

Supporting Information for

Assembly of Indolenines, 3-Amino Oxindoles, and Aldehydes into

Indolenine-substituted Spiro[pyrrolidin-2,3'-oxindoles] via

1,3-Dipolar Cycloaddition with Divergent Diastereoselectivities

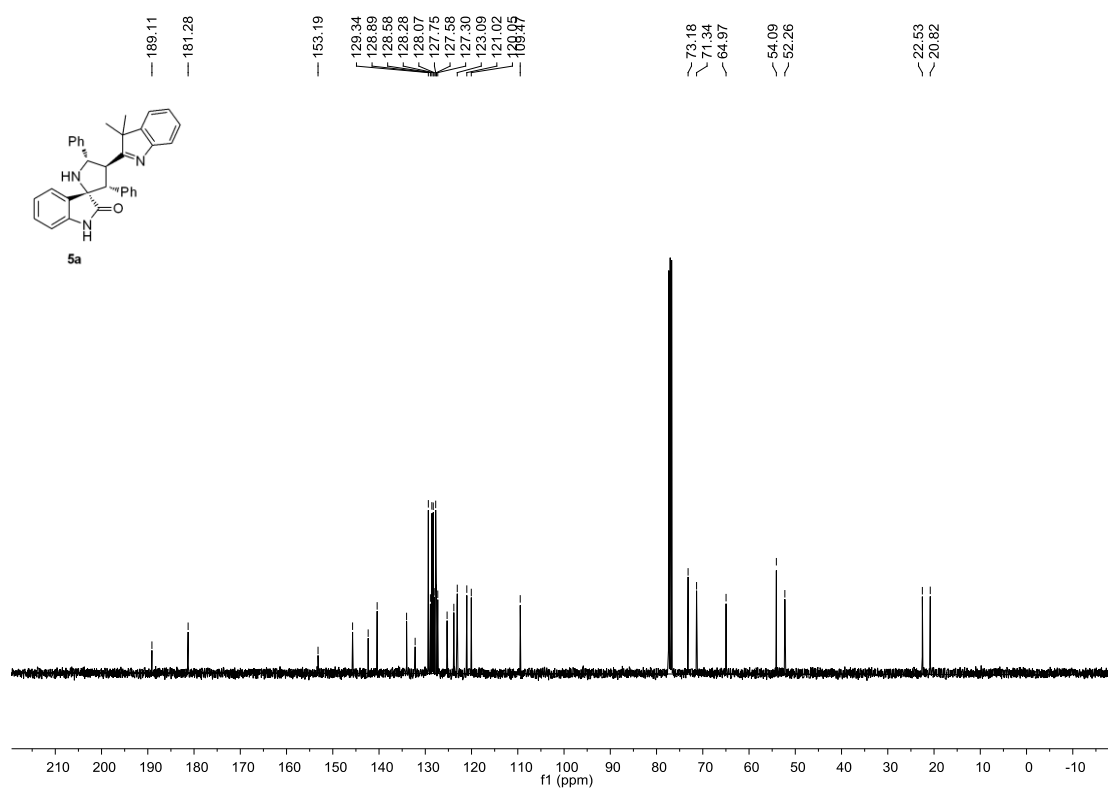
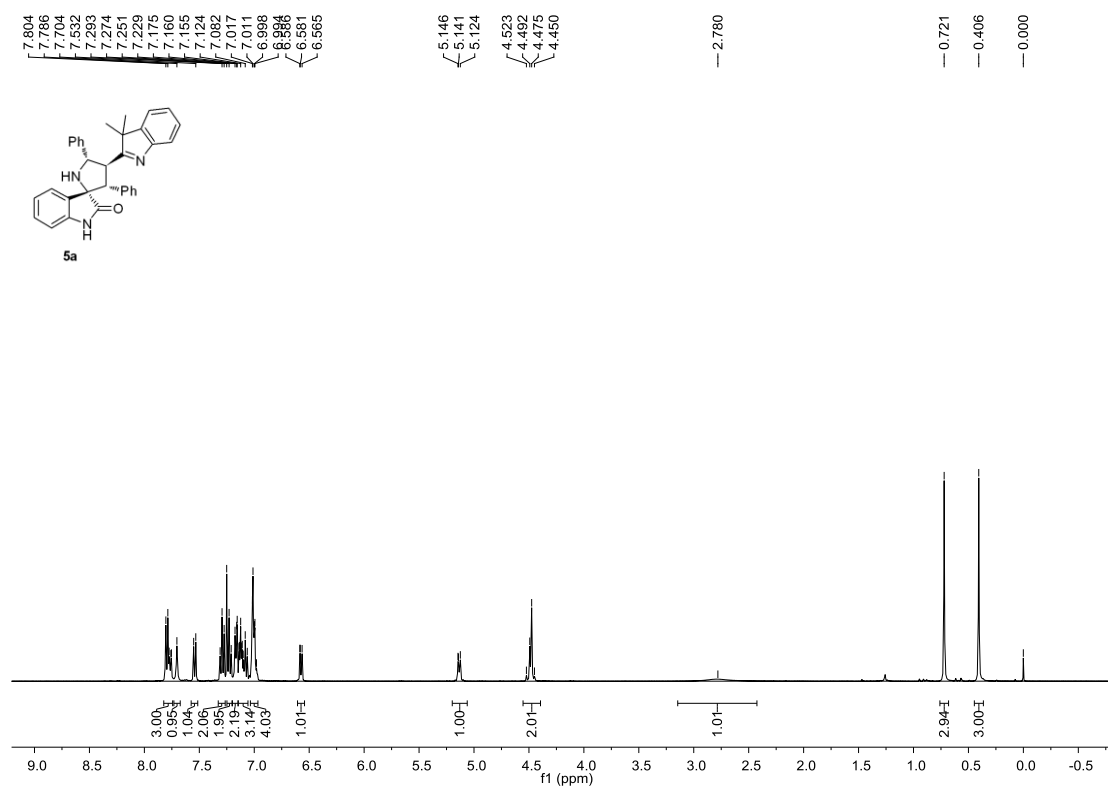
Guodong Zhu, Siyuan Liu, Shiqi Wu, Lianghong Peng, Jingping Qu, and Baomin Wang*

