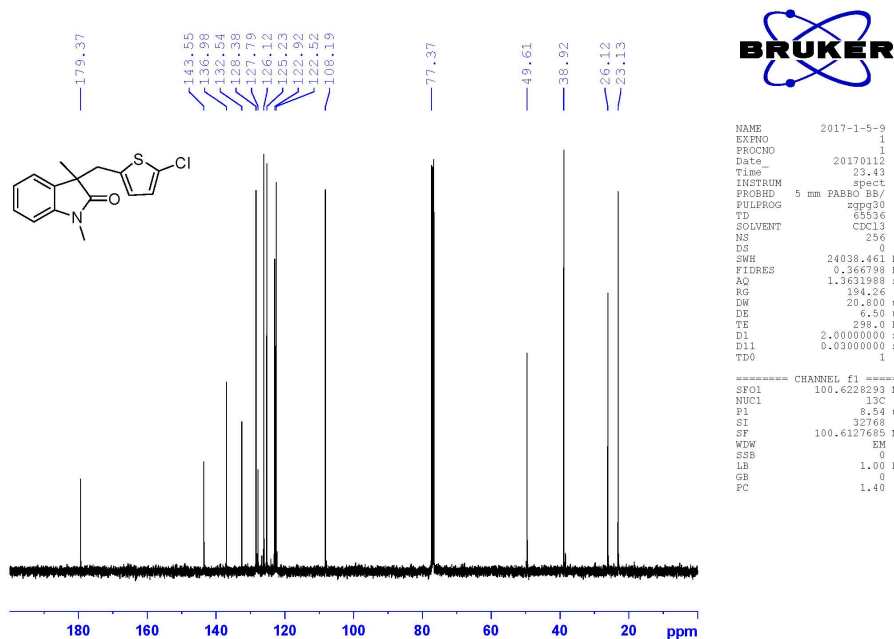
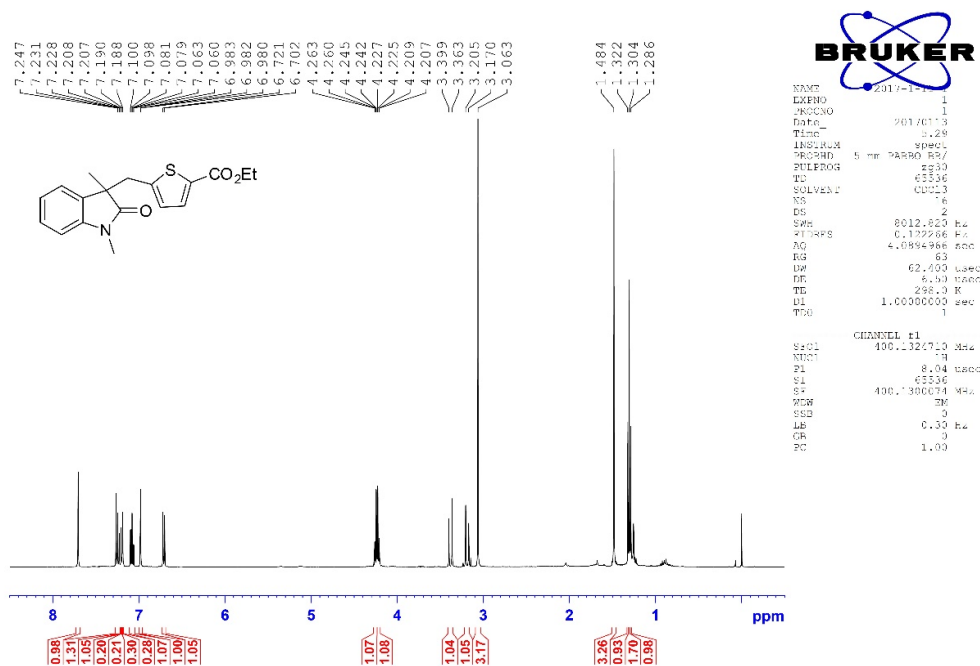
NAME      2016-9-8-1
EXPNO     1
PROCNO    1
Date_     2016/09/09
Time      17.59
INSTRUM   spect
PROBHD    5 mm PABBO BB/
PULPROG   zgpg30
TD         65536
SOLVENT   CDCl3
NS         16
DS         2
SWH        8012.820 Hz
FIDRES     0.122260 Hz
AQ         4.0894668 sec
RG         70.38
FR         62.400 usec
DE         6.50 usec
TE         298.0 K
D1         1.0000000 sec