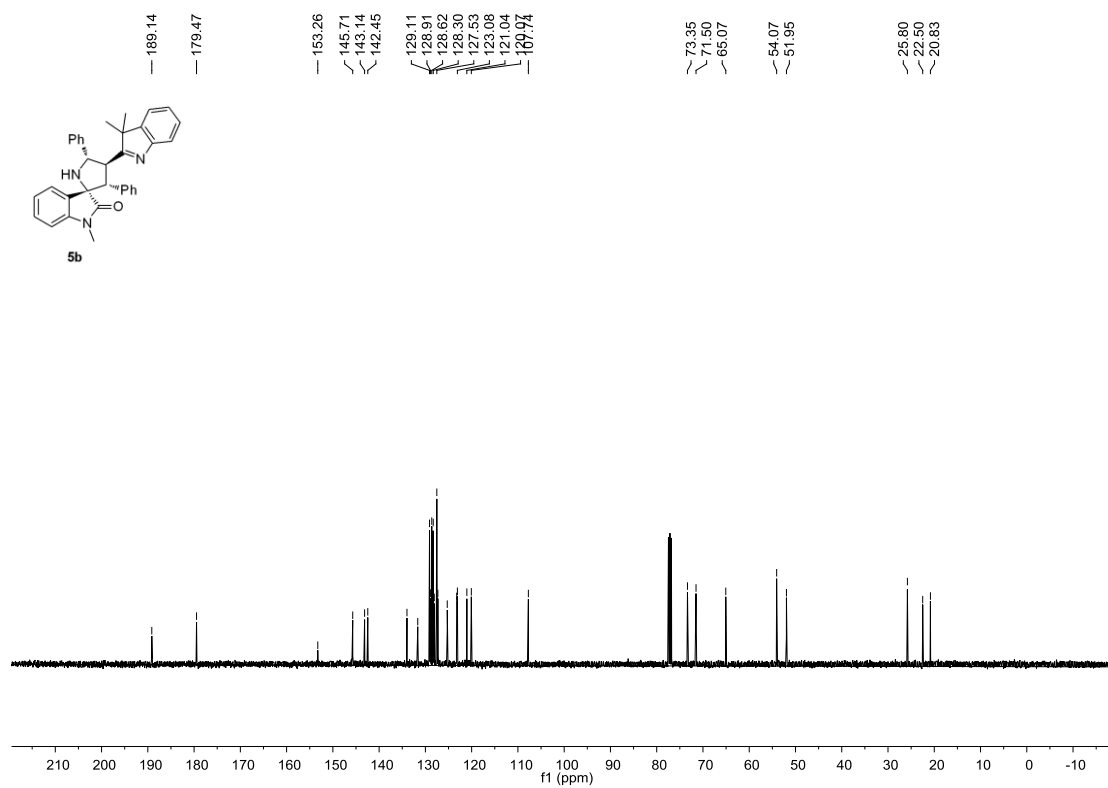
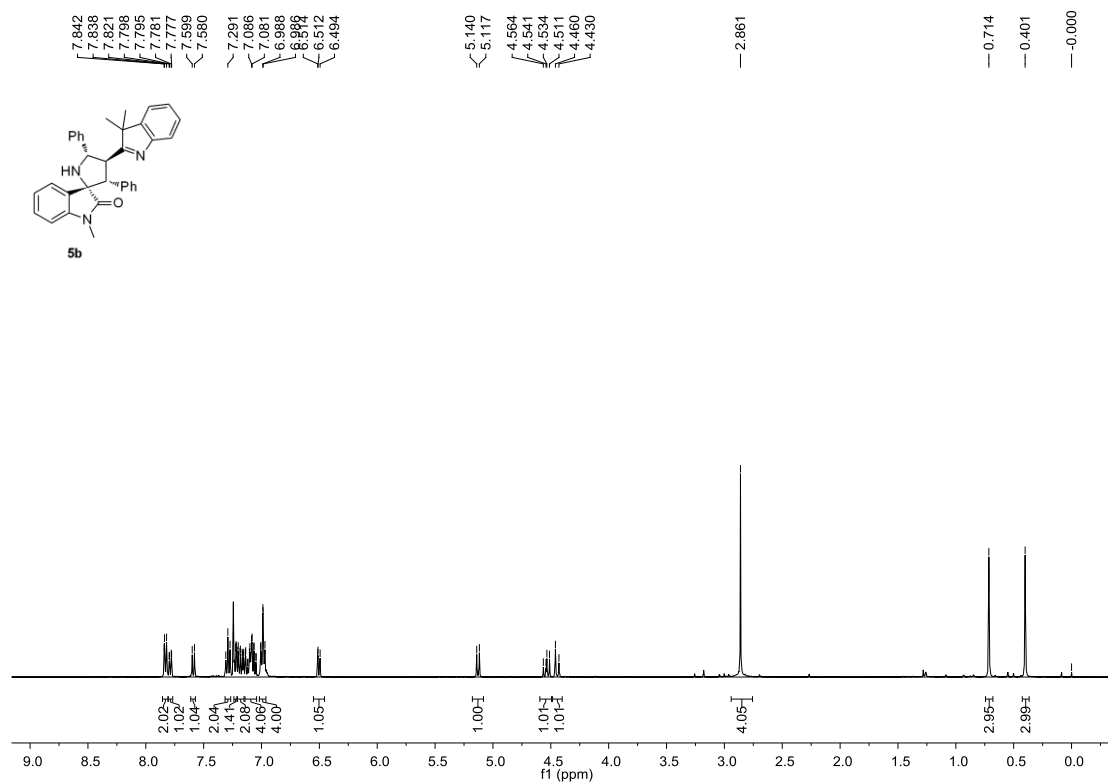
State Key Laboratory of Fine Chemicals, School of Pharmaceutical Science and Technology, Dalian University of Technology, Dalian 116024, P. R. China
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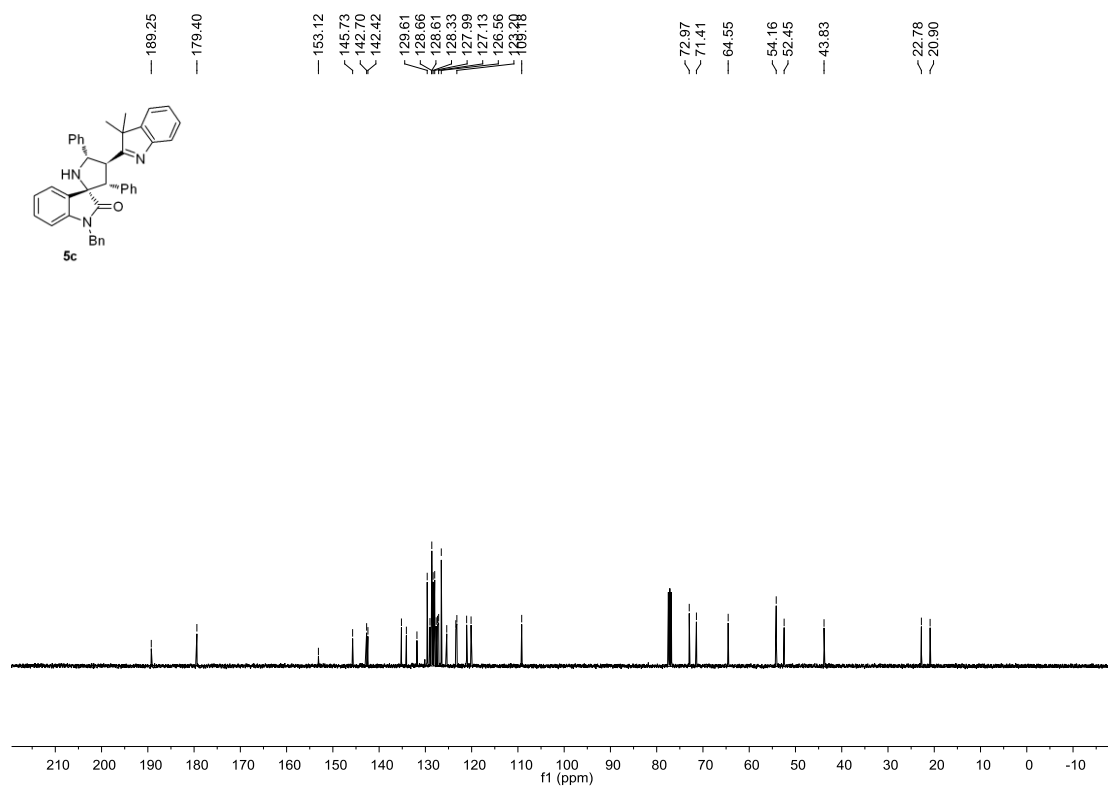
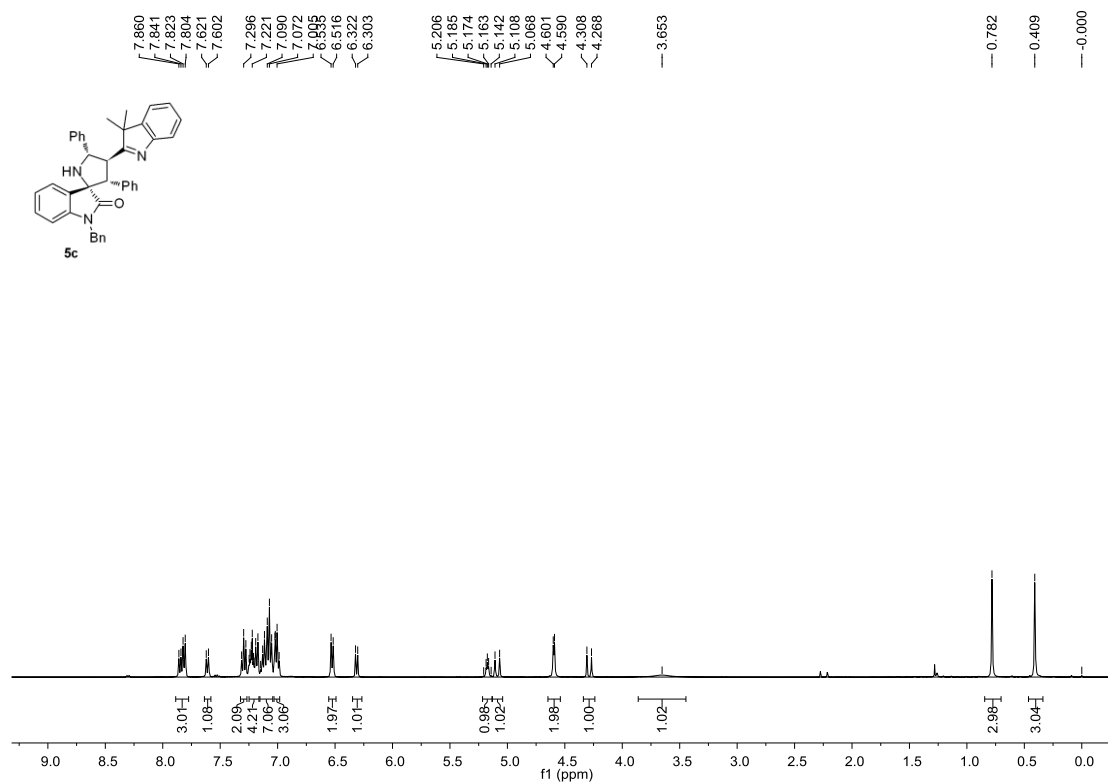
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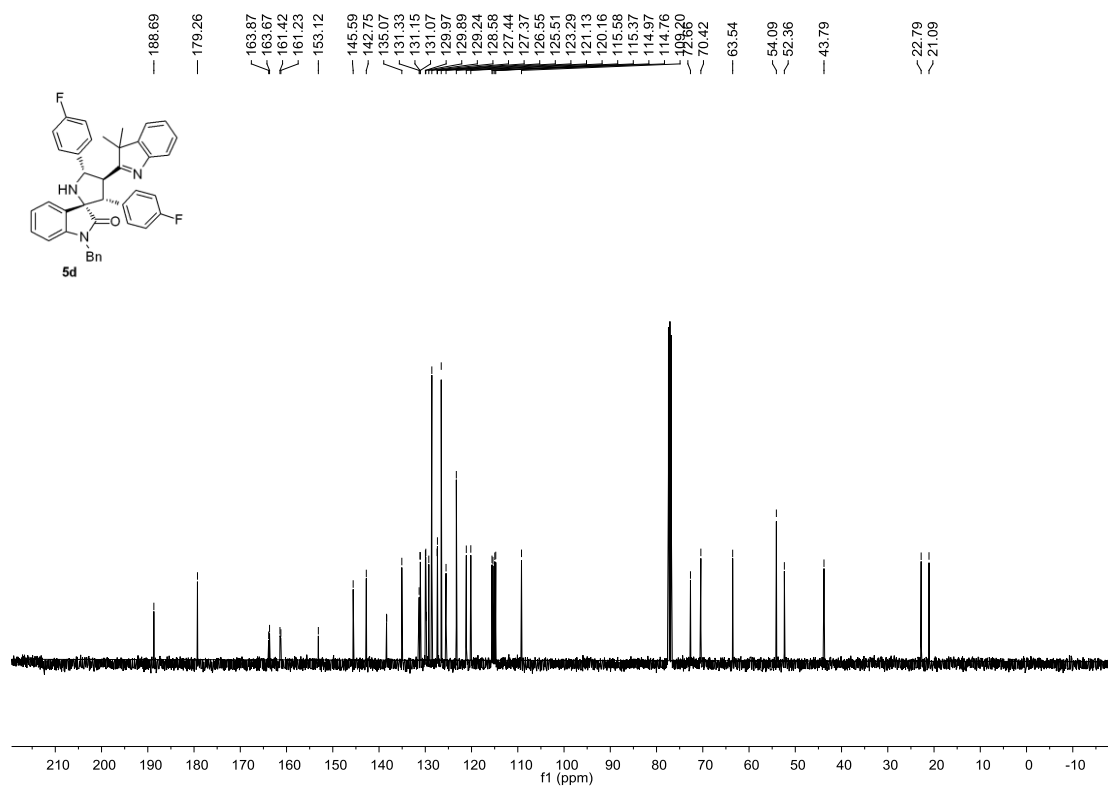
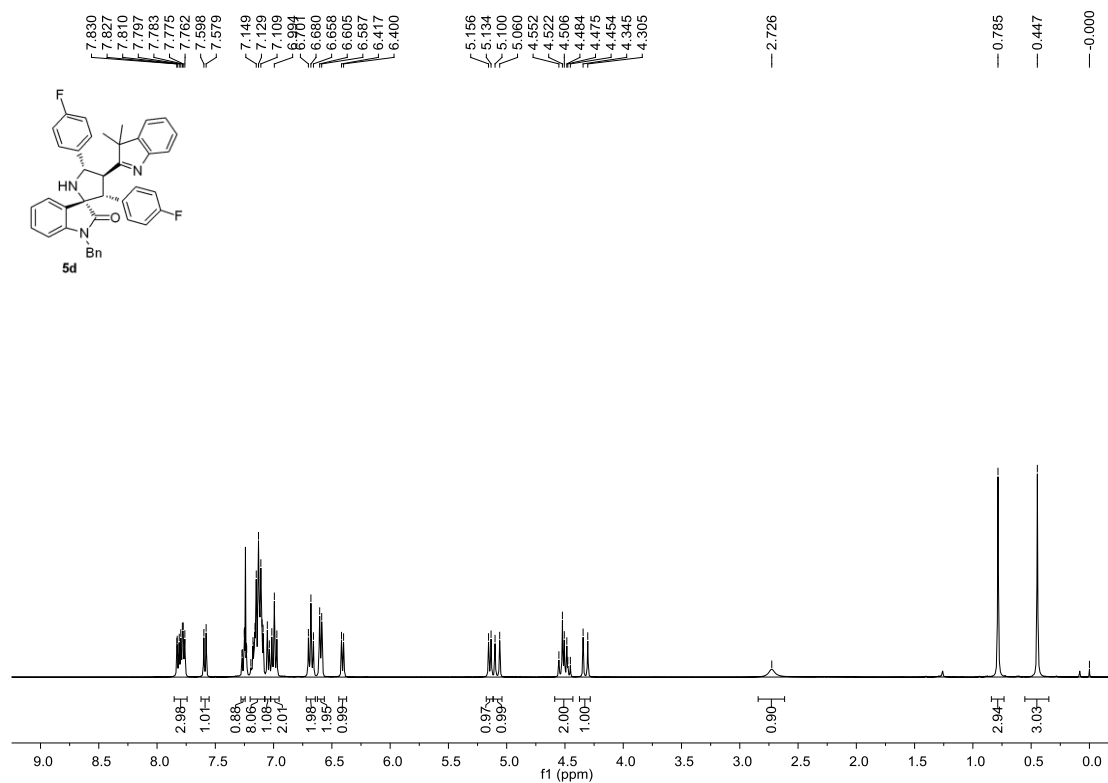
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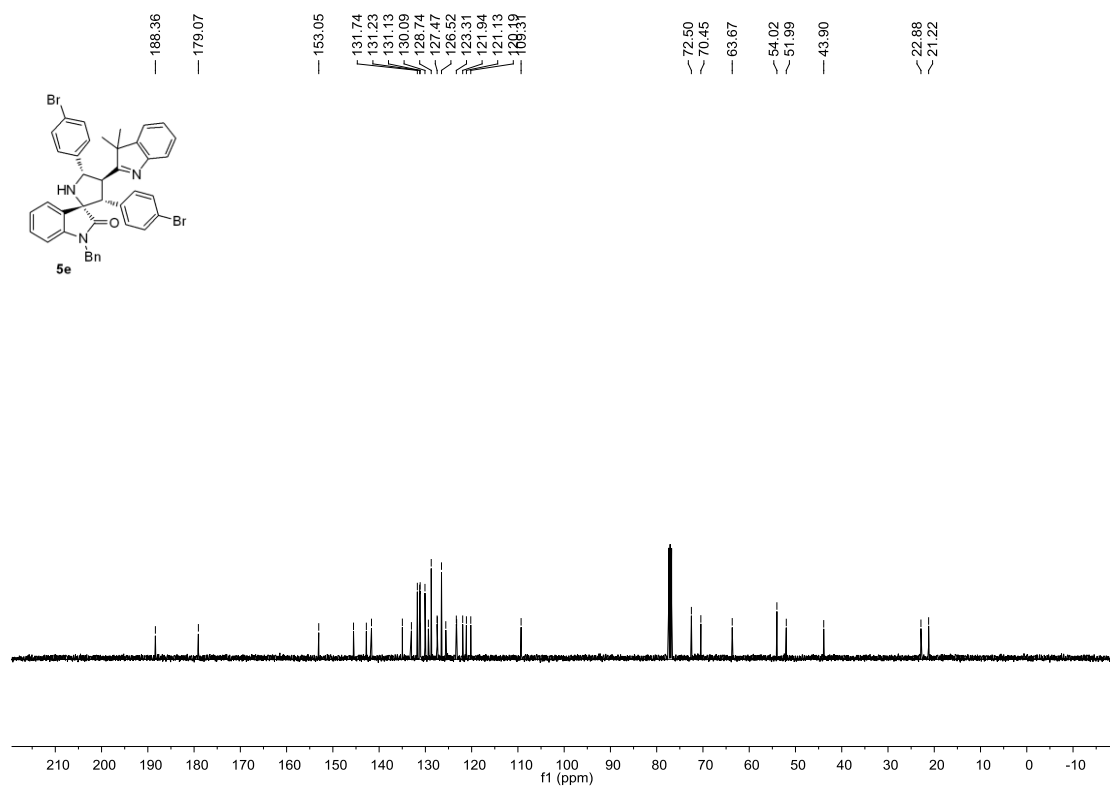
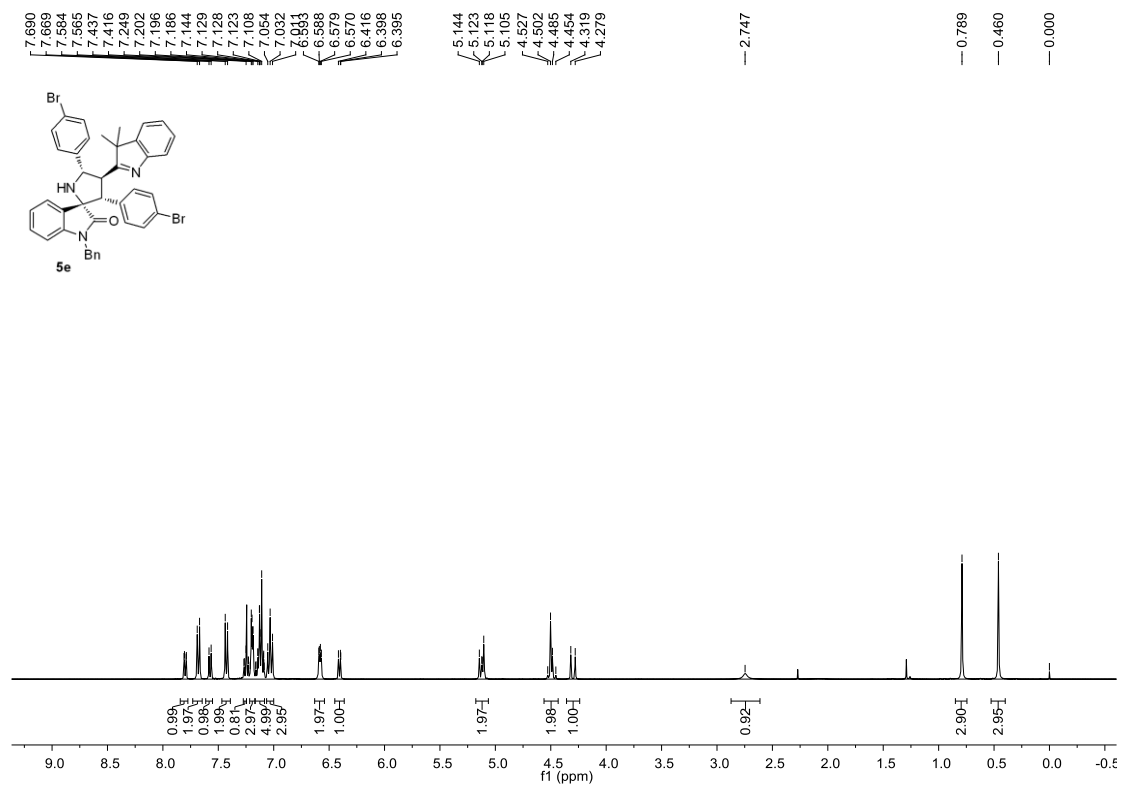
1. Copies of ^1H NMR and ^{13}C NMR spectra

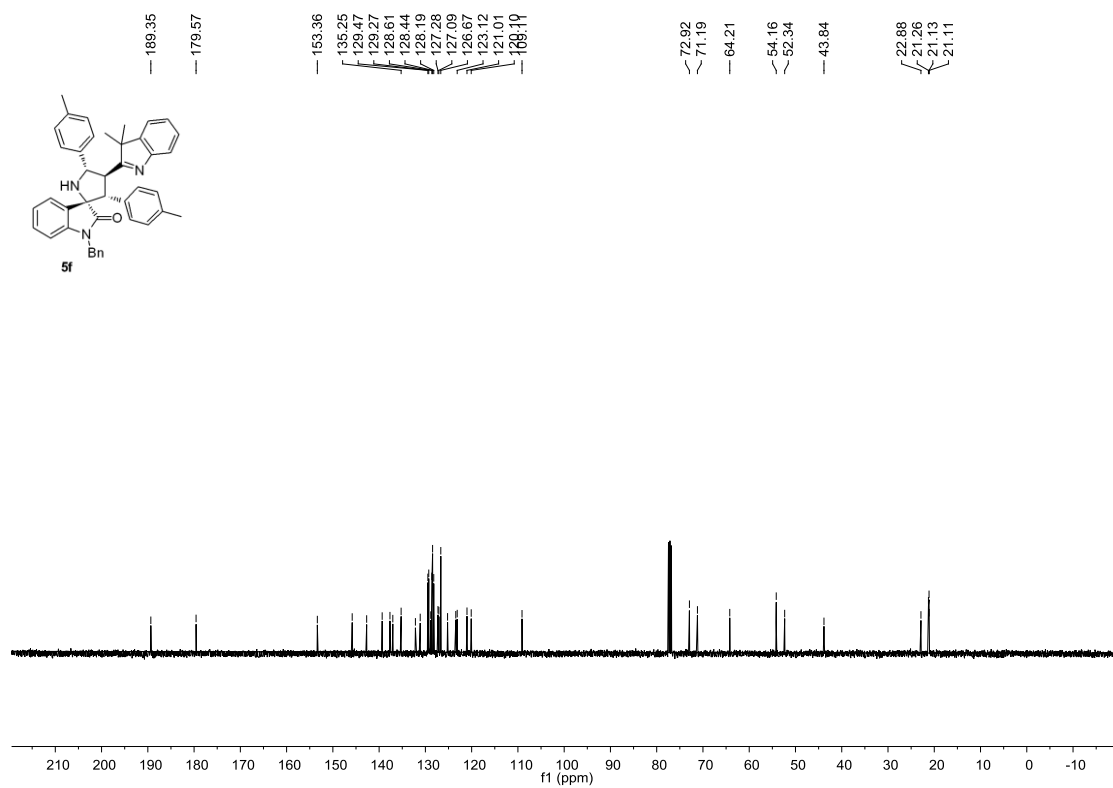
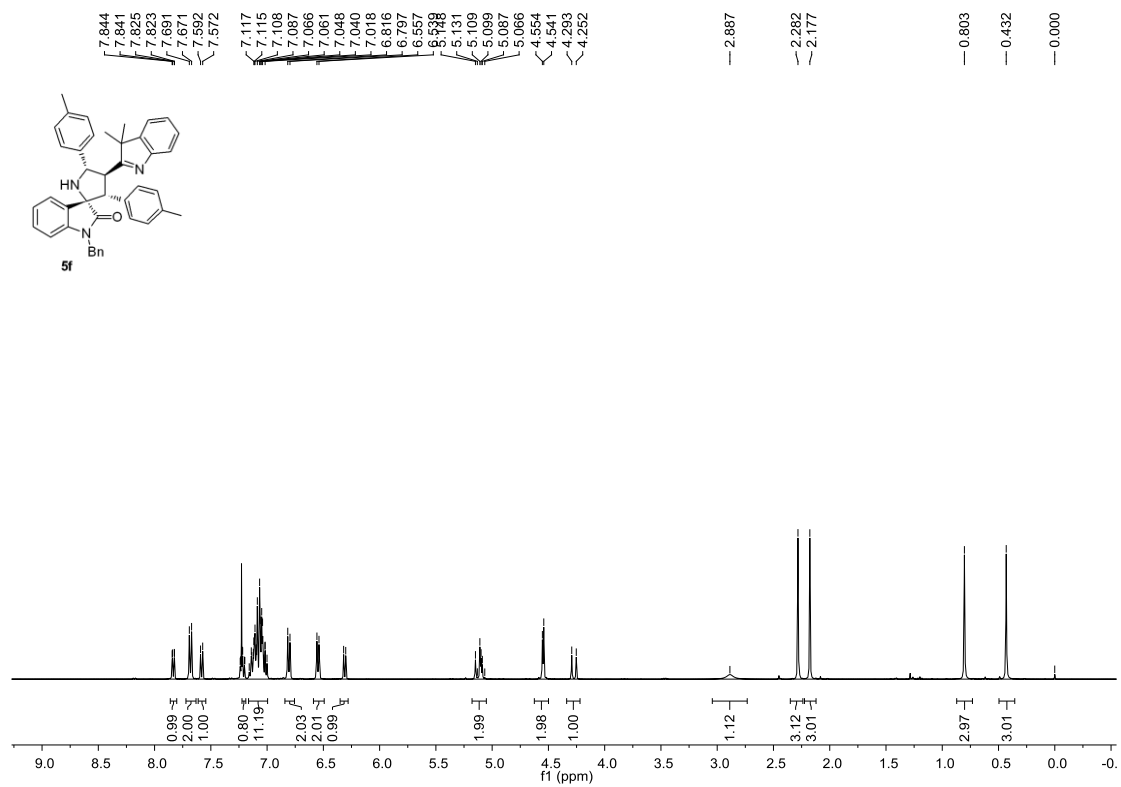


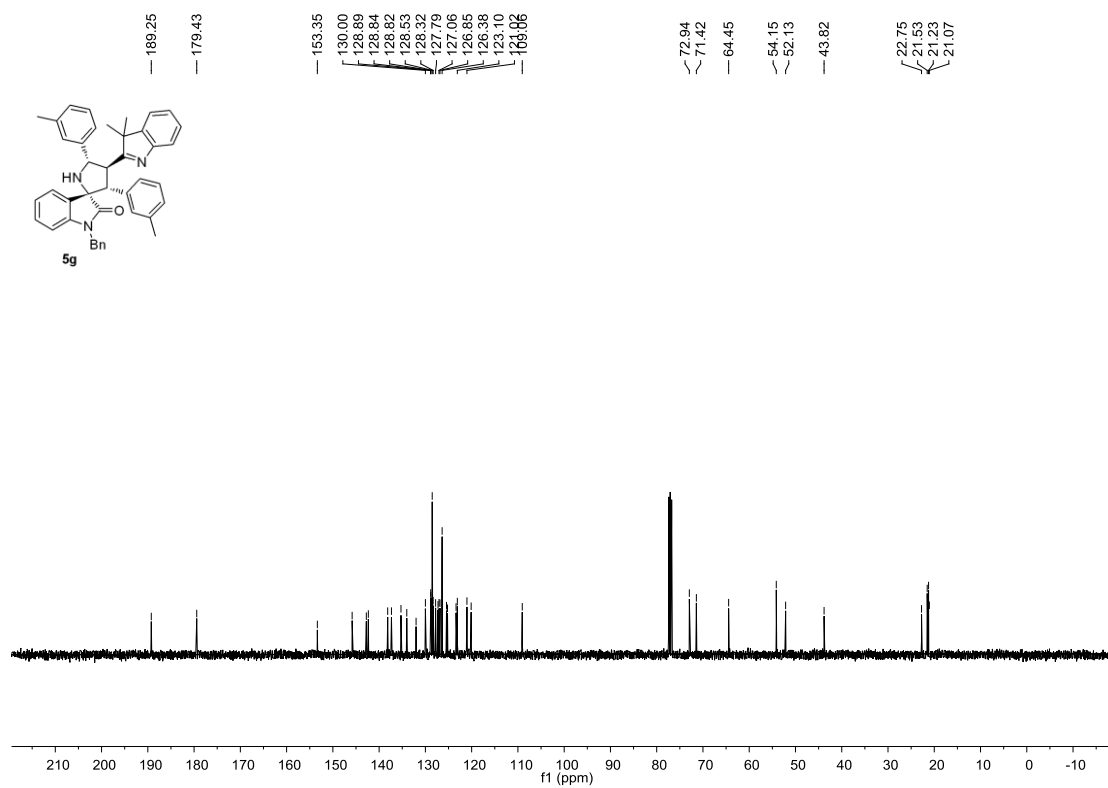
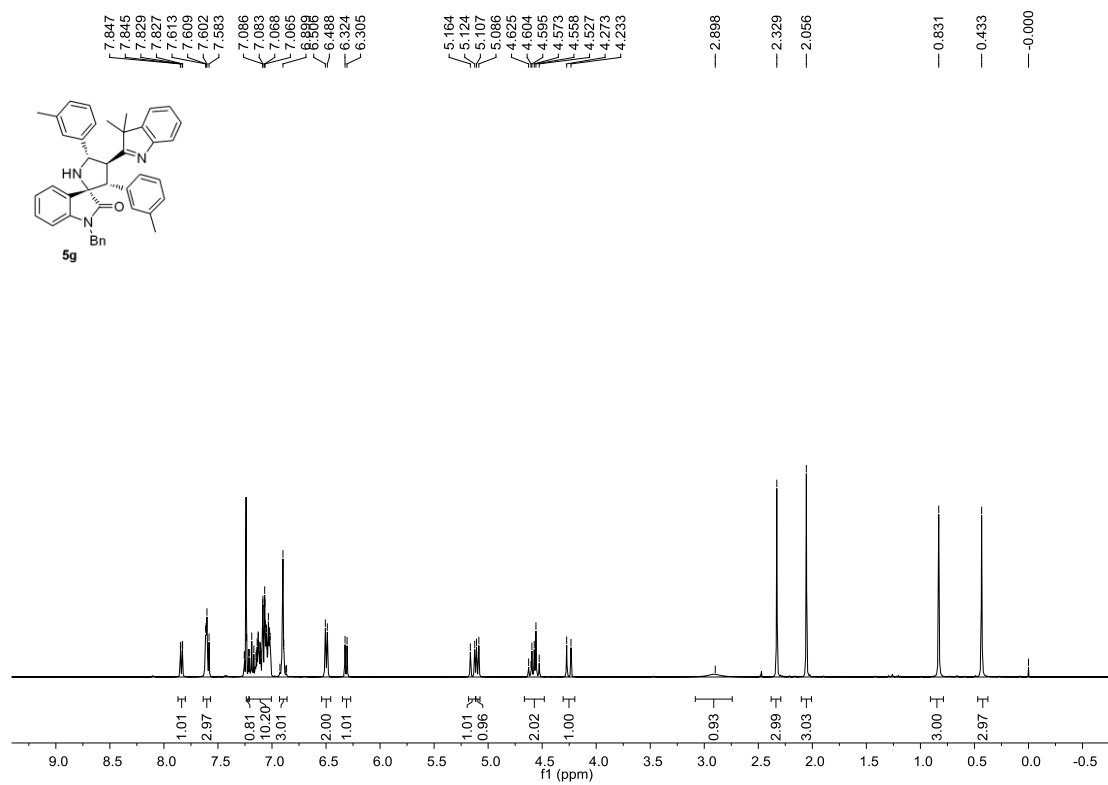