D11        0
TDC        1
  
```

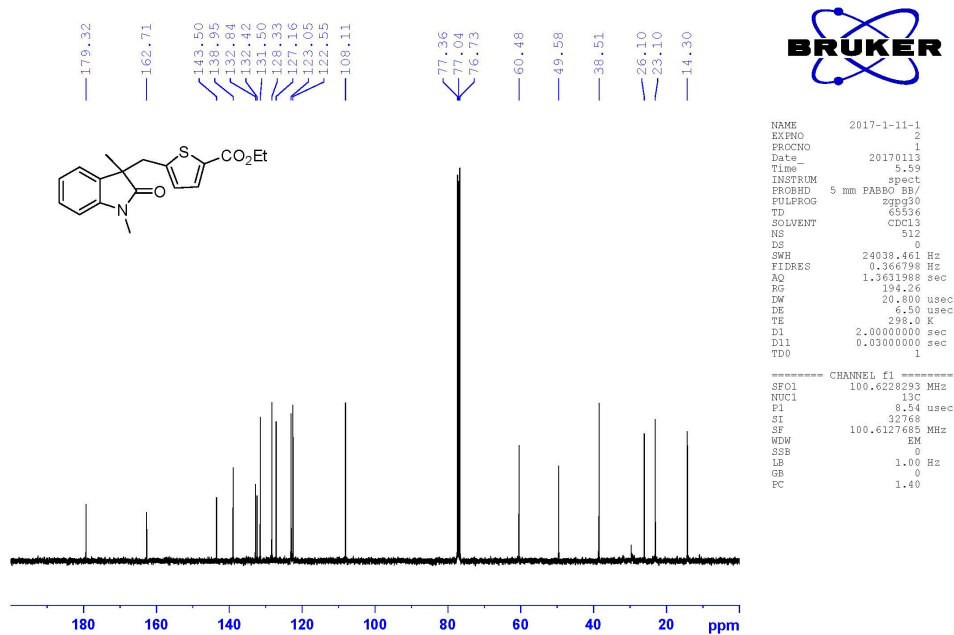
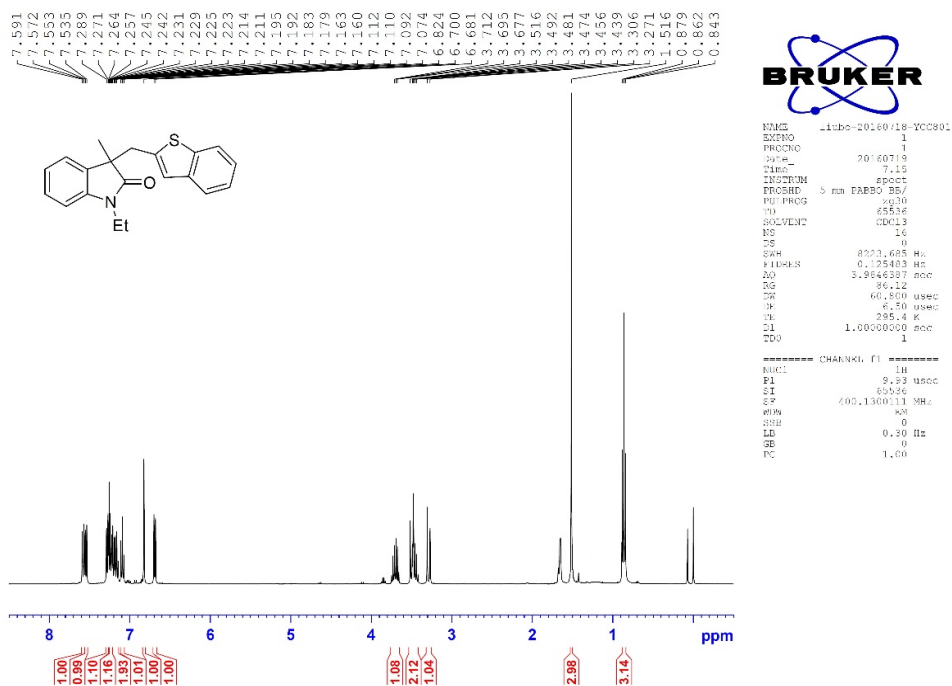
```

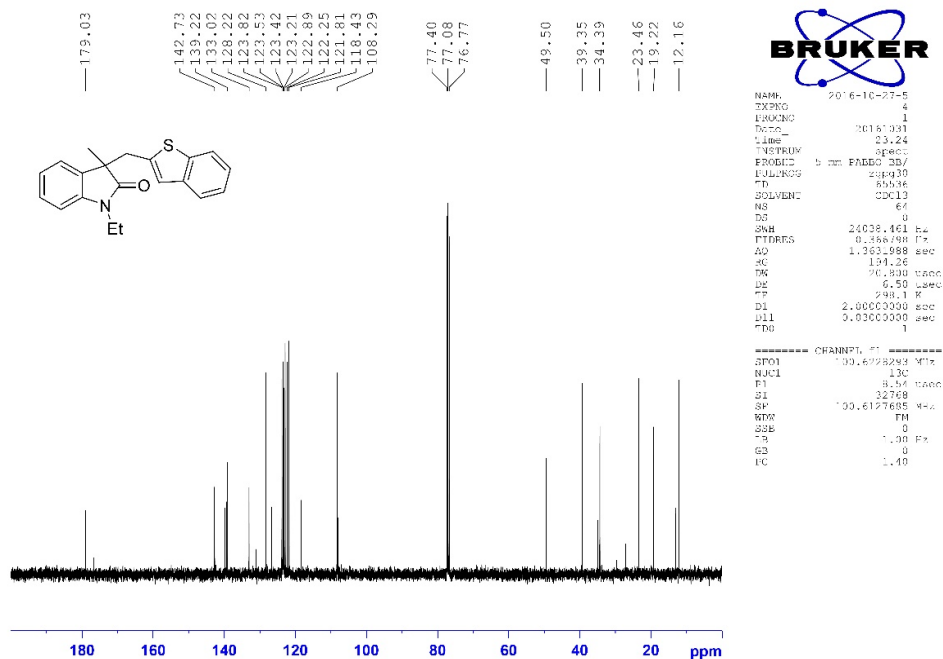
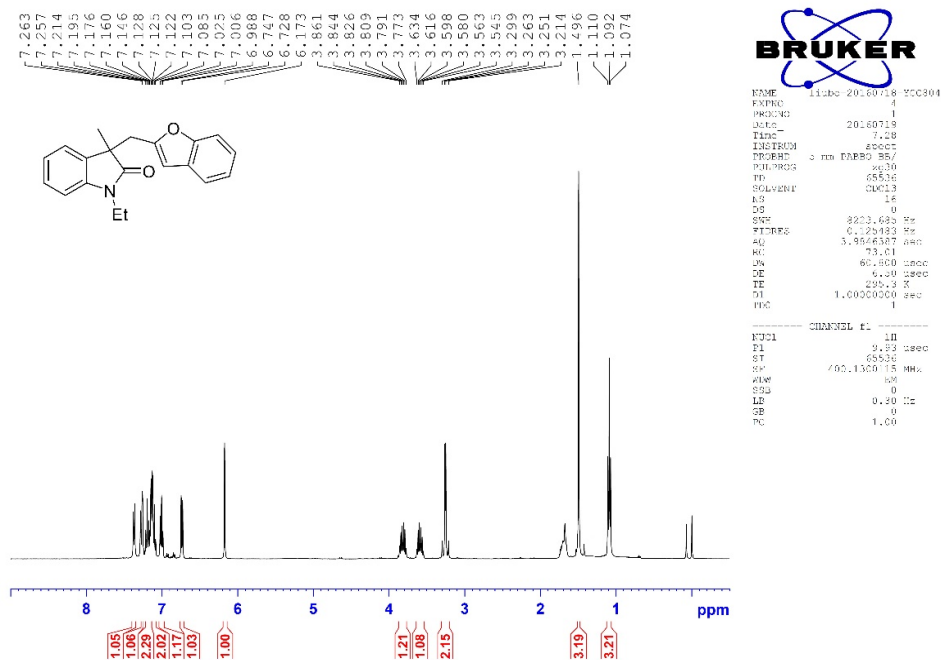
===== CHANNEL f1 =====
NUC1       1H, 400.132410 MHz
P1         130
PL         0.04 usec
PT         80.38
SF         400.130310 MHz
WDW        EM
SSB         0
GB         0
CB         0
PC         1.00
  
```

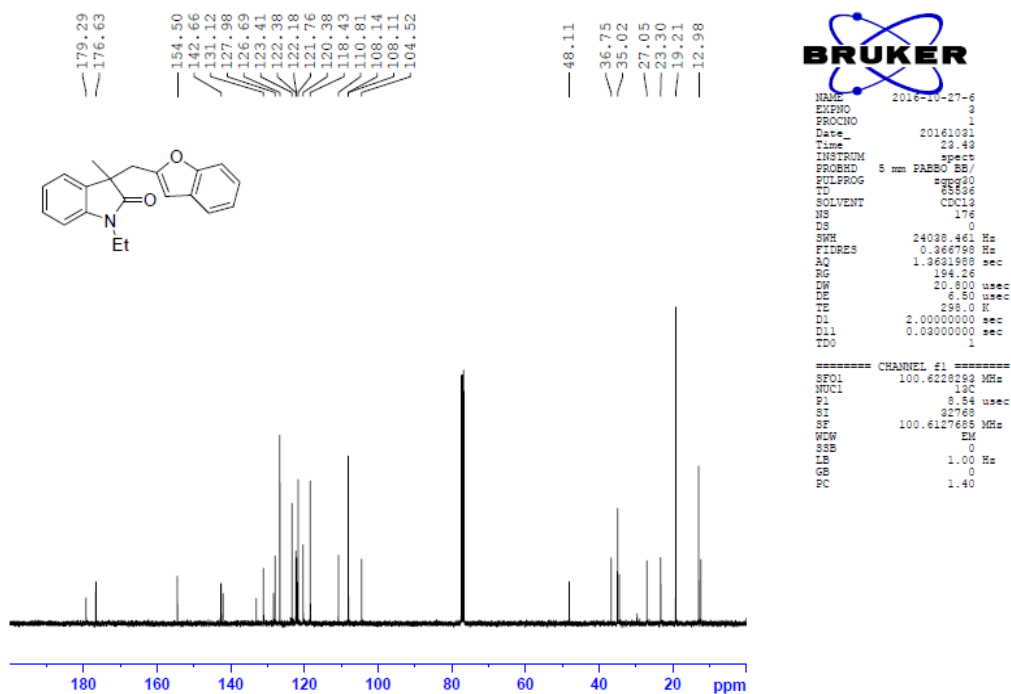