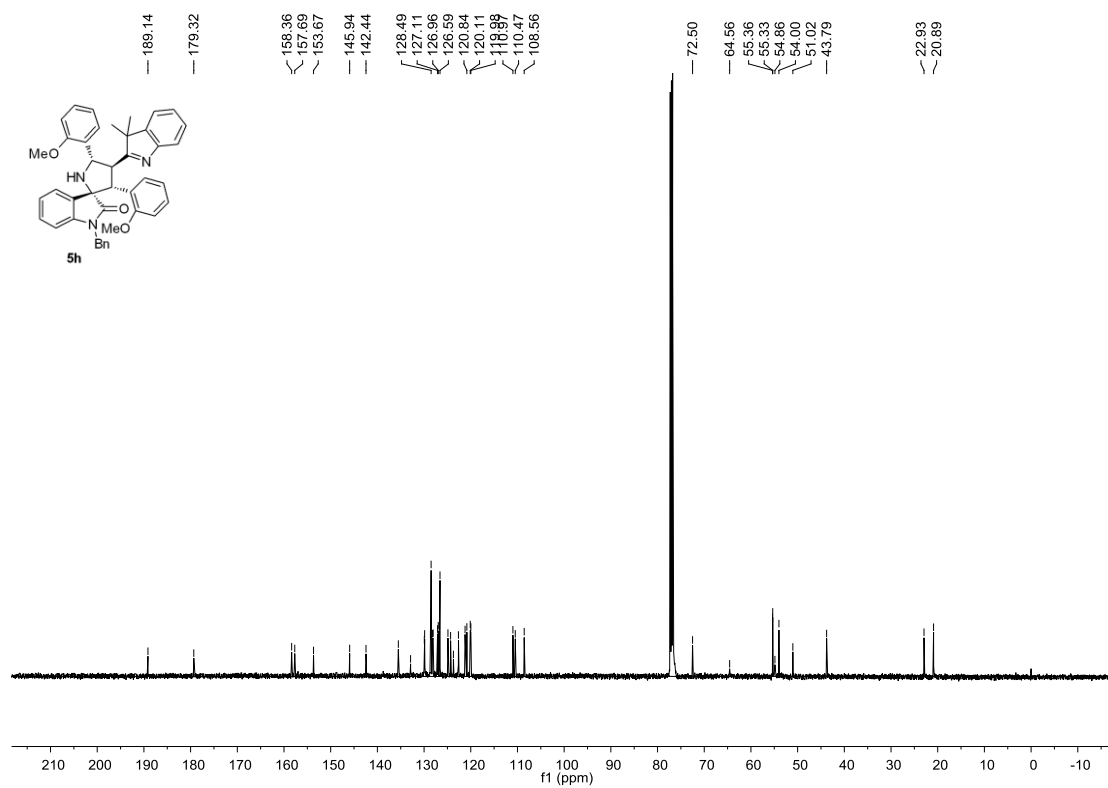
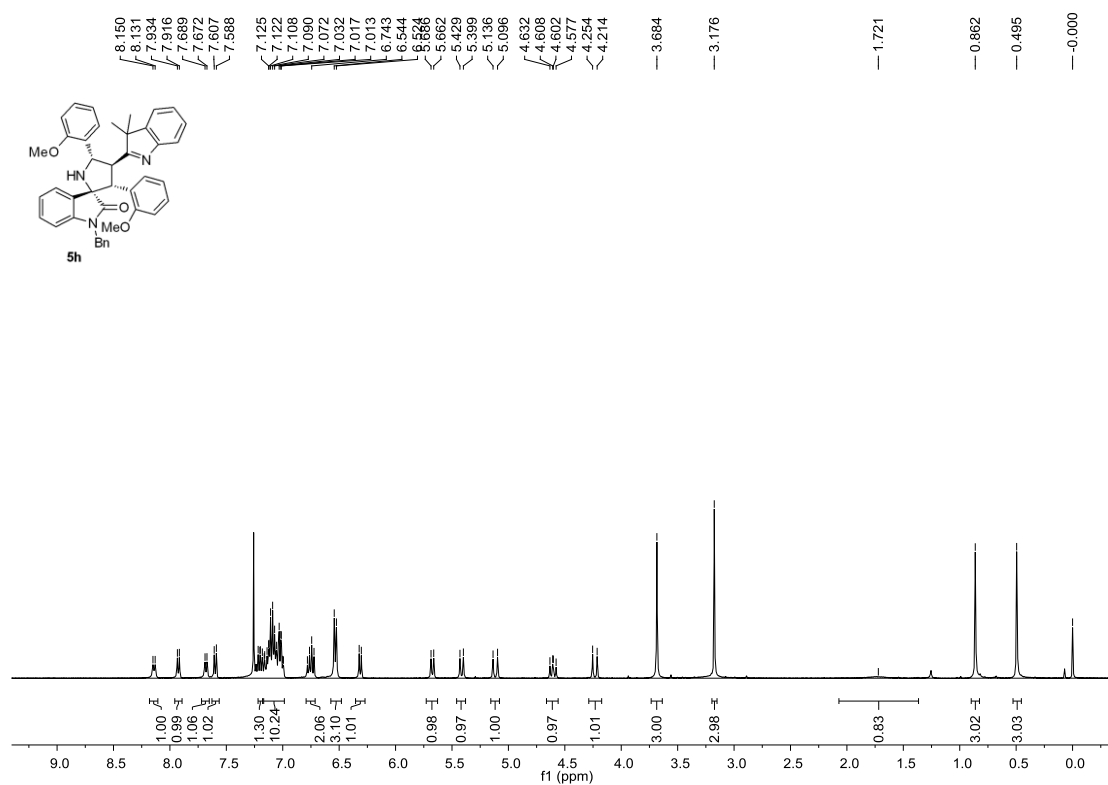


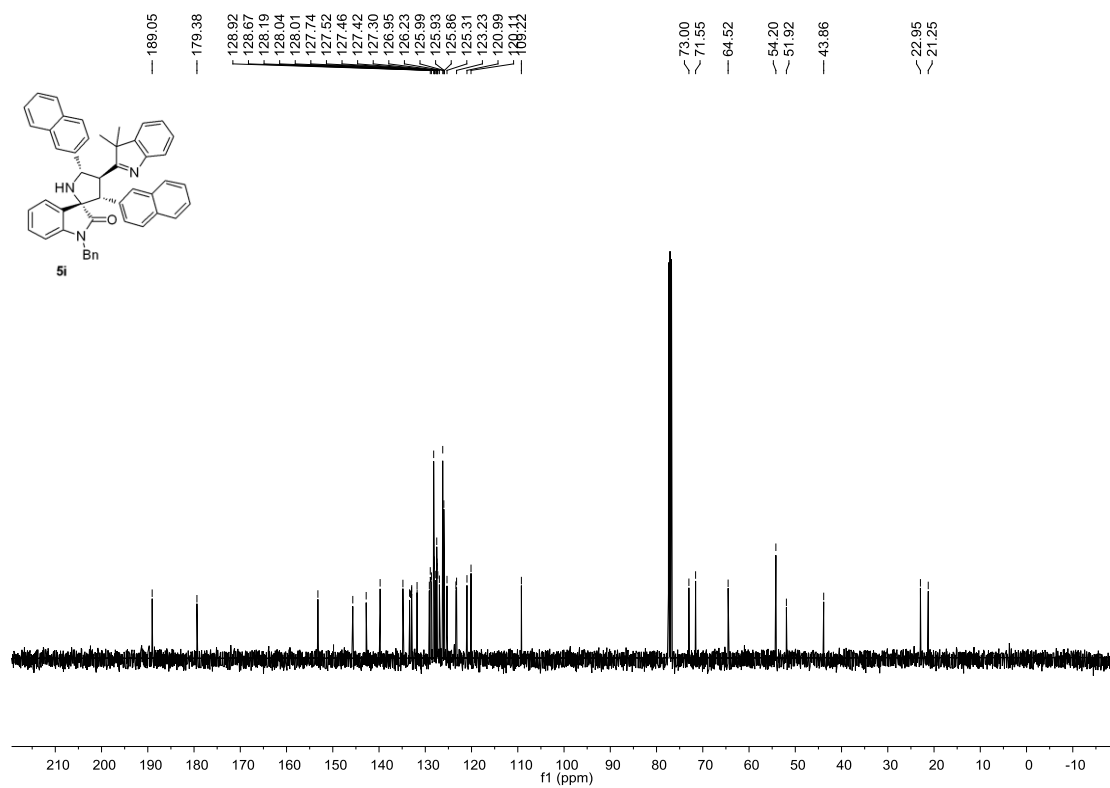
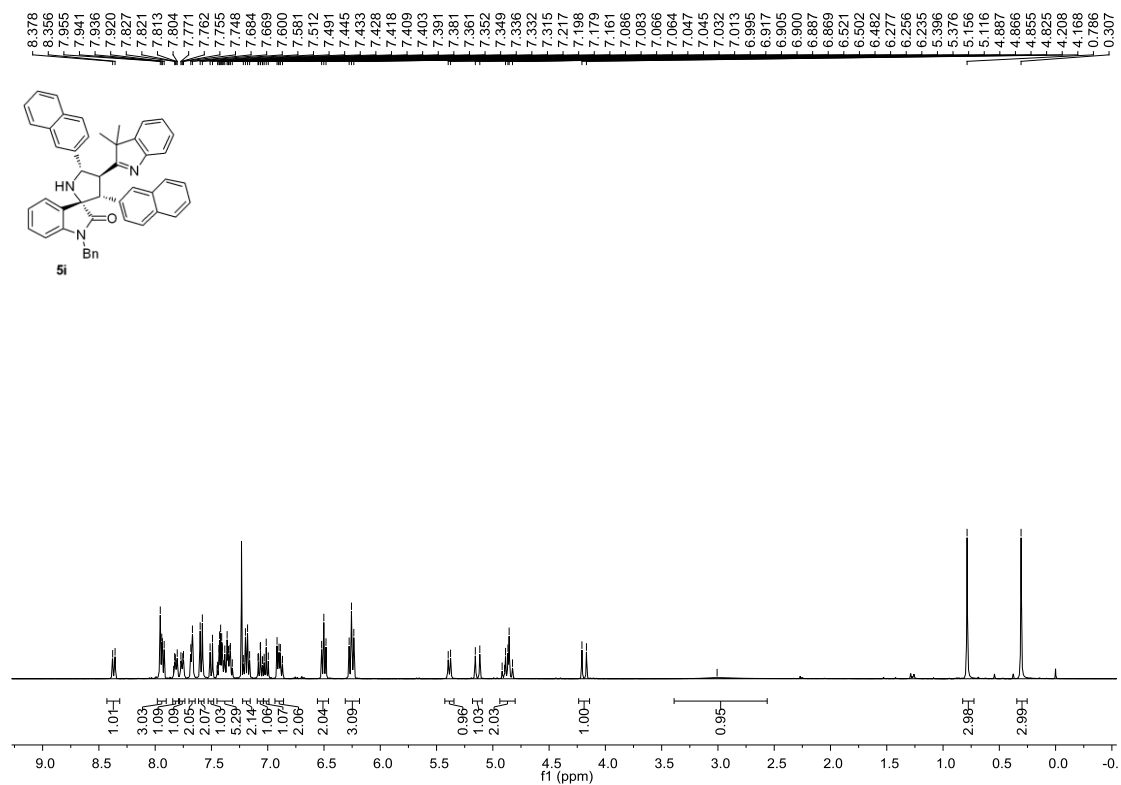