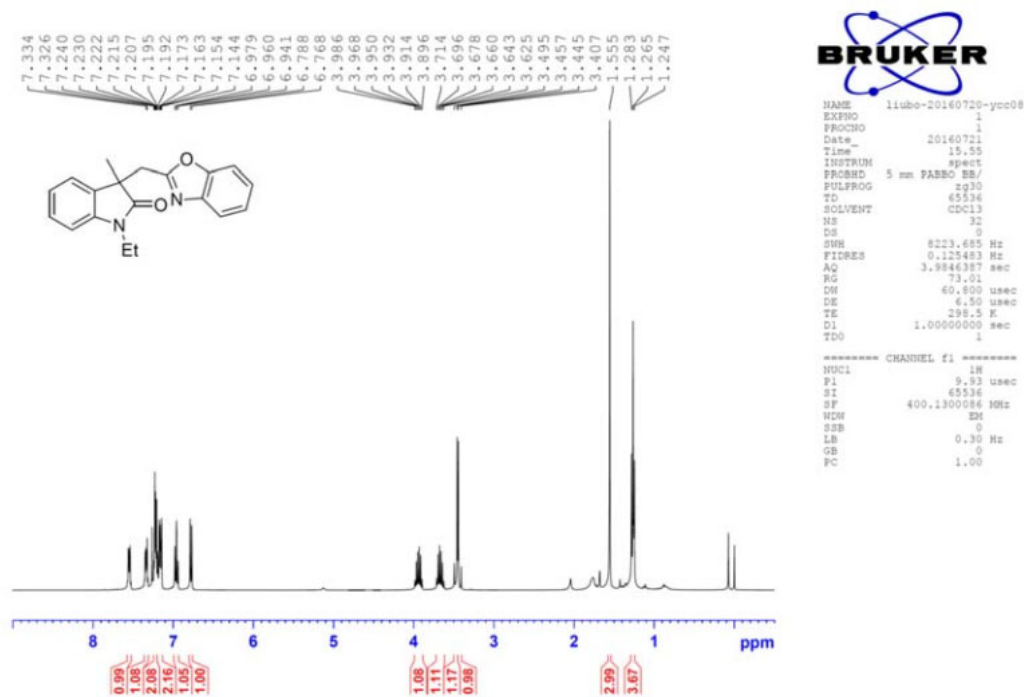
¹H and ¹³C NMR Spectra for Compound 2ec

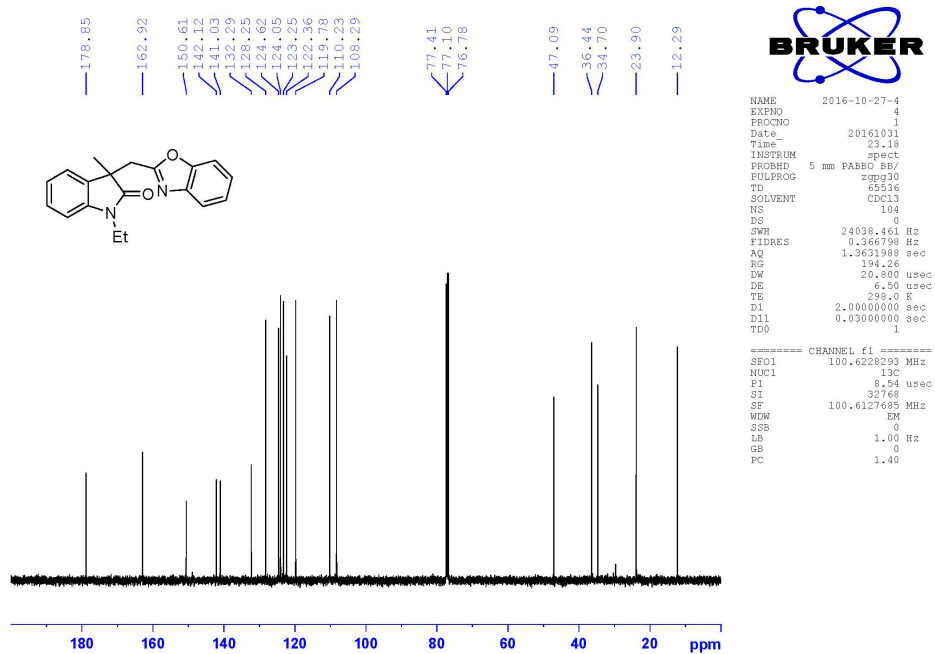
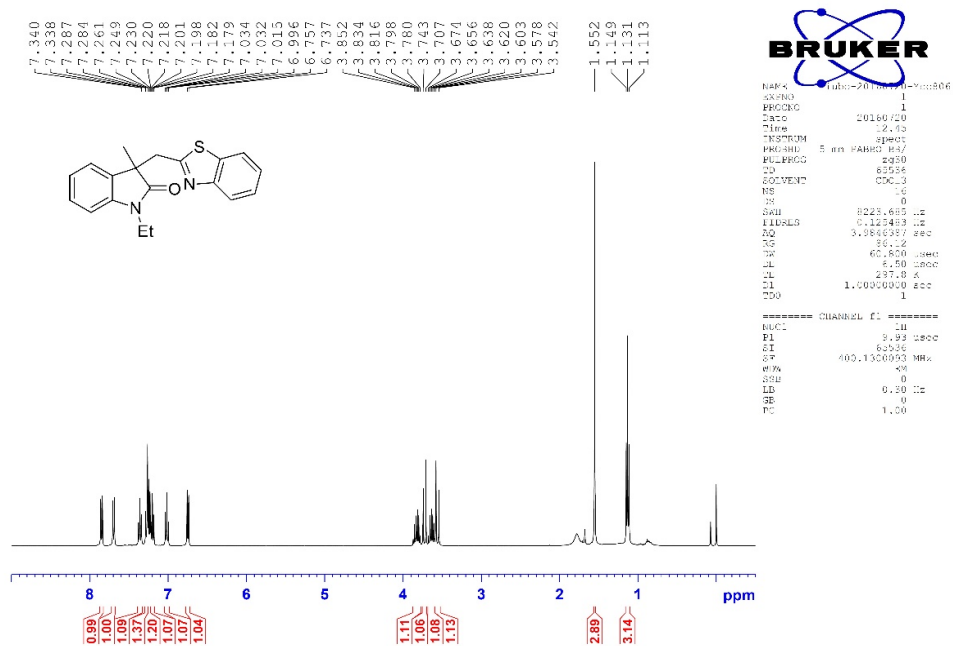
¹H and ¹³C NMR Spectra for Compound 2ed

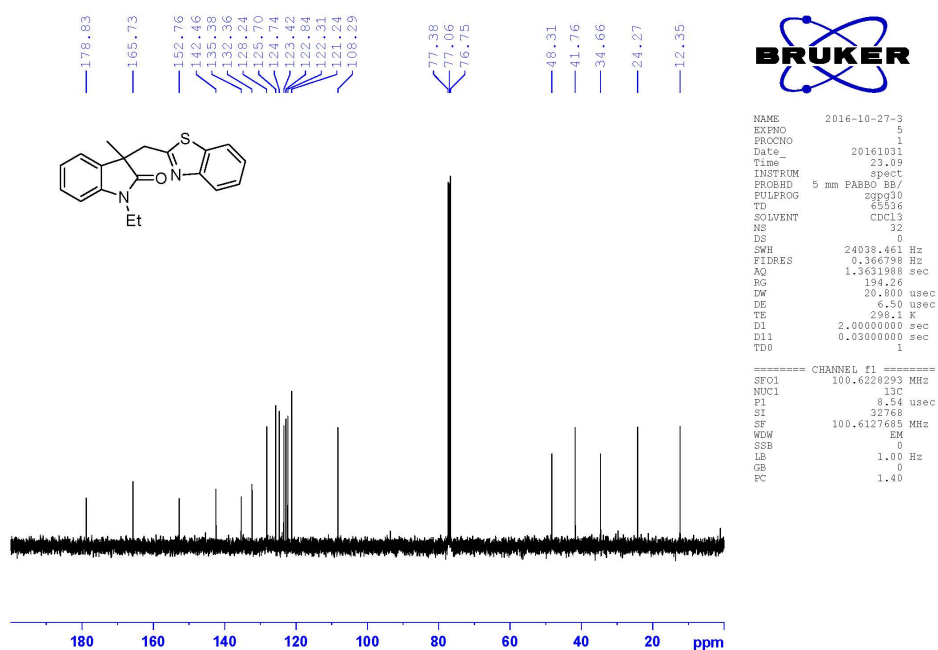
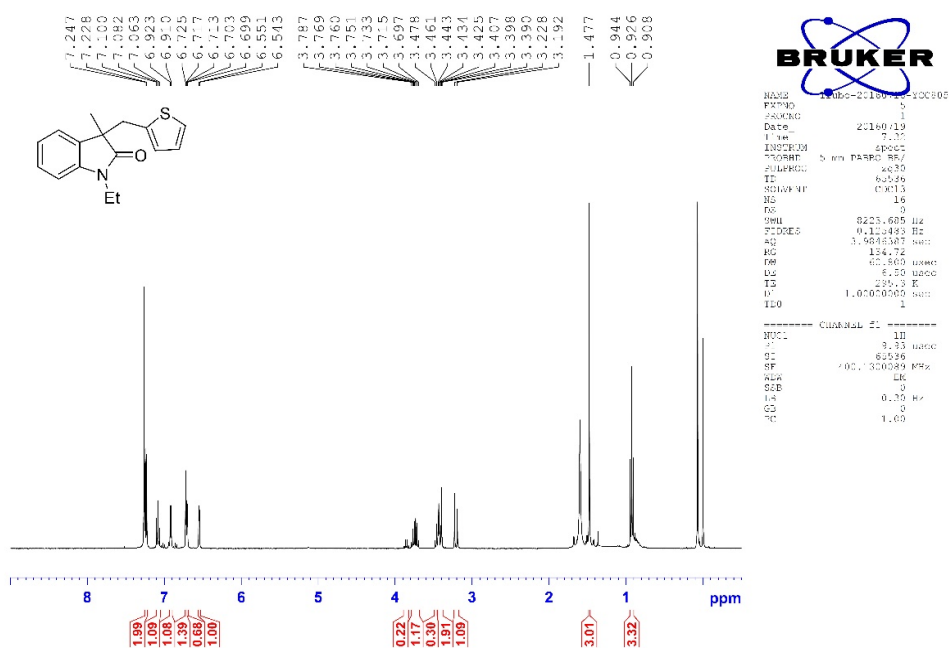
¹H and ¹³C NMR Spectra for Compound 2ee

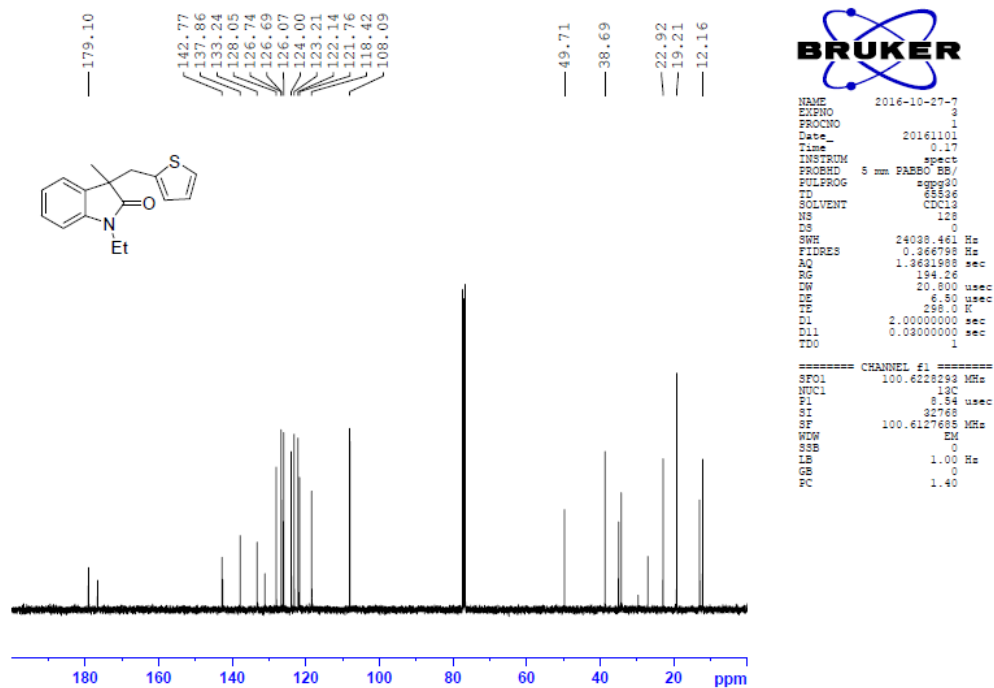
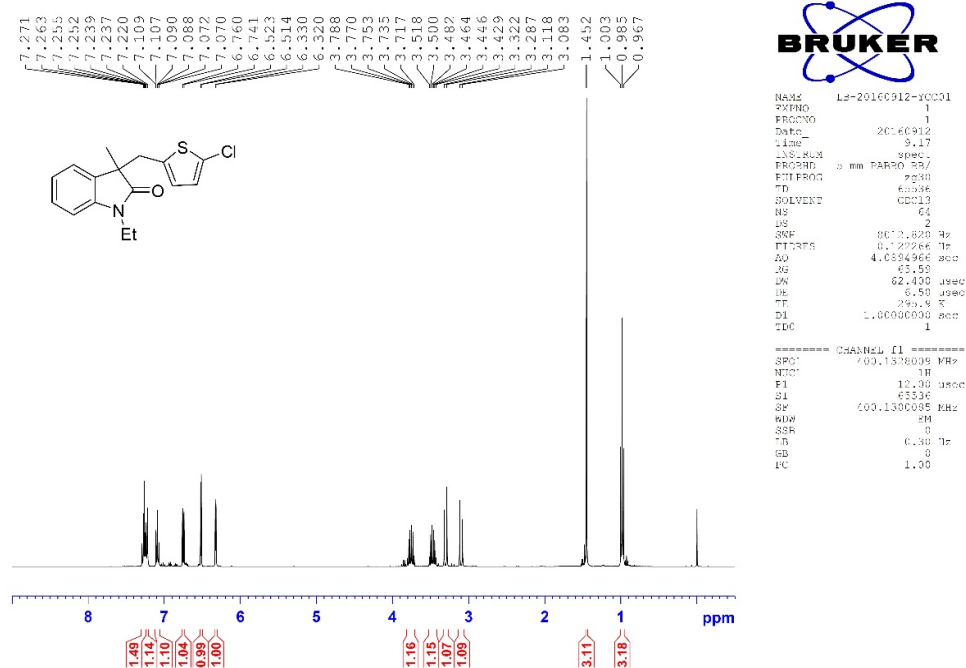
¹H and ¹³C NMR Spectra for Compound 2fa

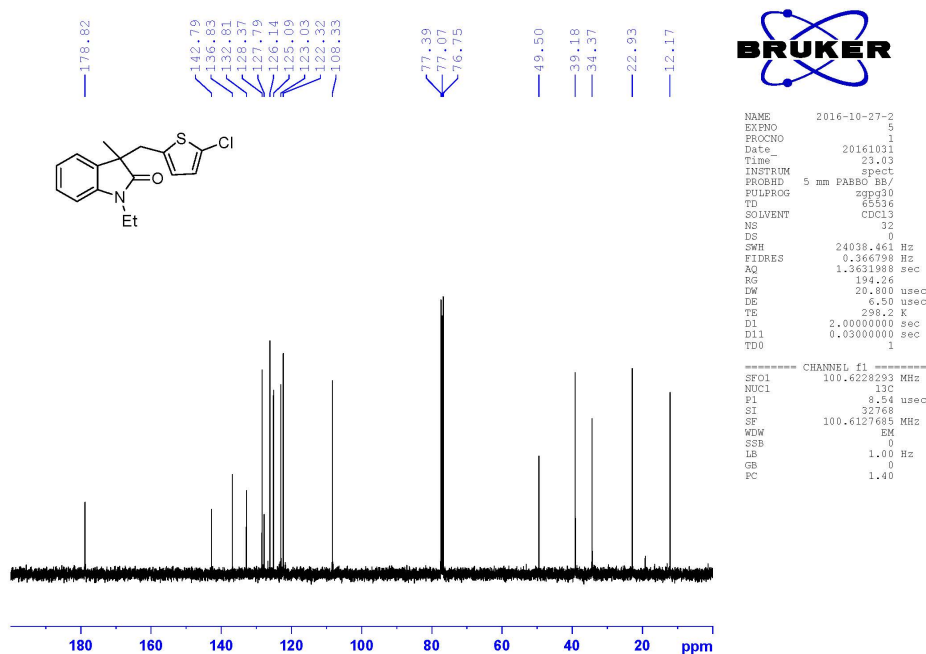
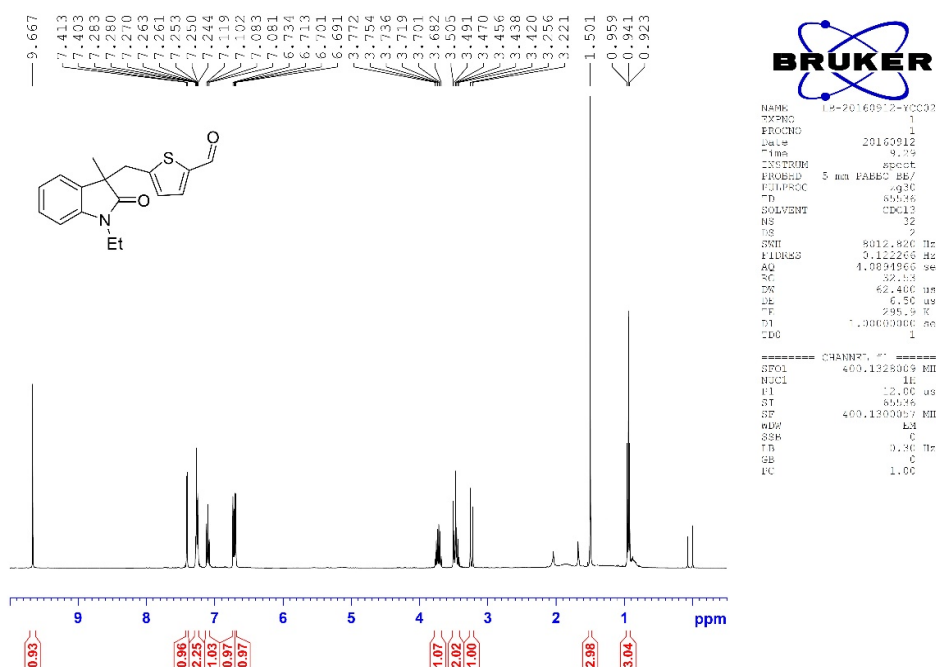
¹H and ¹³C NMR Spectra for Compound 2fb

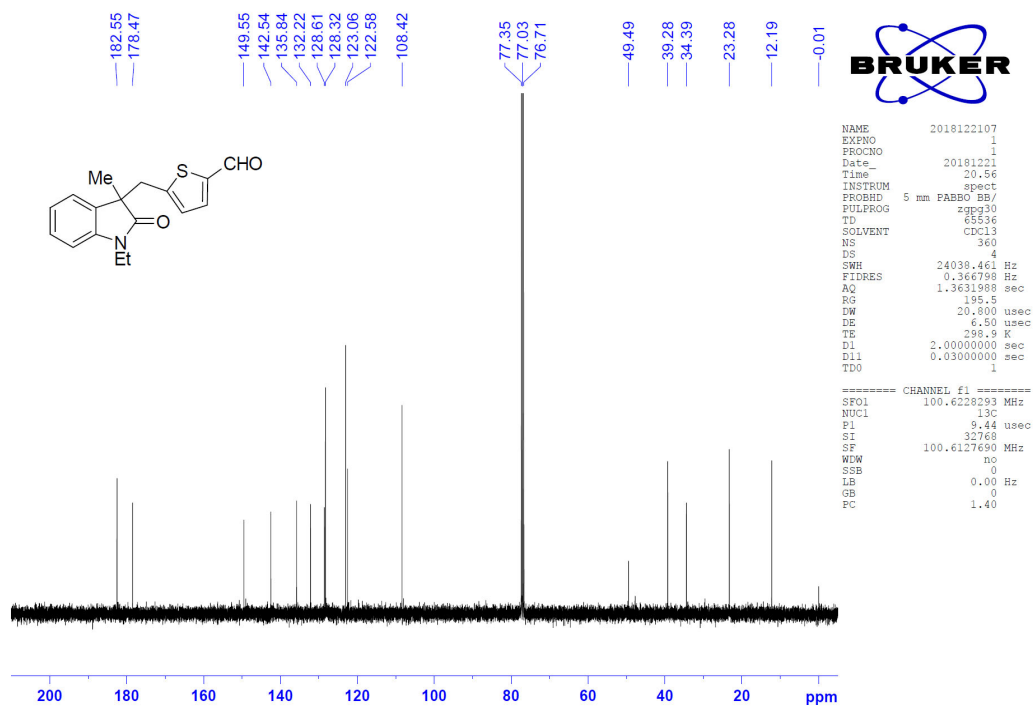
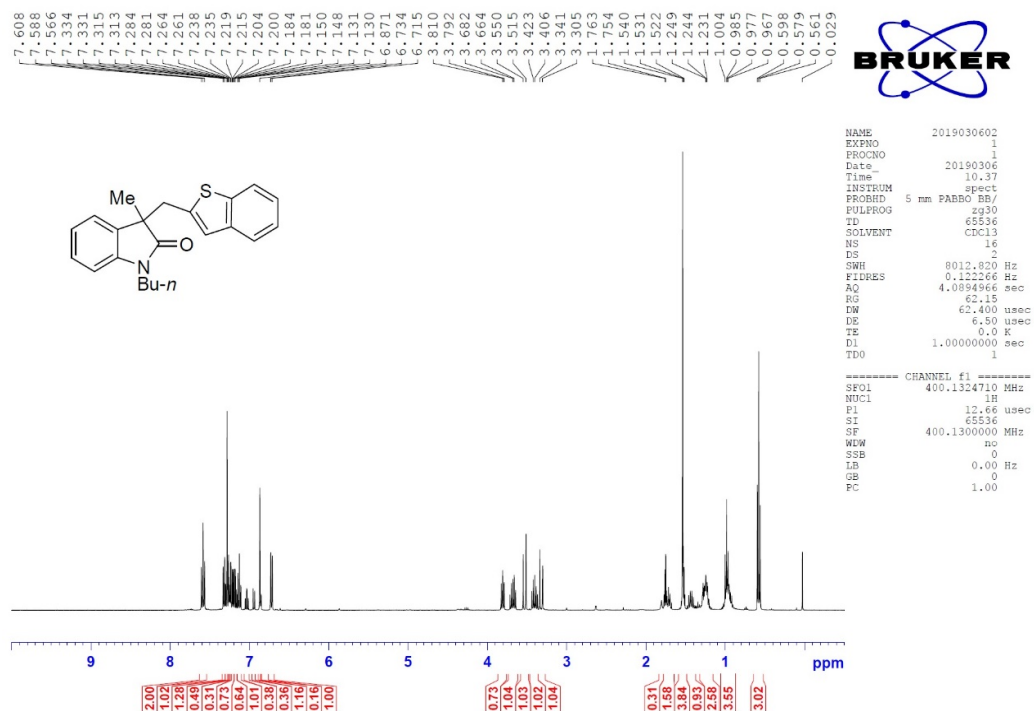
¹H and ¹³C NMR Spectra for Compound 2fc

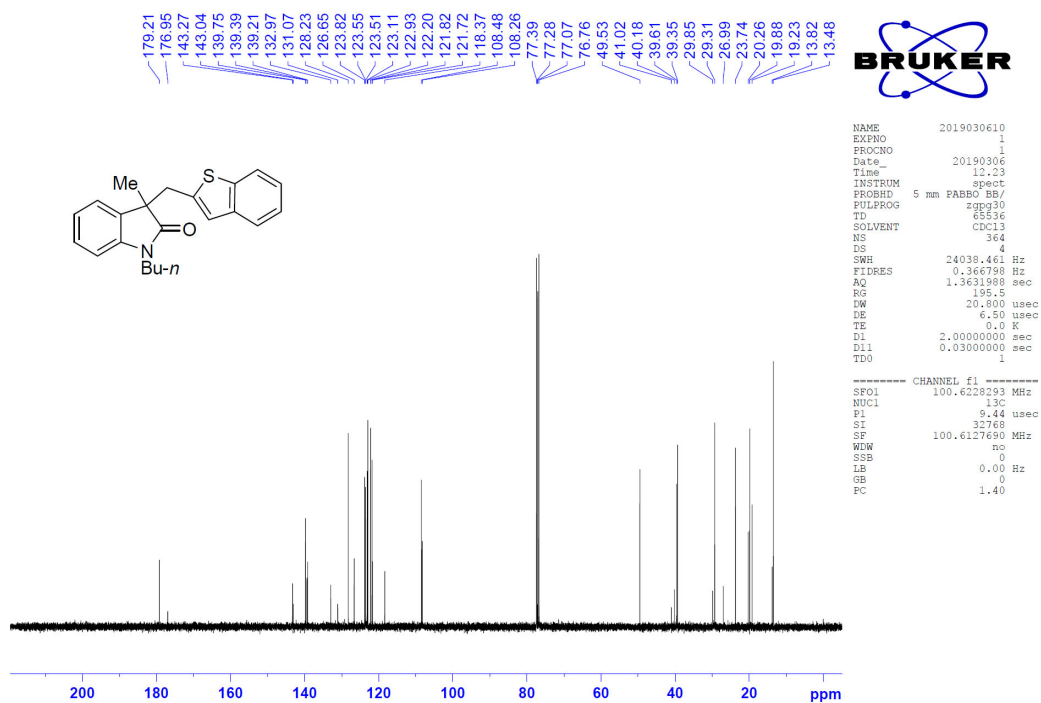
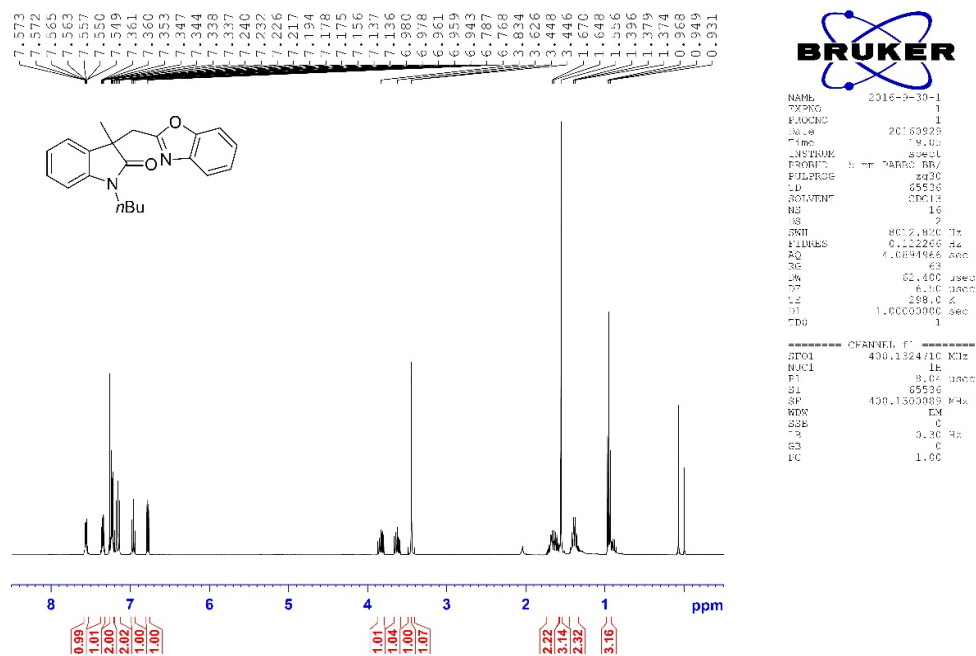
¹H and ¹³C NMR Spectra for Compound 2fd

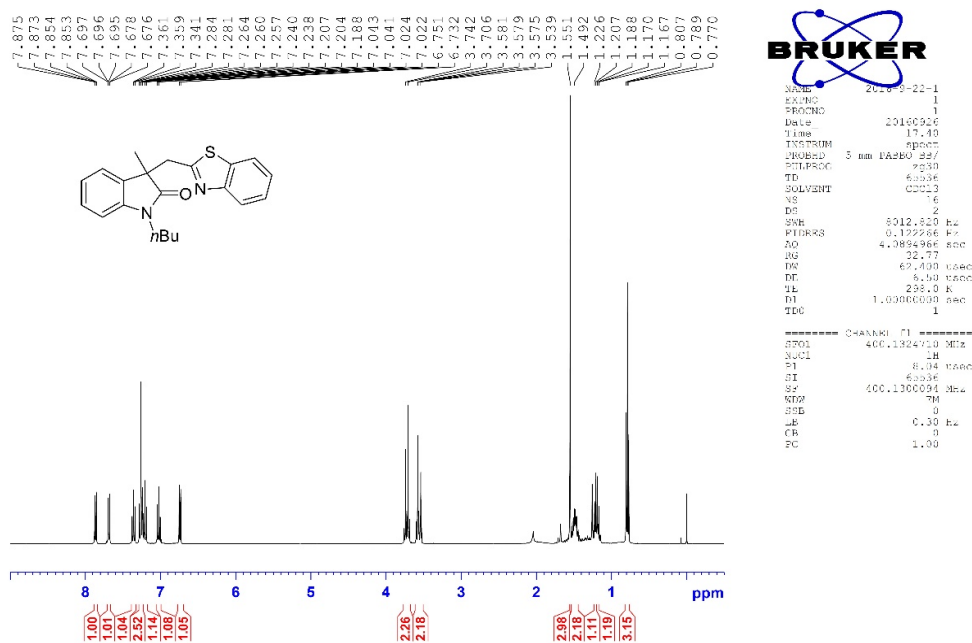
¹H and ¹³C NMR Spectra for Compound 2f

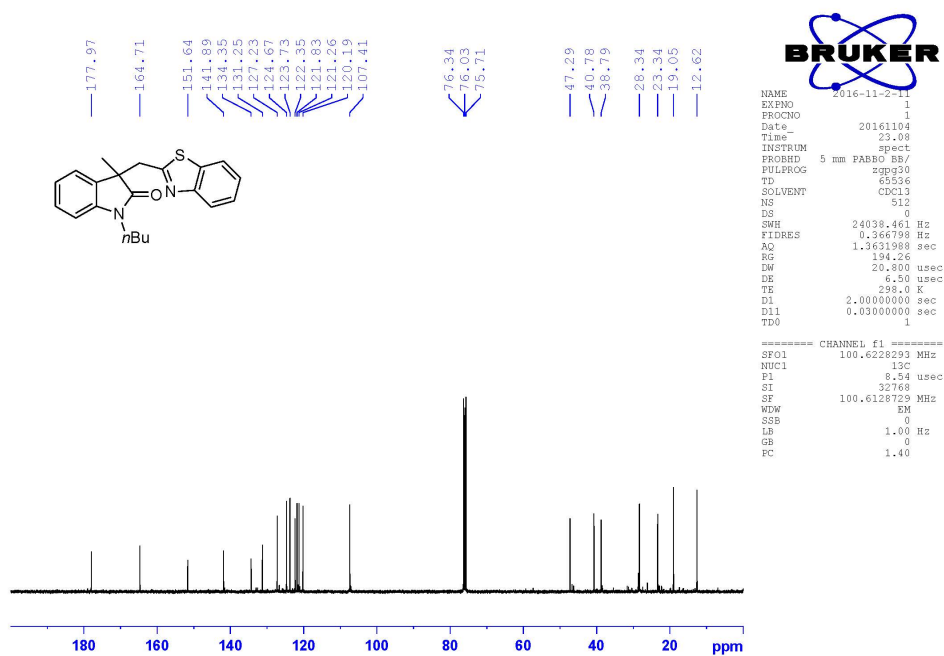