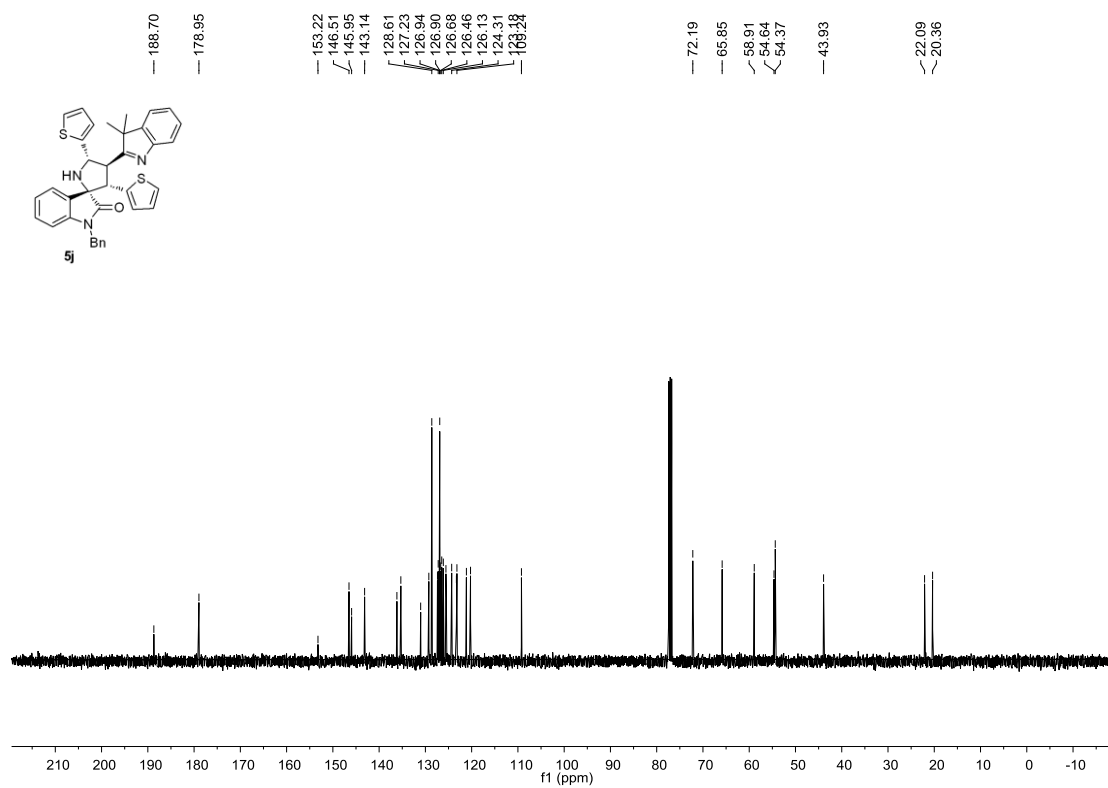
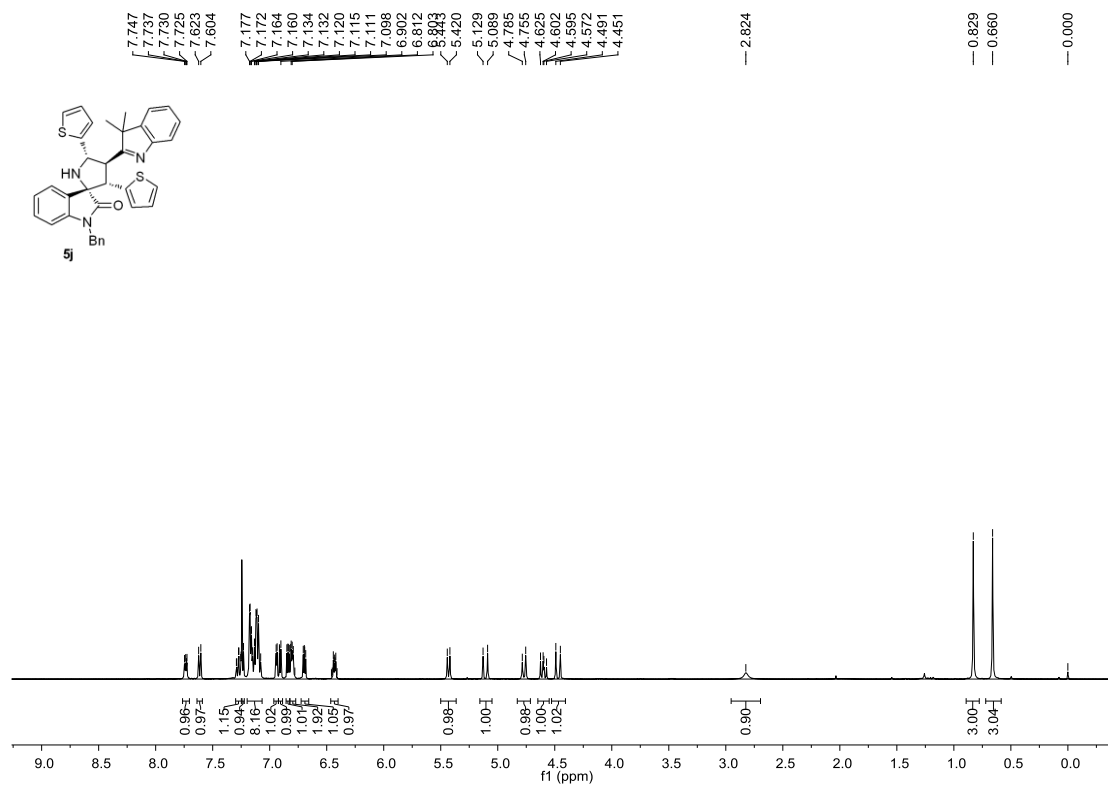


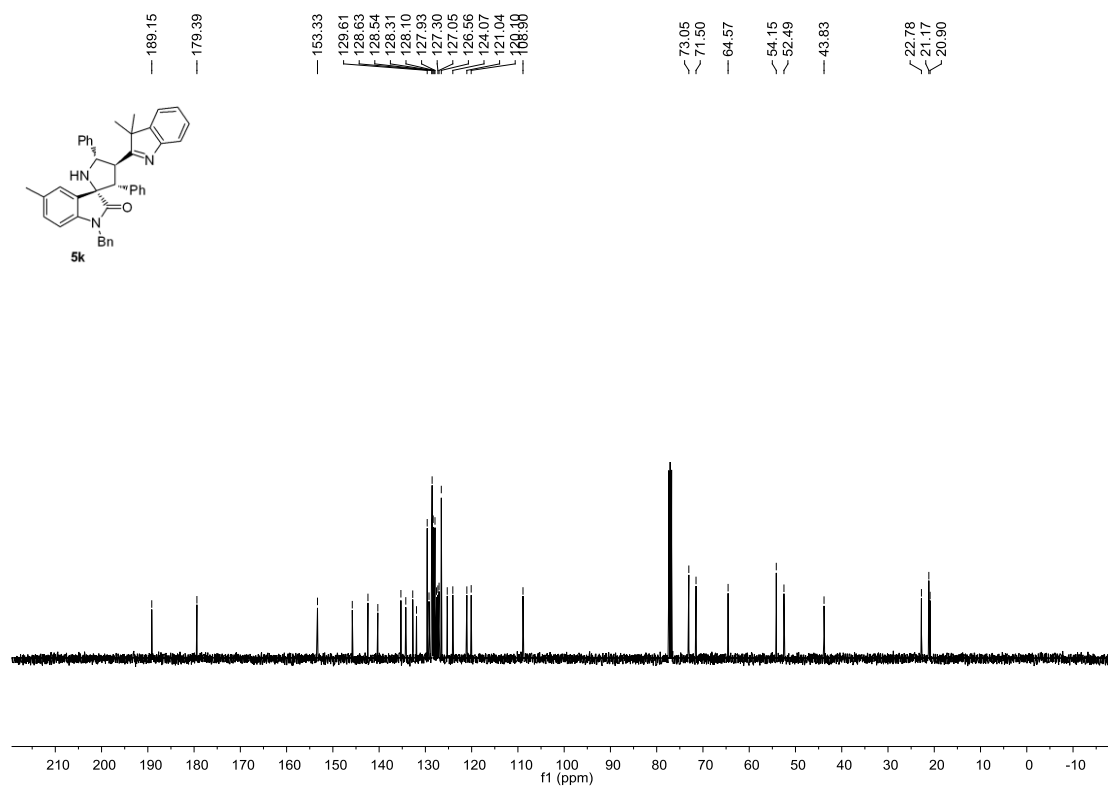
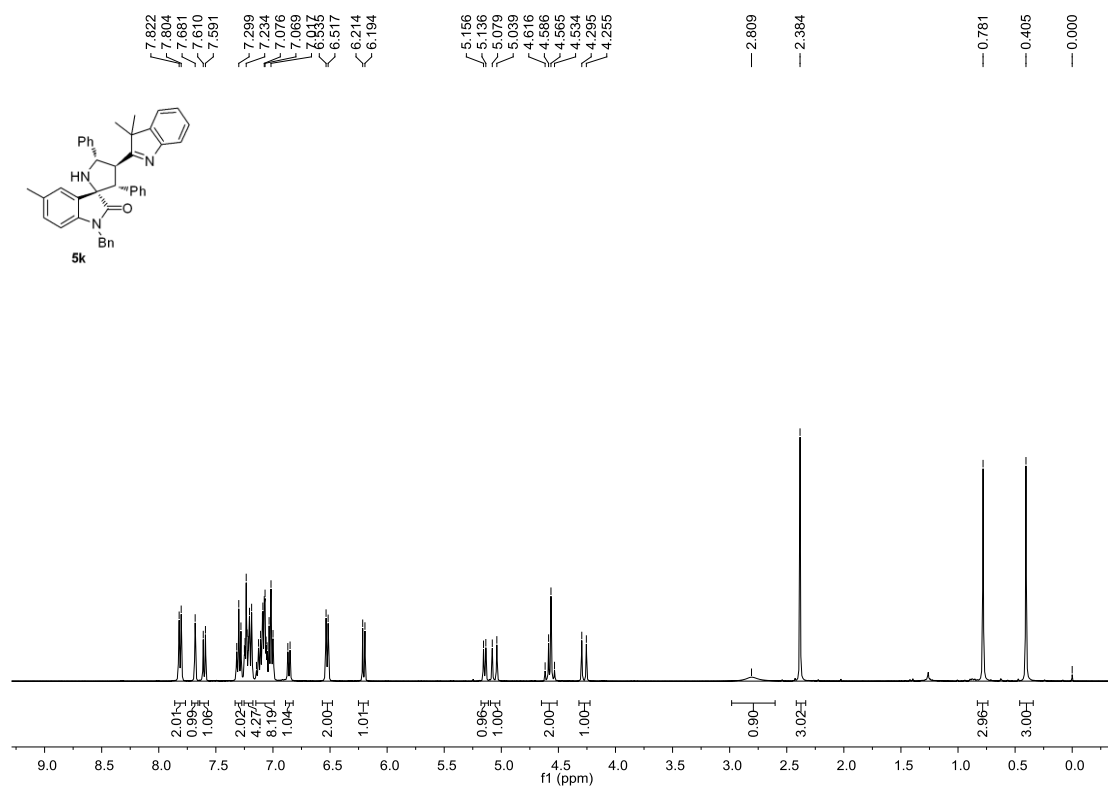