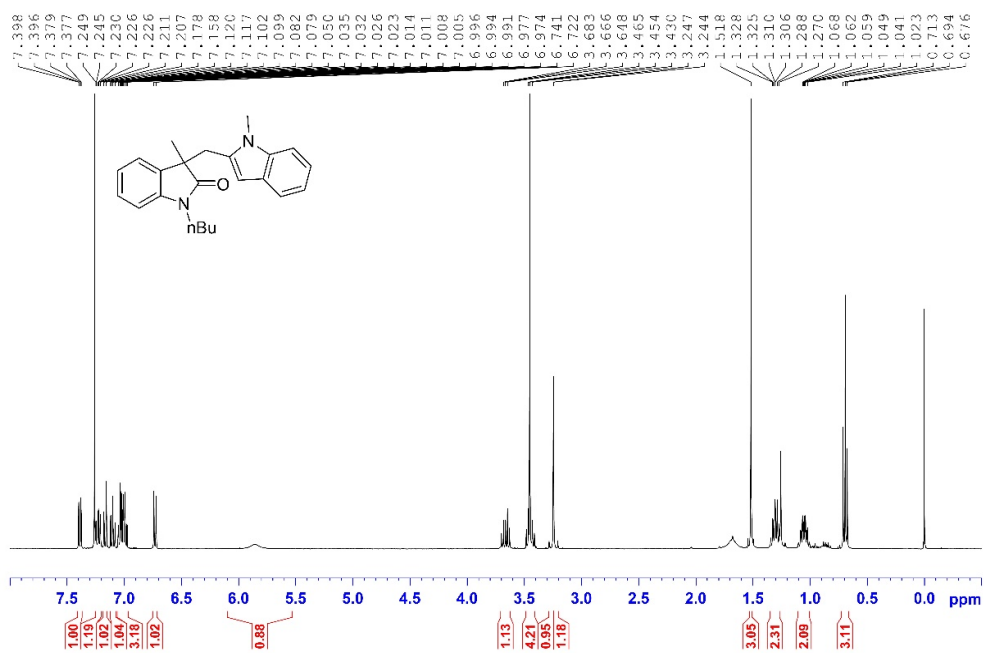
¹H and ¹³C NMR Spectra for Compound 2ff

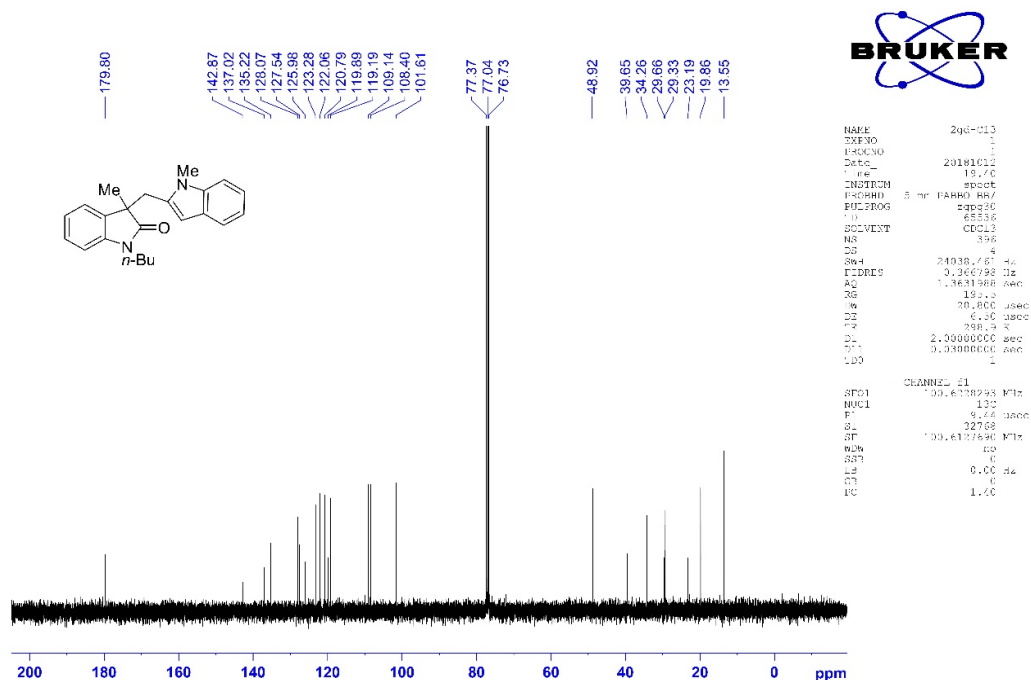
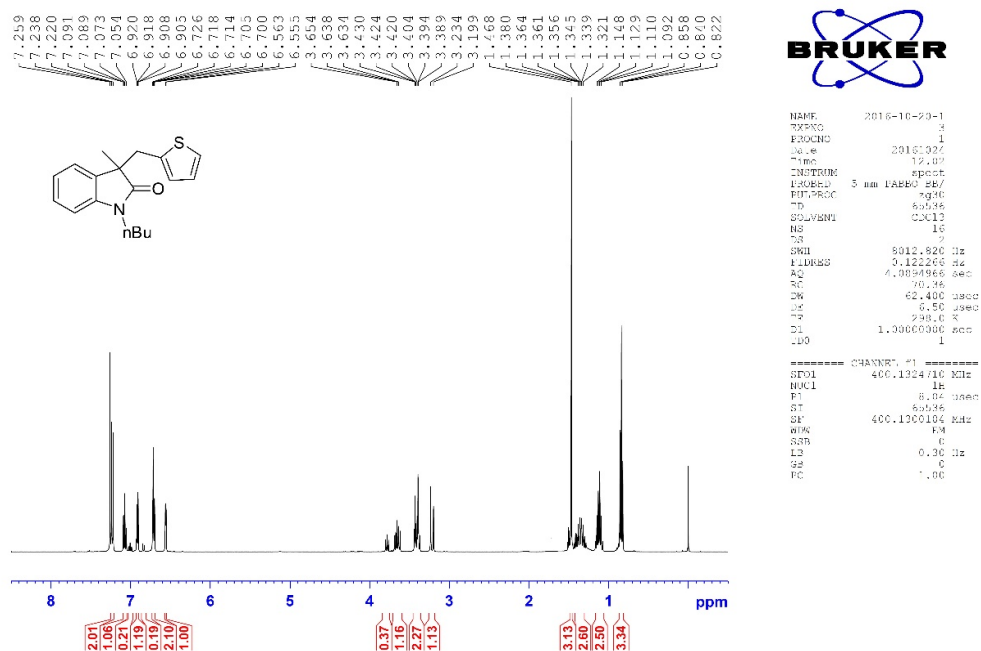
¹H and ¹³C NMR Spectra for Compound 2fg

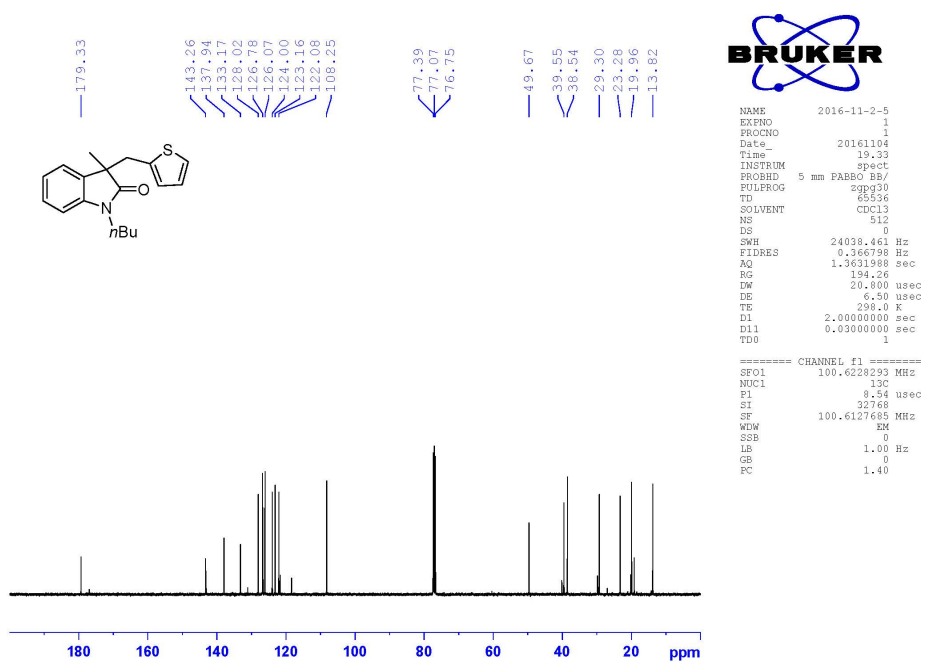
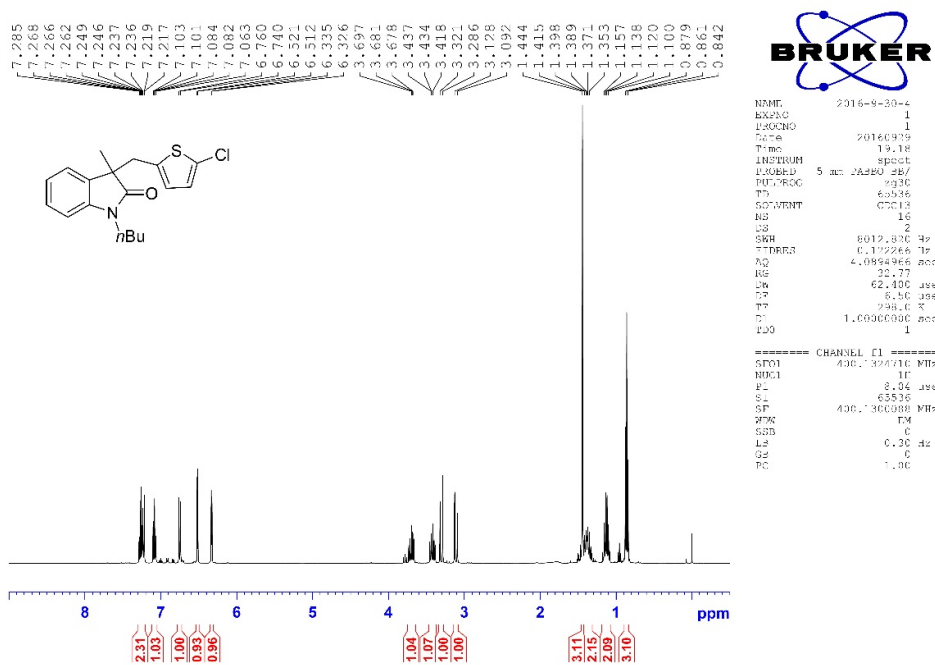
¹H and ¹³C NMR Spectra for Compound 2ga

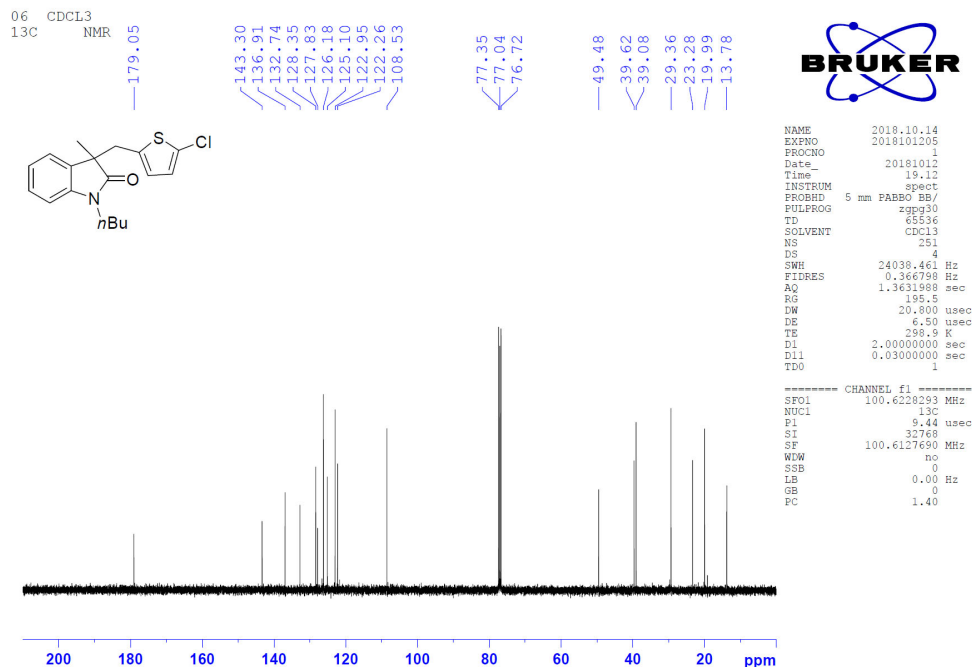
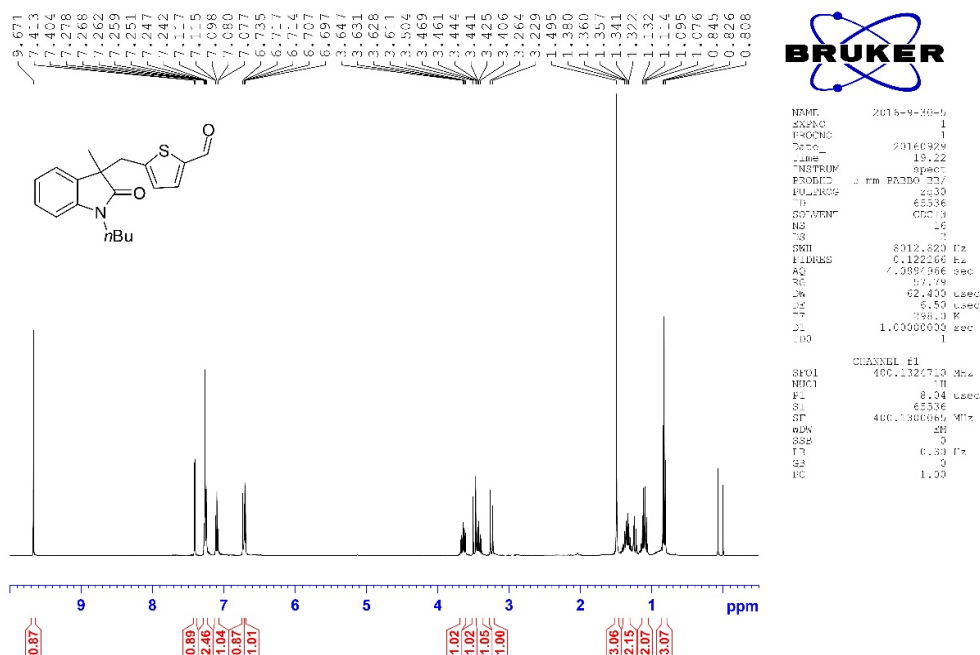
¹H and ¹³C NMR Spectra for Compound 2gb

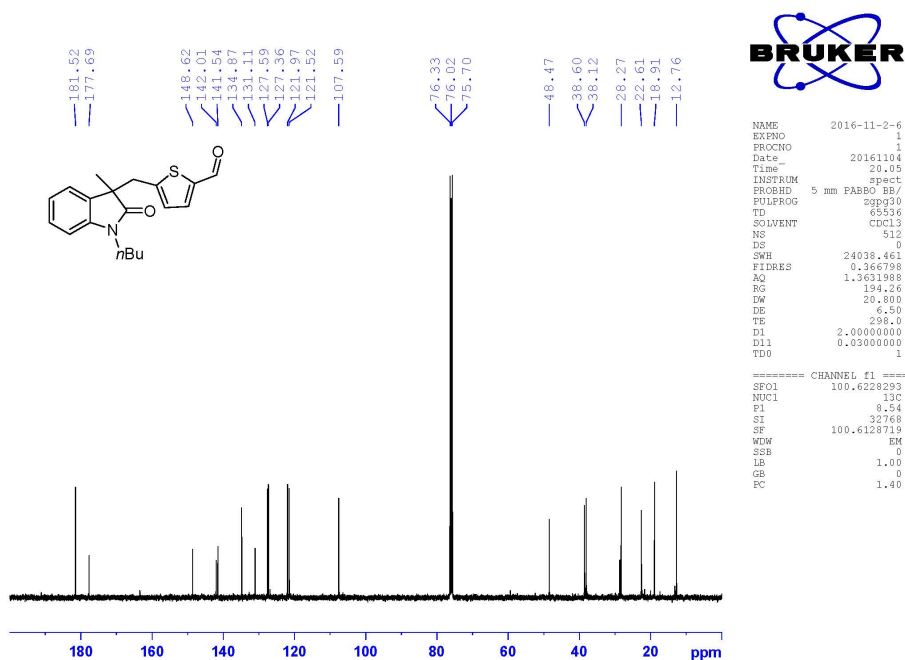
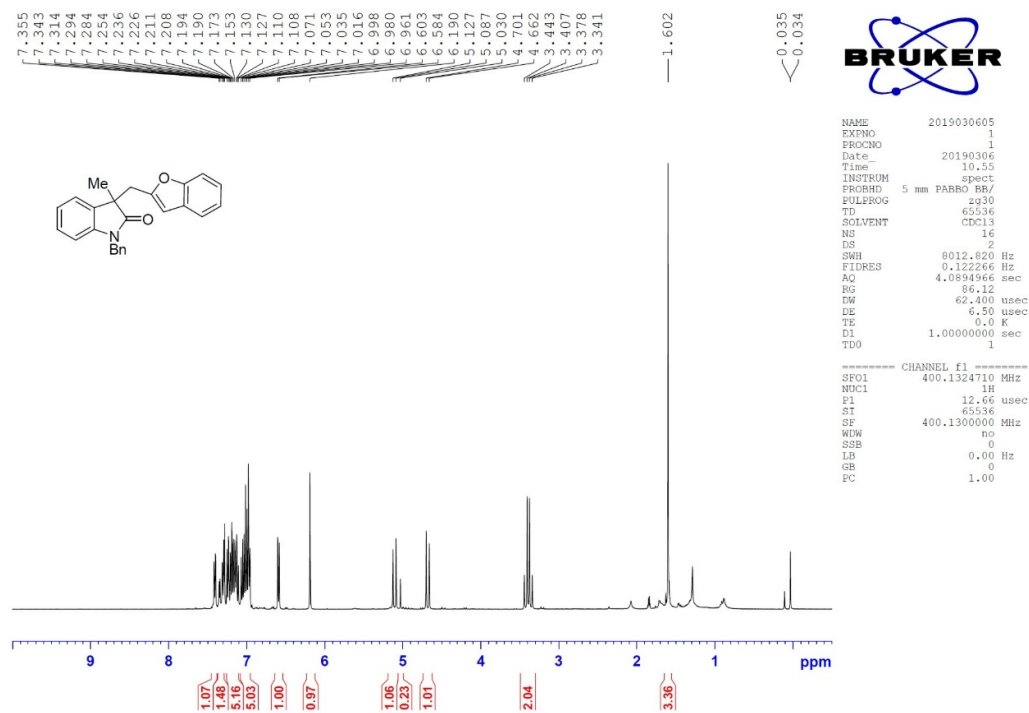


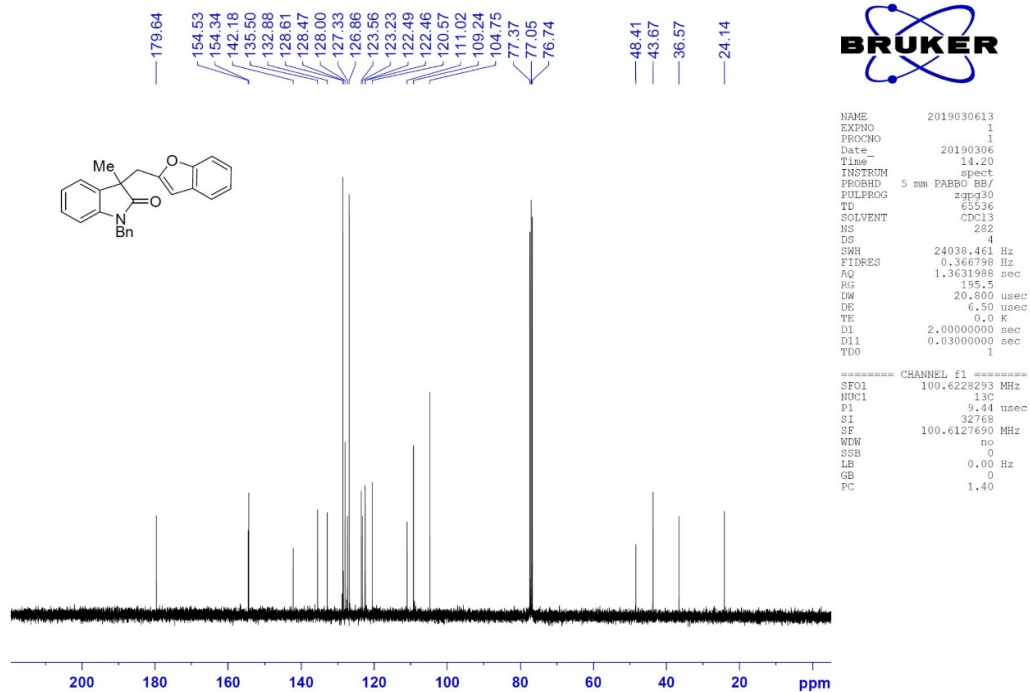
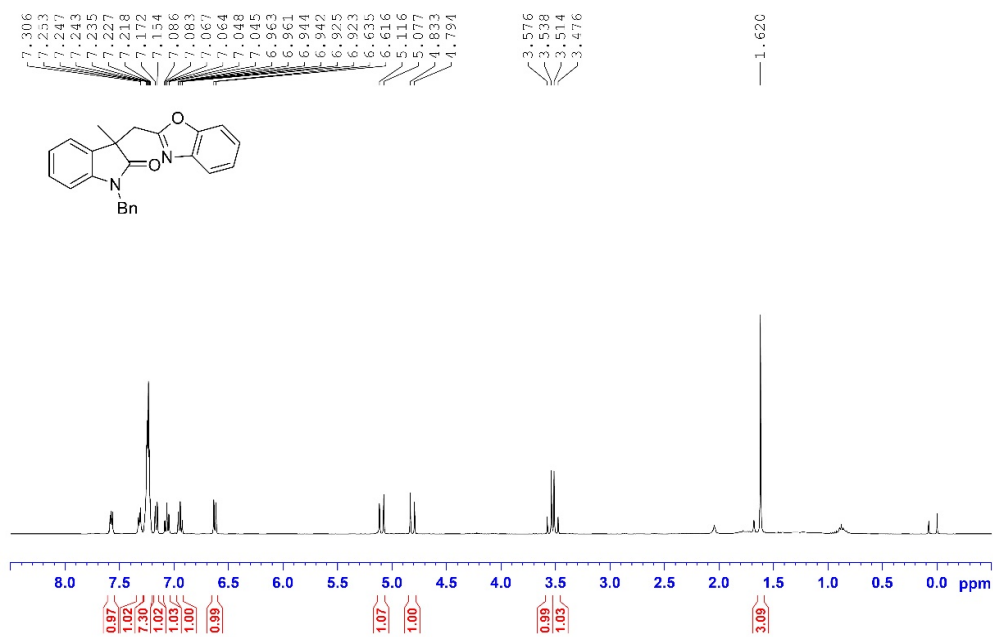
¹H and ¹³C NMR Spectra for Compound 2gd

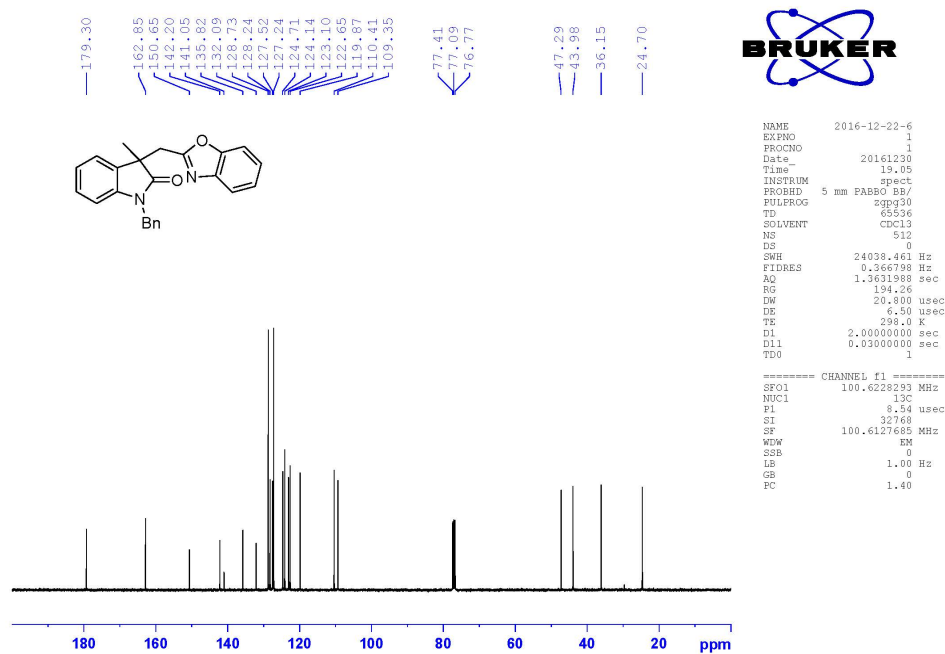
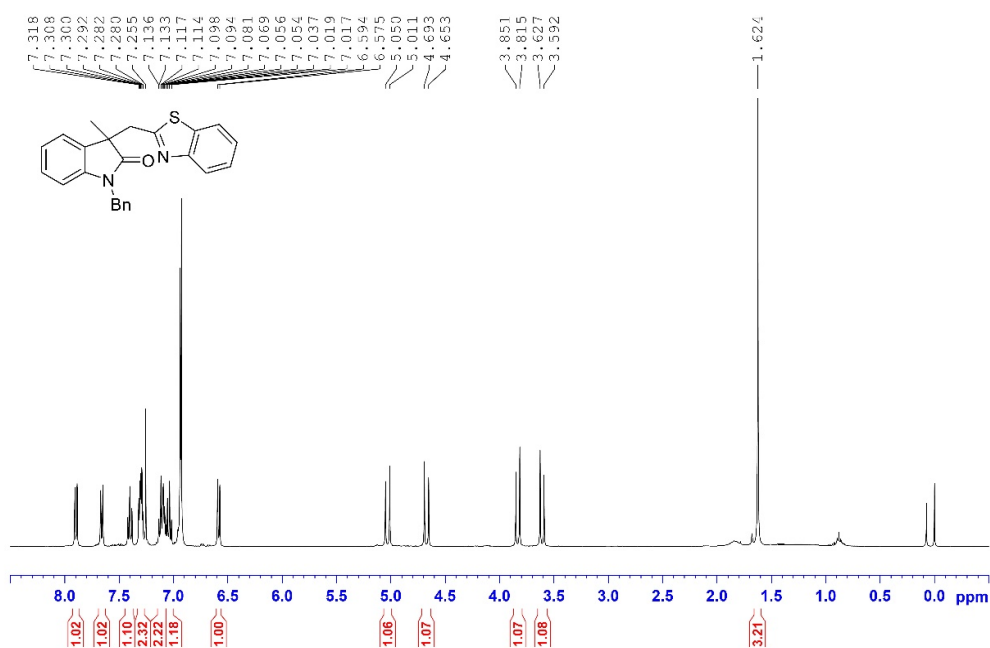
¹H and ¹³C NMR Spectra for Compound 2ge

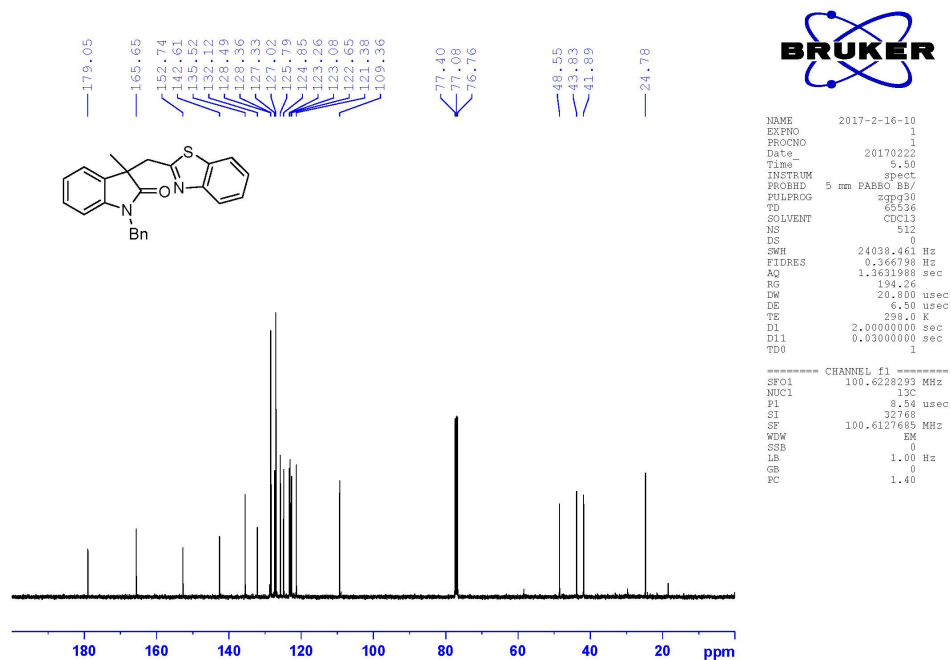
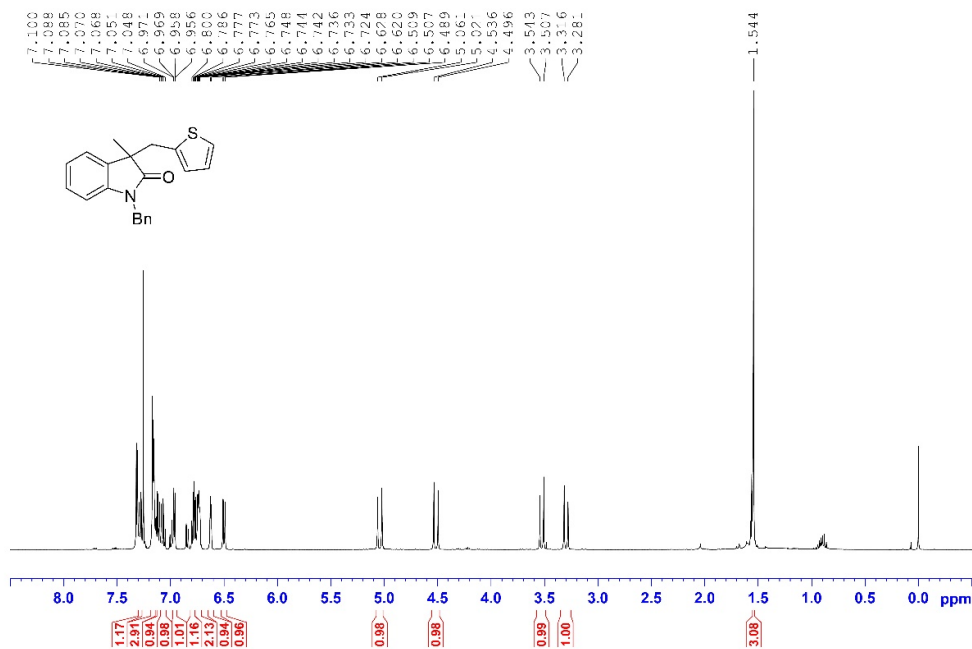
¹H and ¹³C NMR Spectra for Compound 2gf