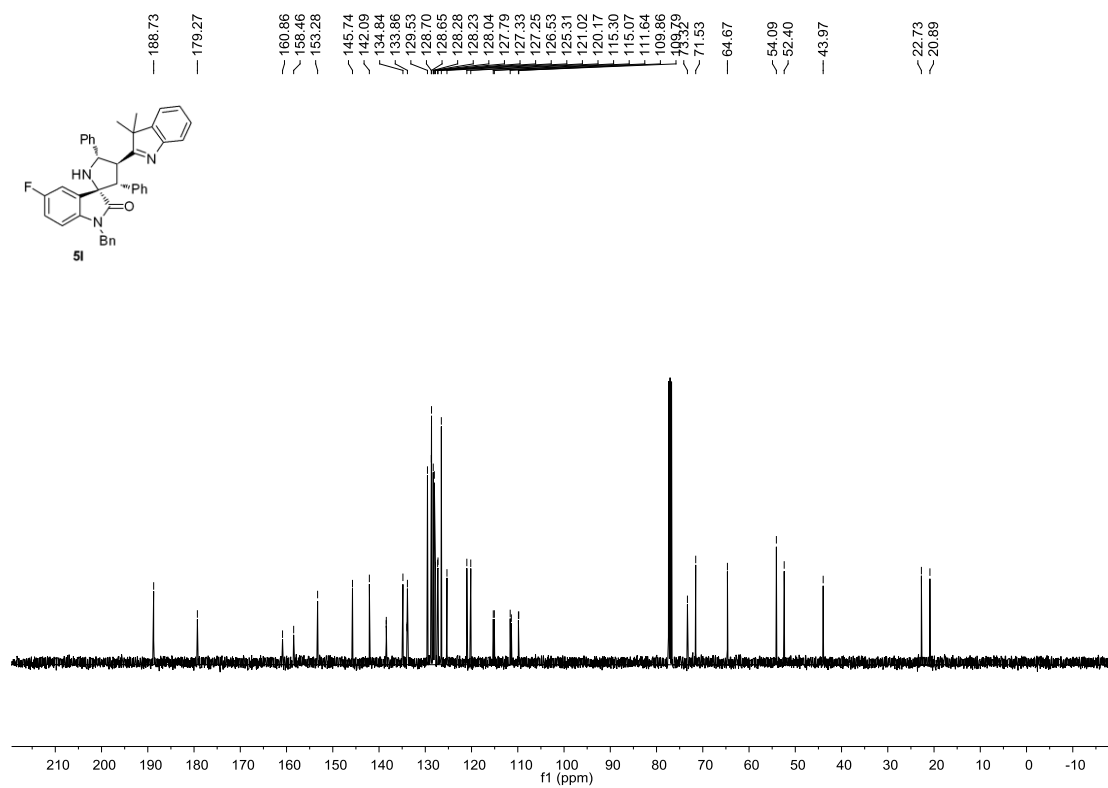
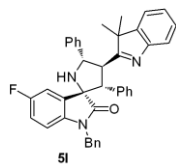


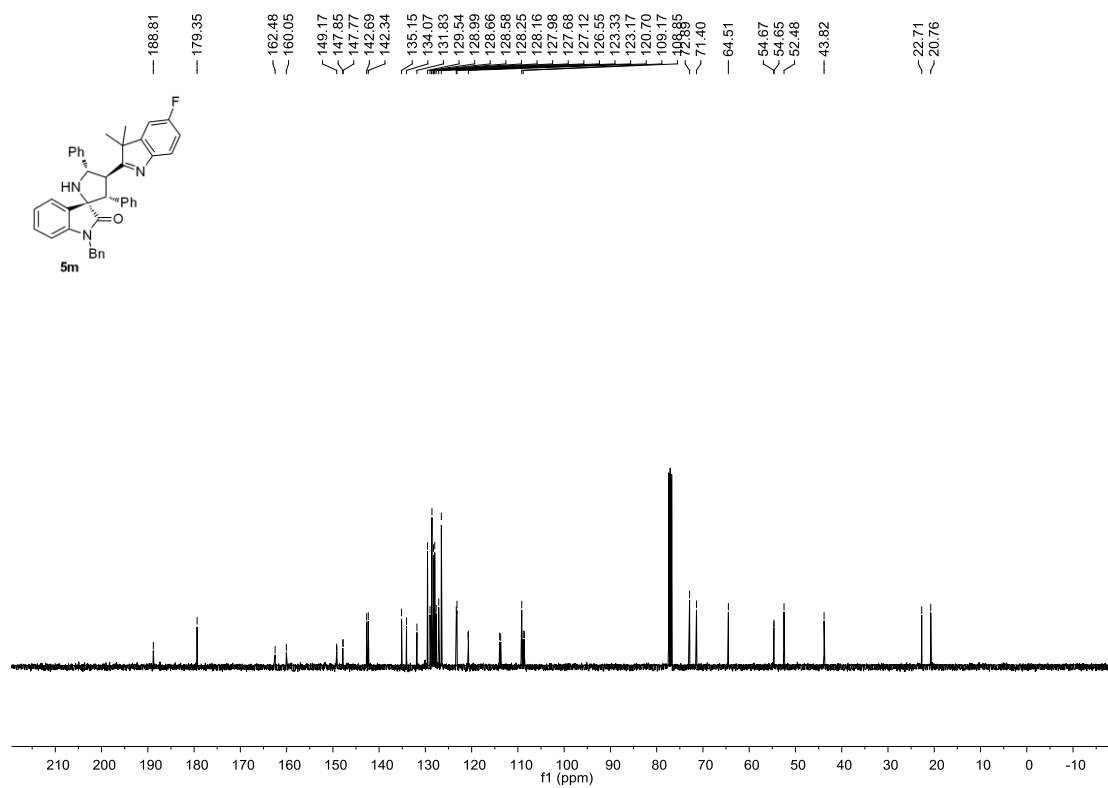
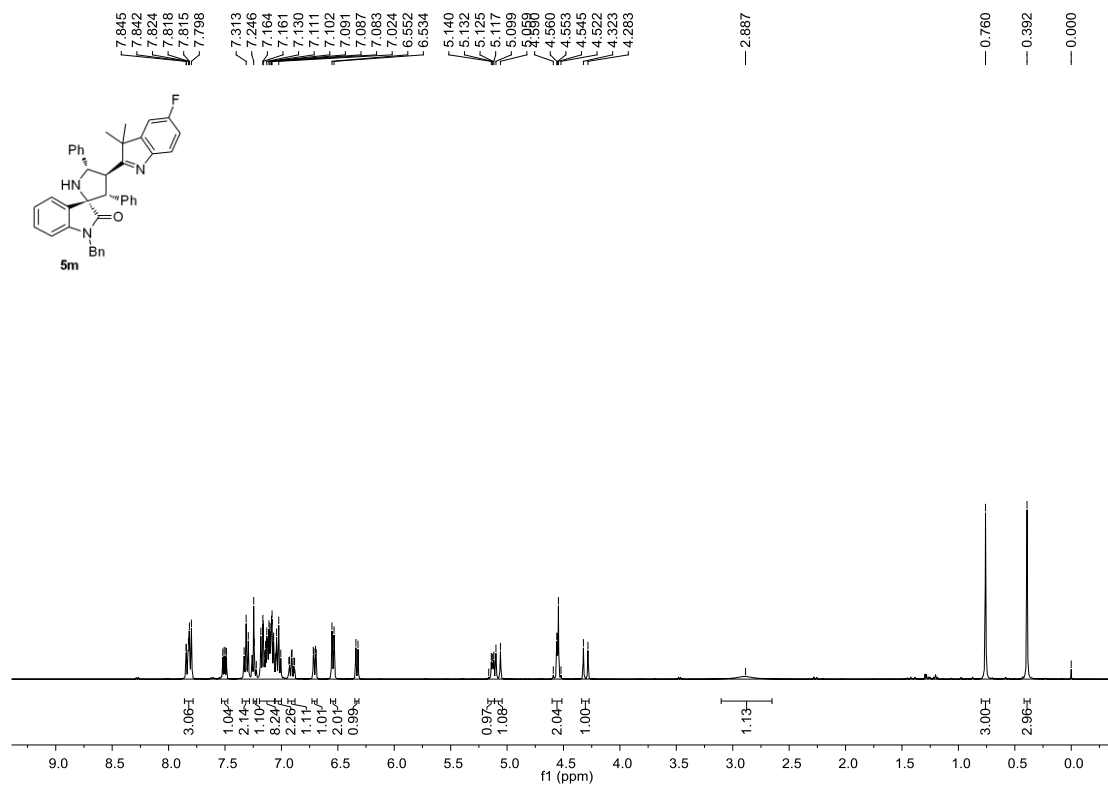


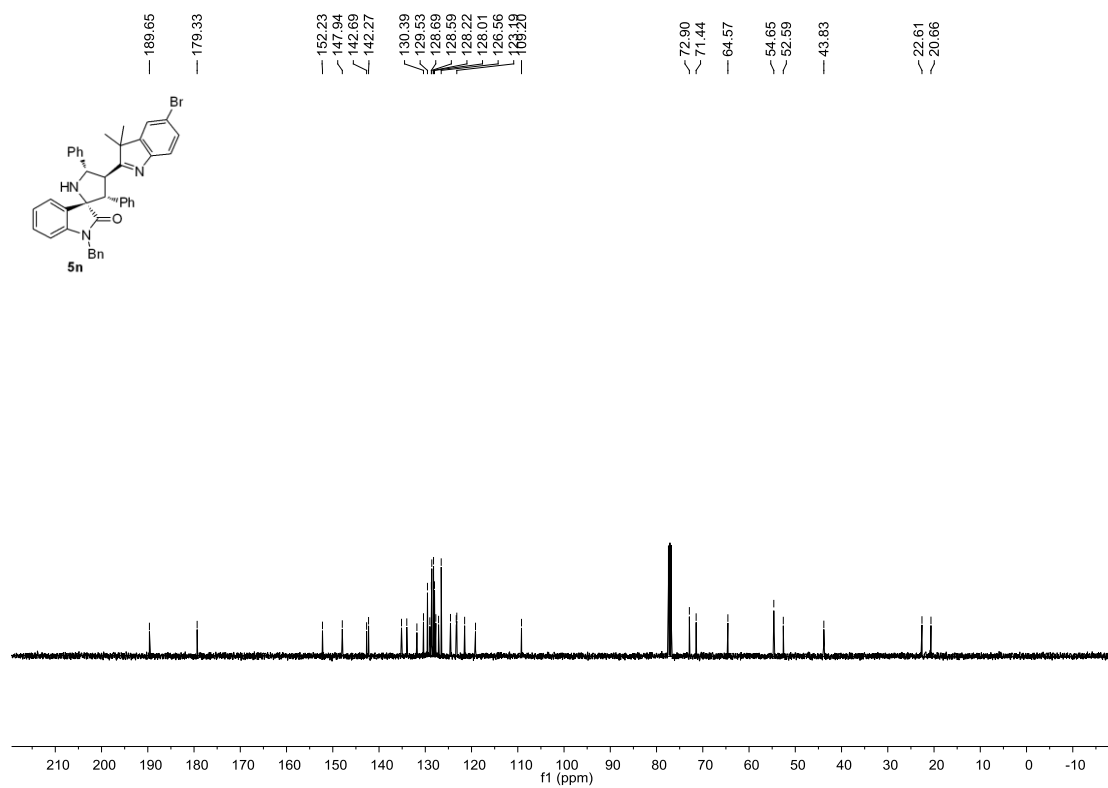
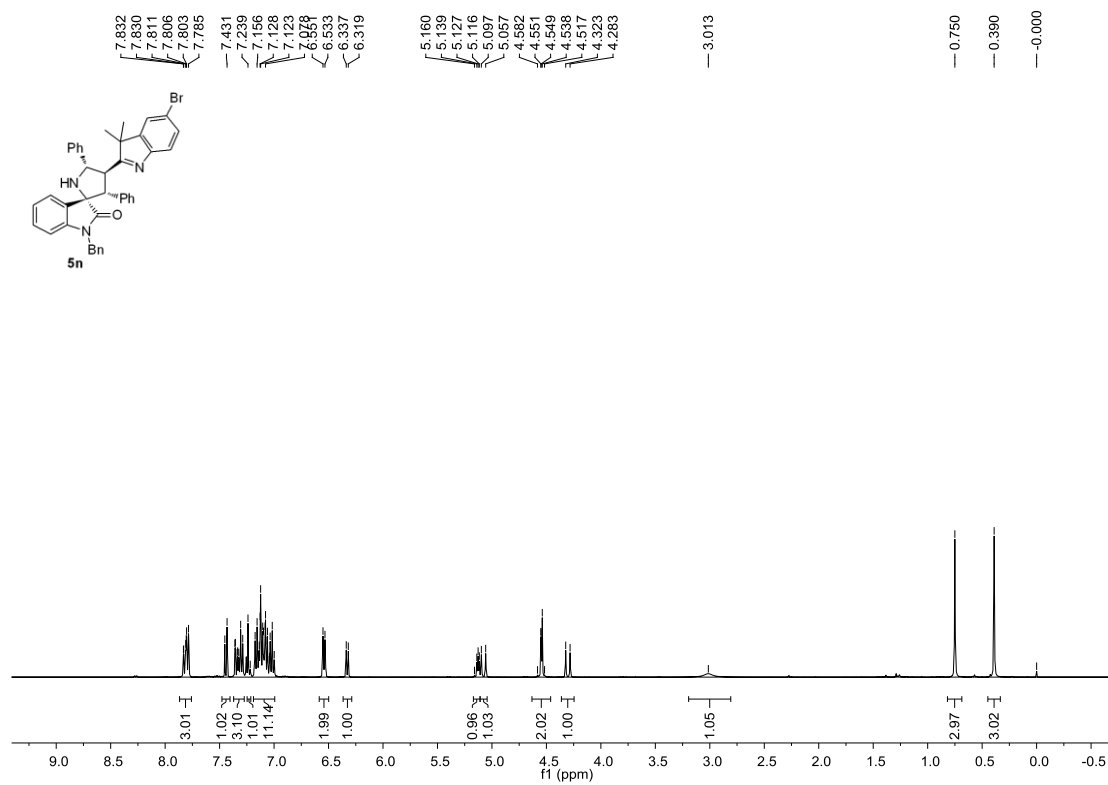


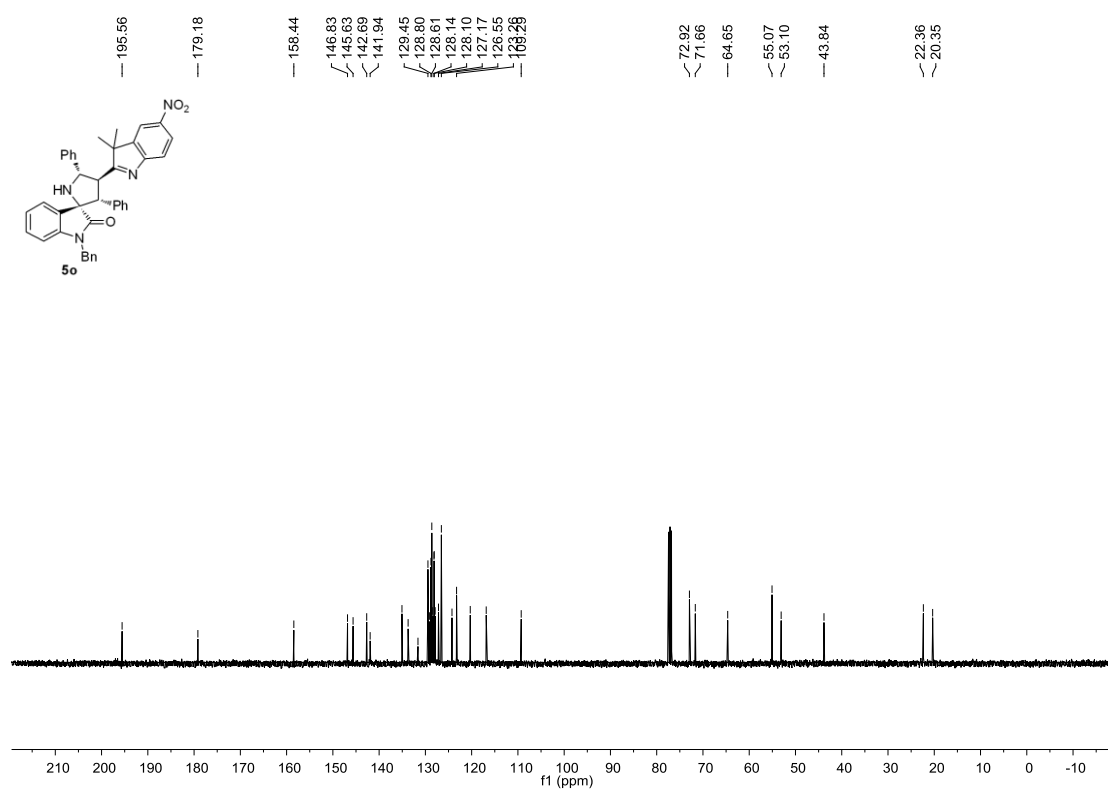
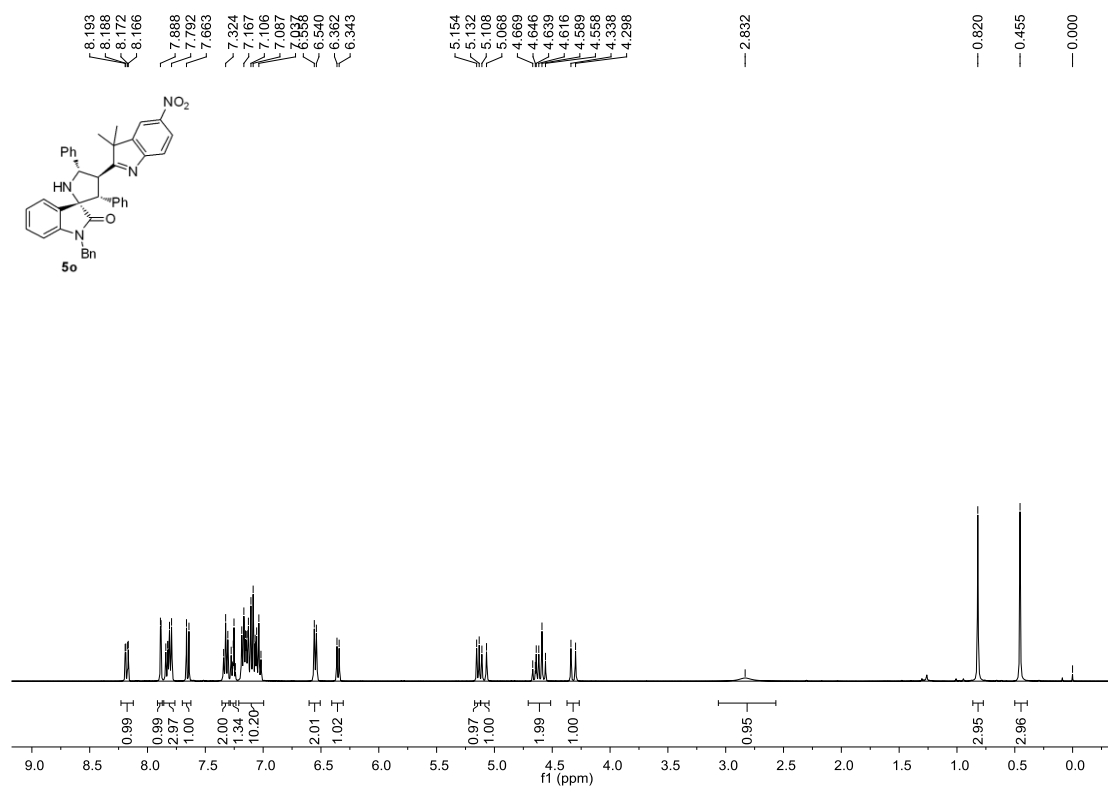


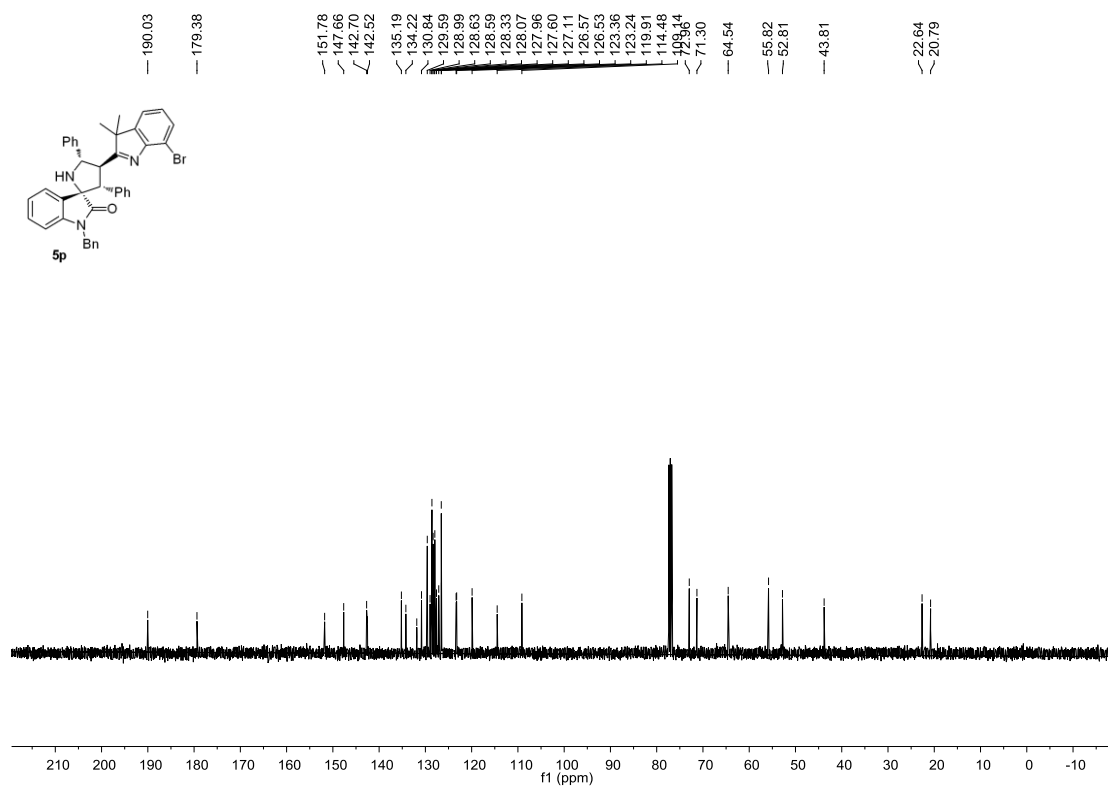
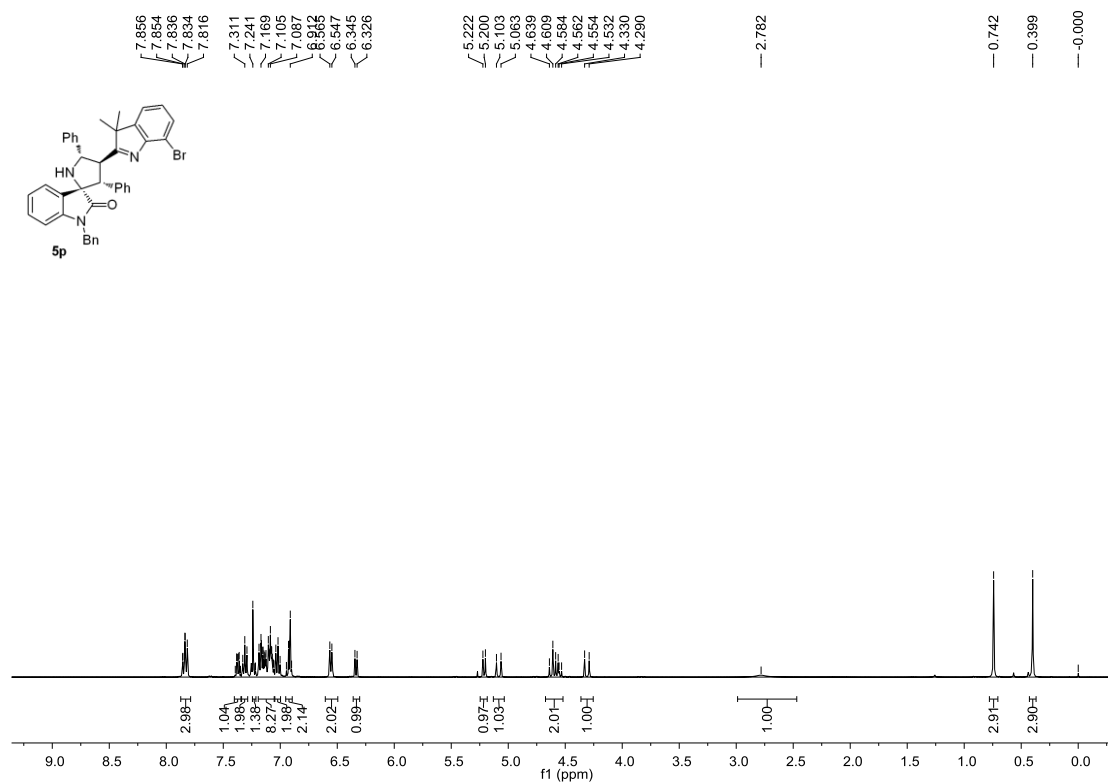