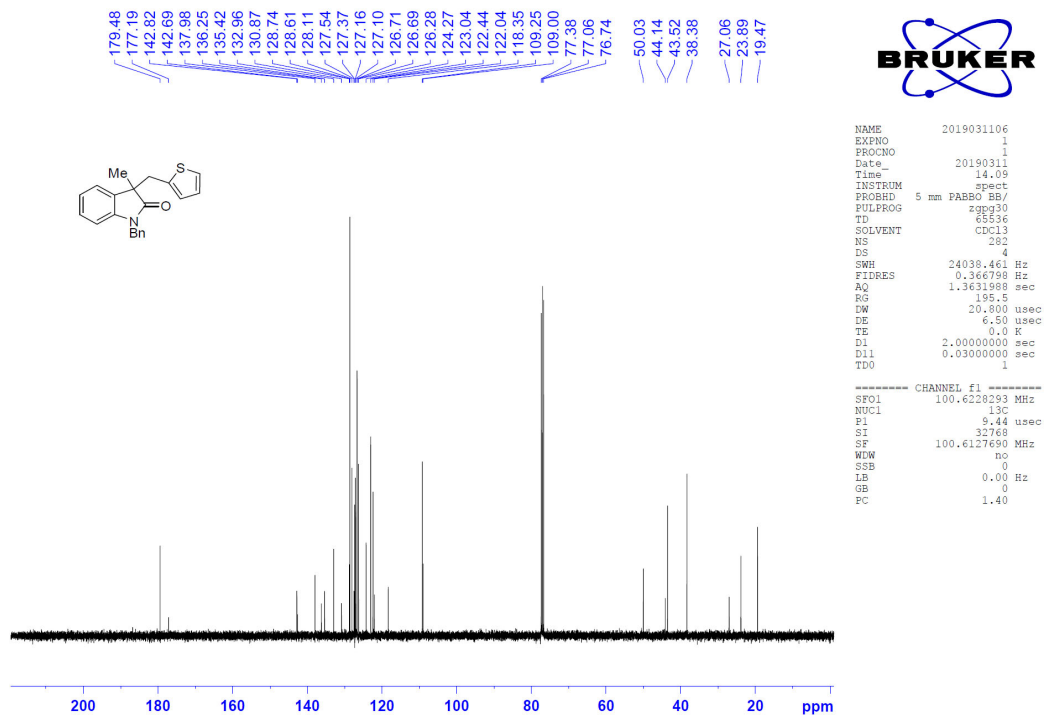
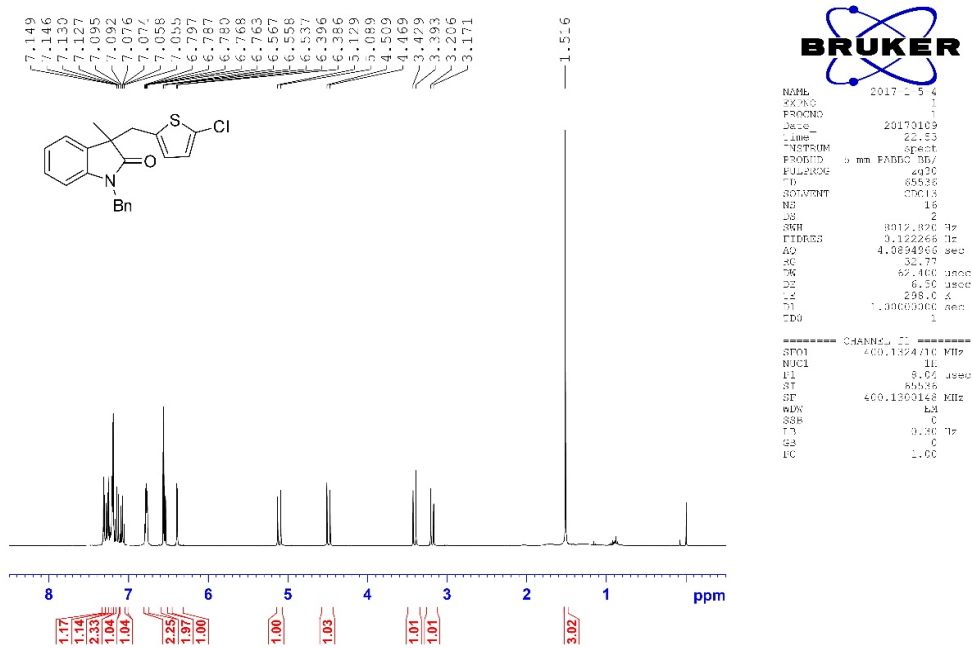
¹H and ¹³C NMR Spectra for Compound 2gg

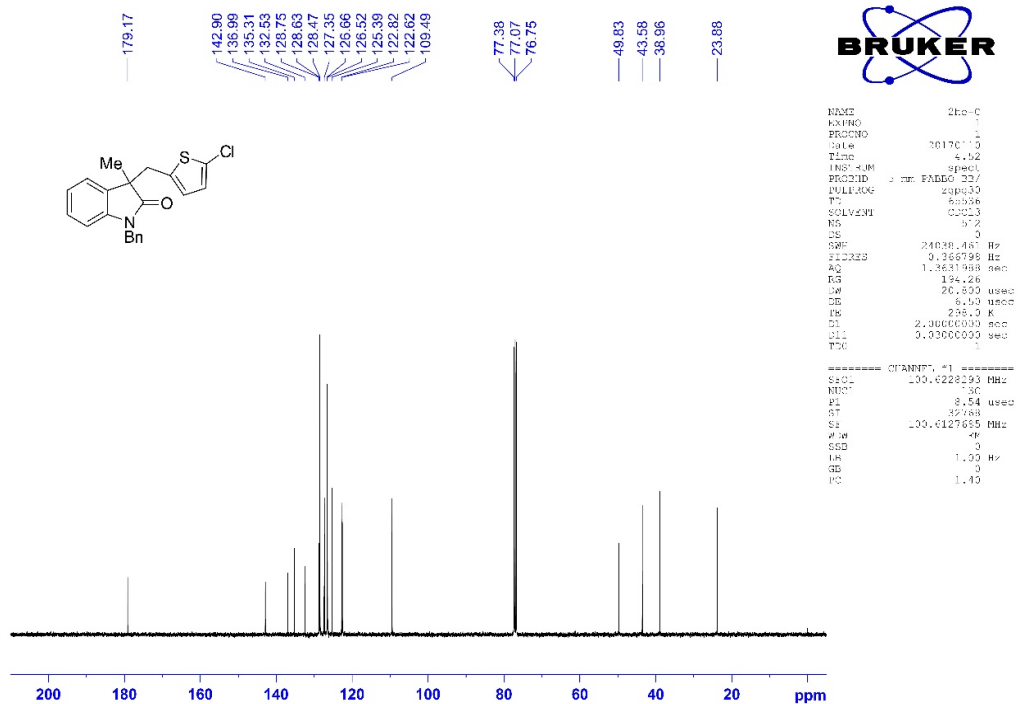
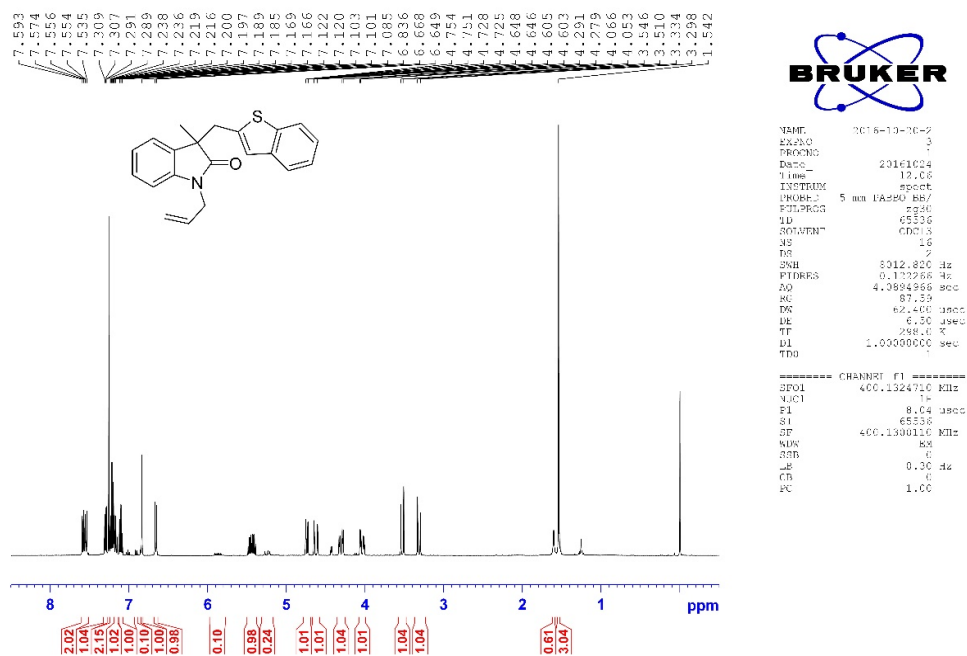
¹H and ¹³C NMR Spectra for Compound 2ha

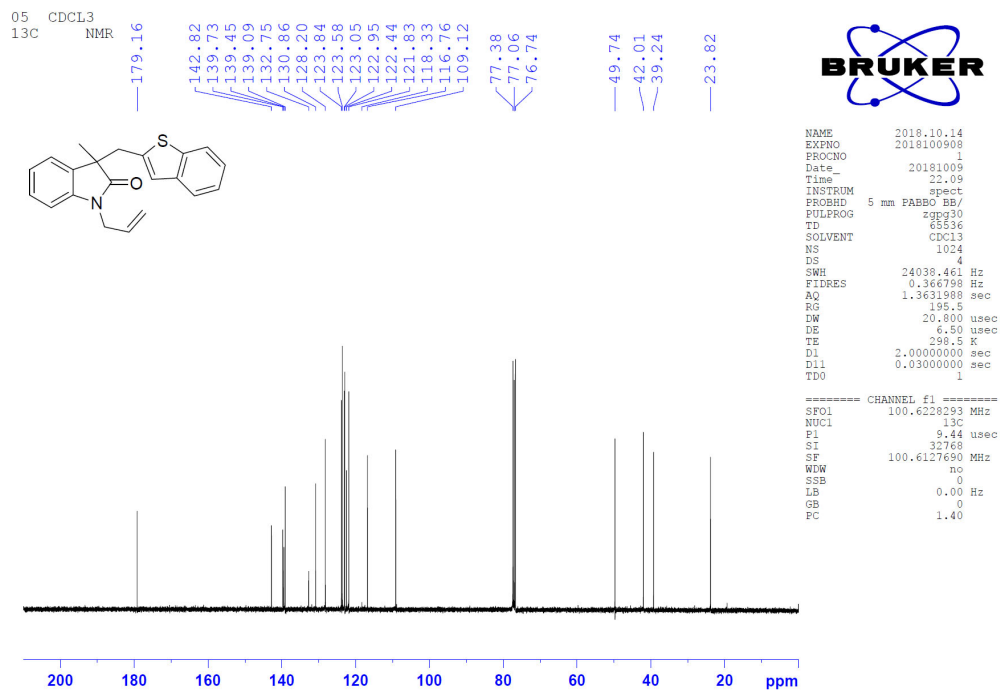
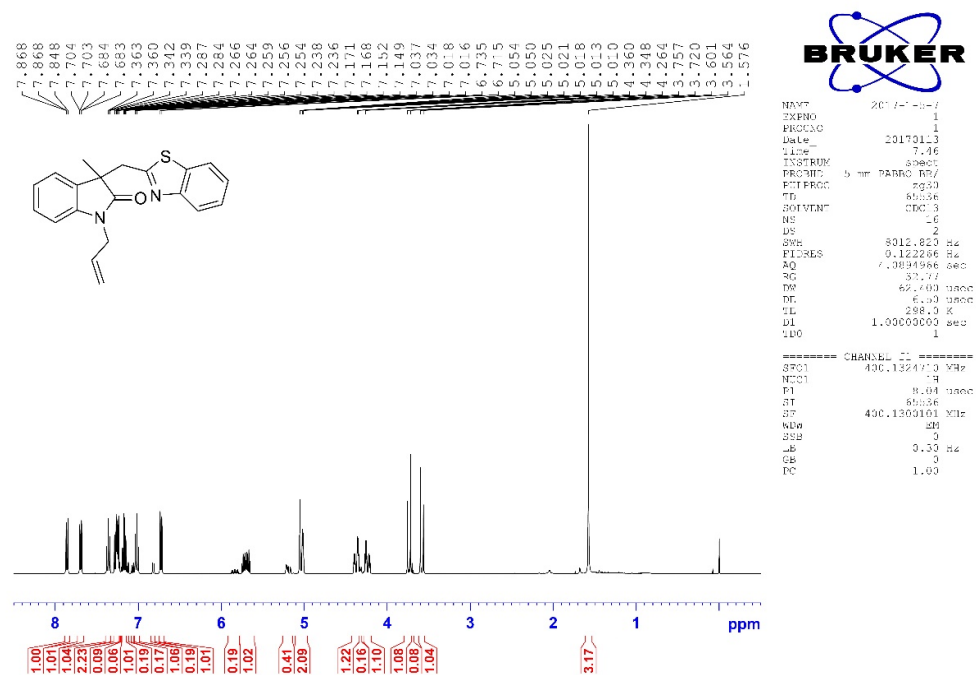
¹H and ¹³C NMR Spectra for Compound 2hb

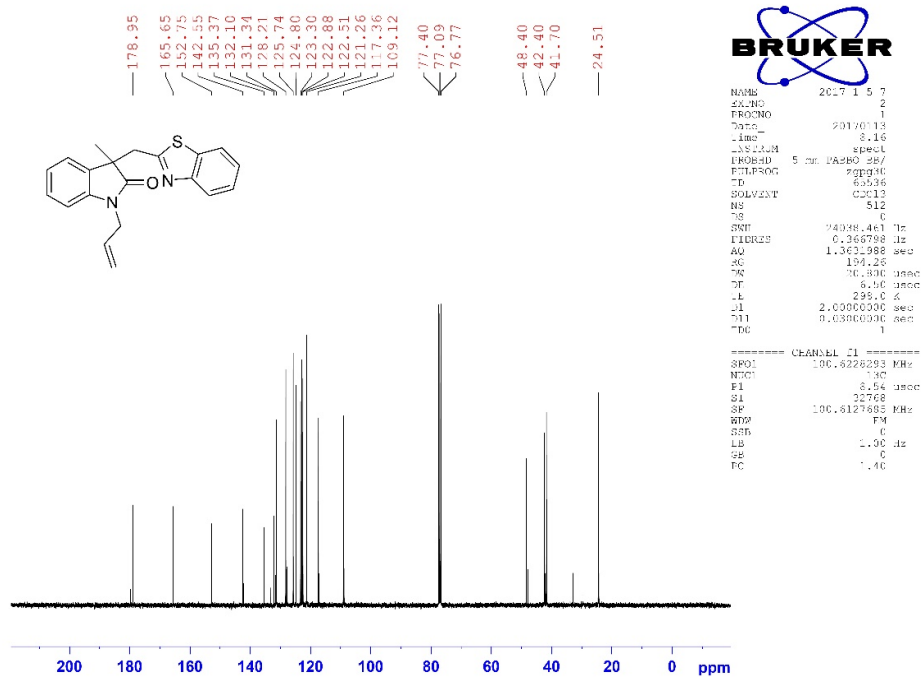
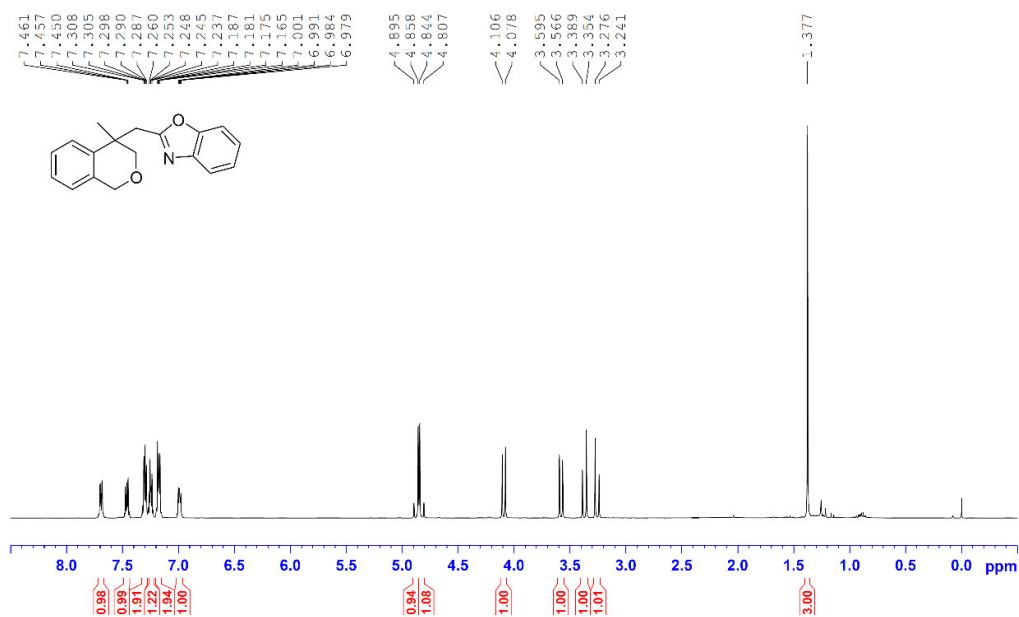
¹H and ¹³C NMR Spectra for Compound 2hc

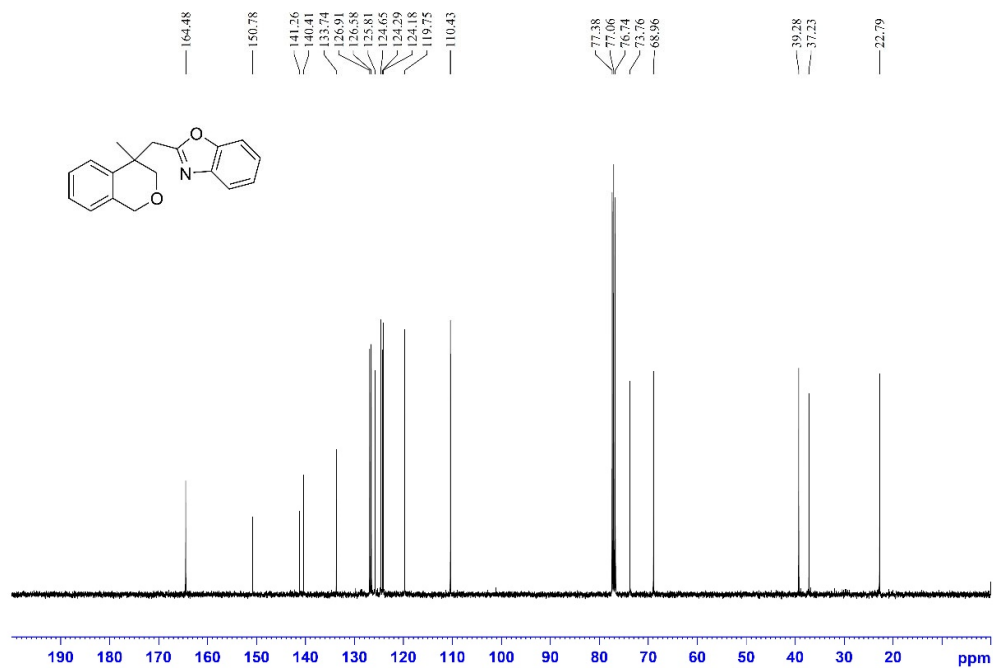
¹H and ¹³C NMR Spectra for Compound 2hd

¹H and ¹³C NMR Spectra for Compound 2he

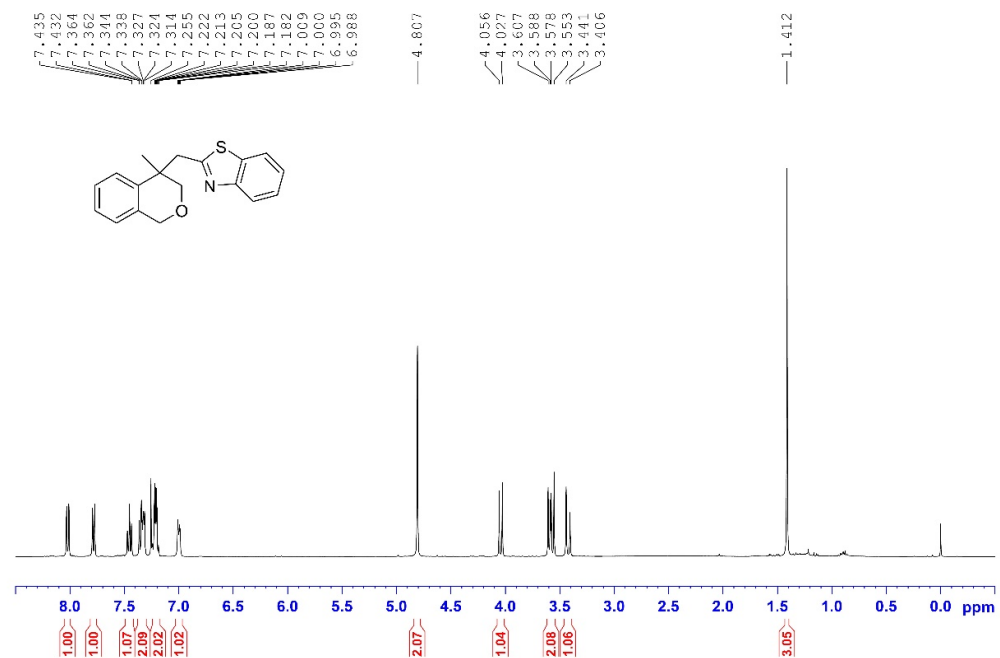
¹H and ¹³C NMR Spectra for Compound 2ia

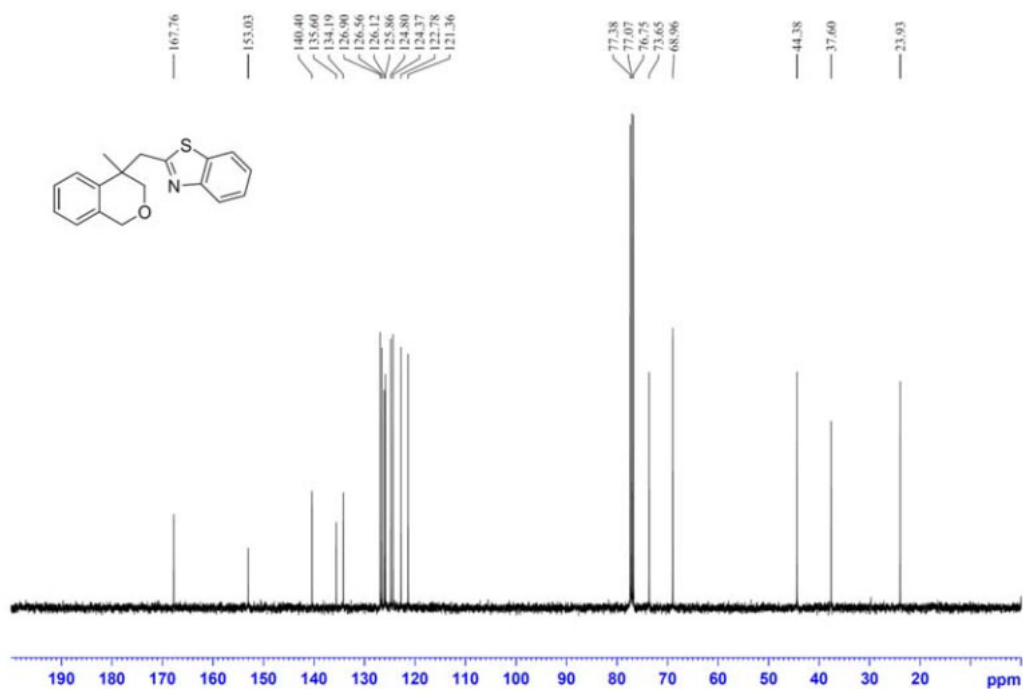
¹H and ¹³C NMR Spectra for Compound 2ib

 ^1H and ^{13}C NMR Spectra for Compound 2ja



¹H and ¹³C NMR Spectra for Compound **2jb**



¹H and ¹³C NMR Spectra for Compound 2jc