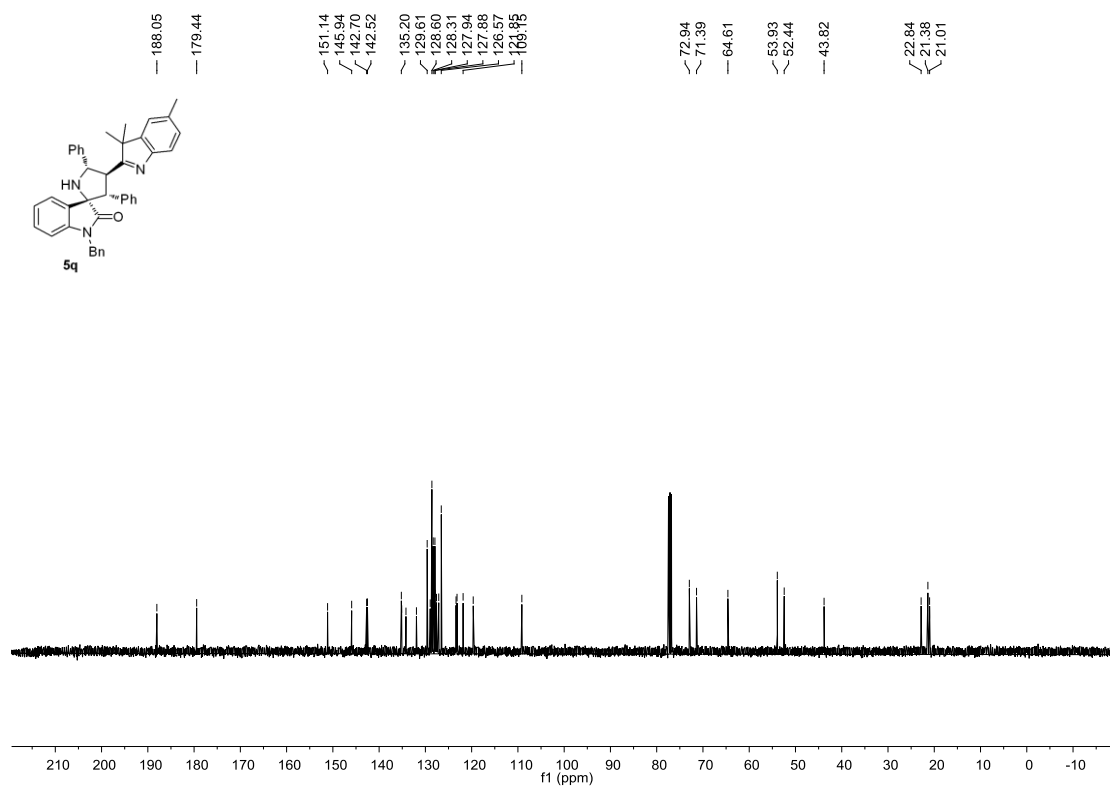
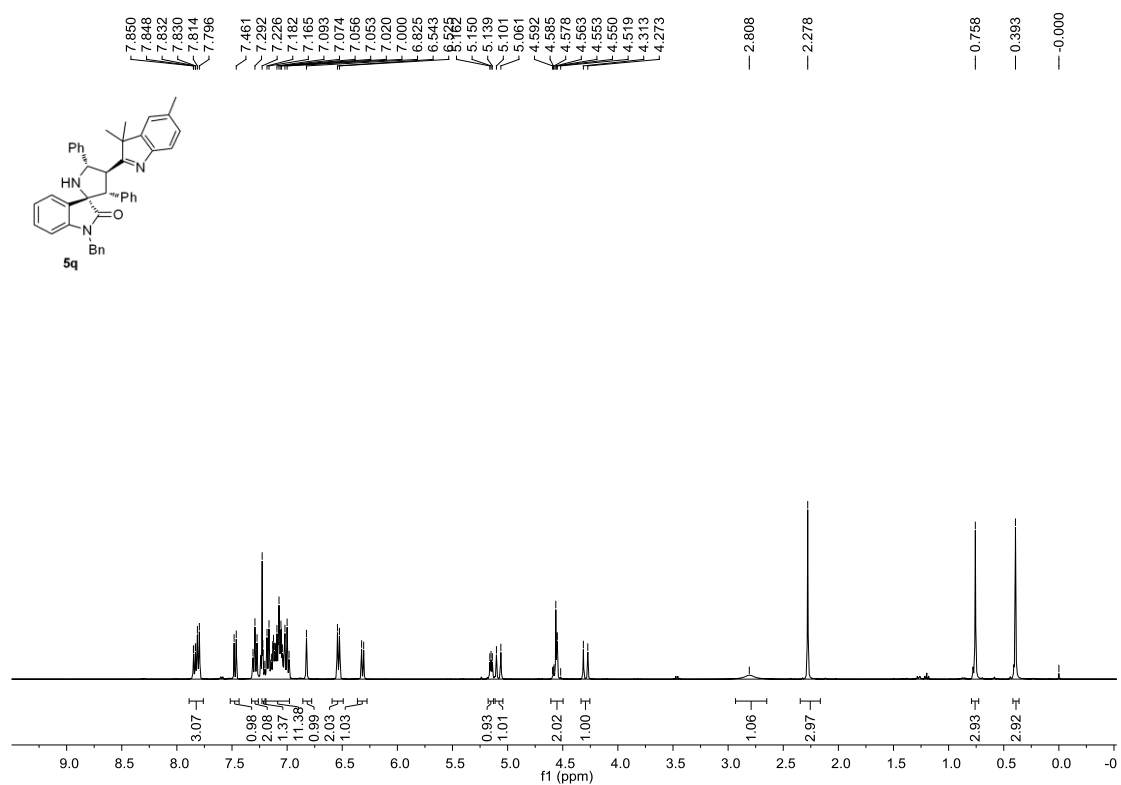


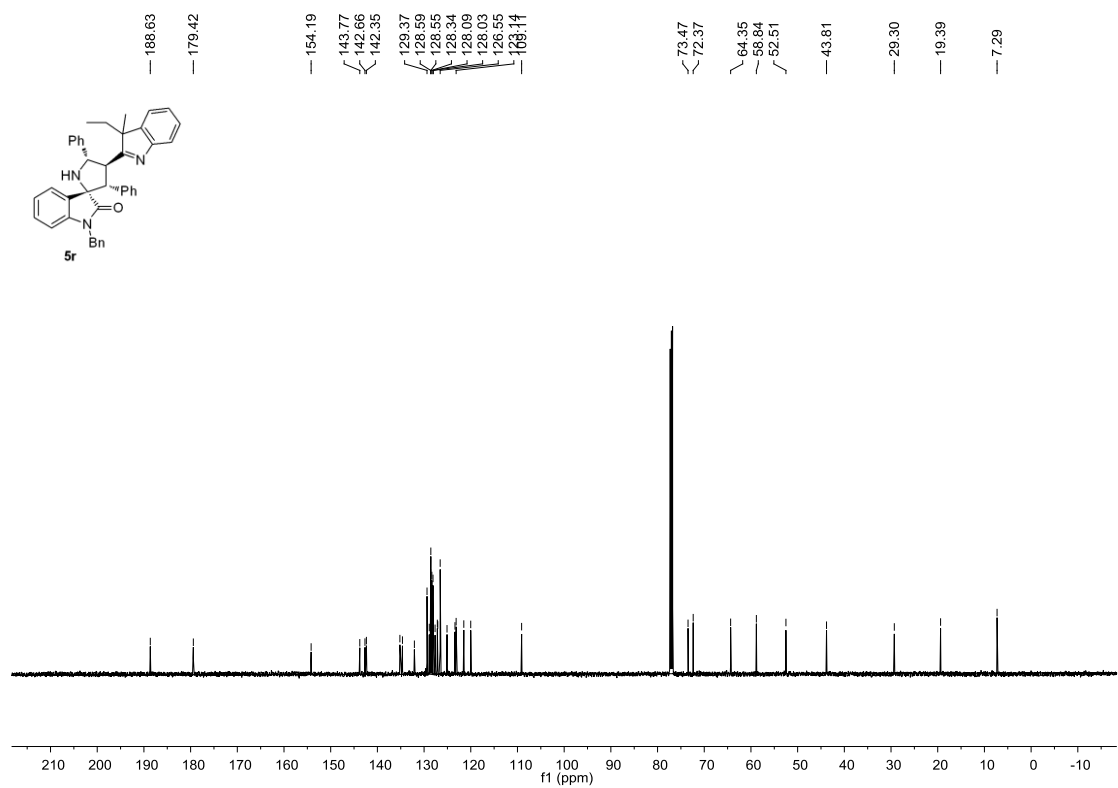
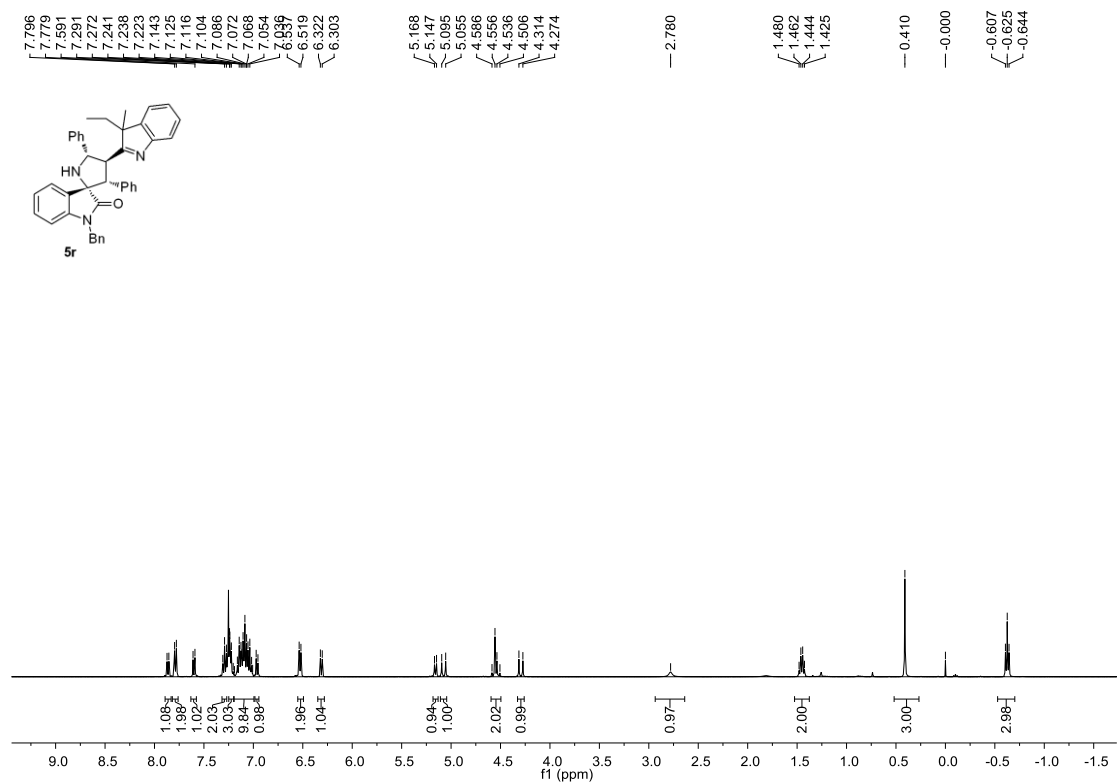


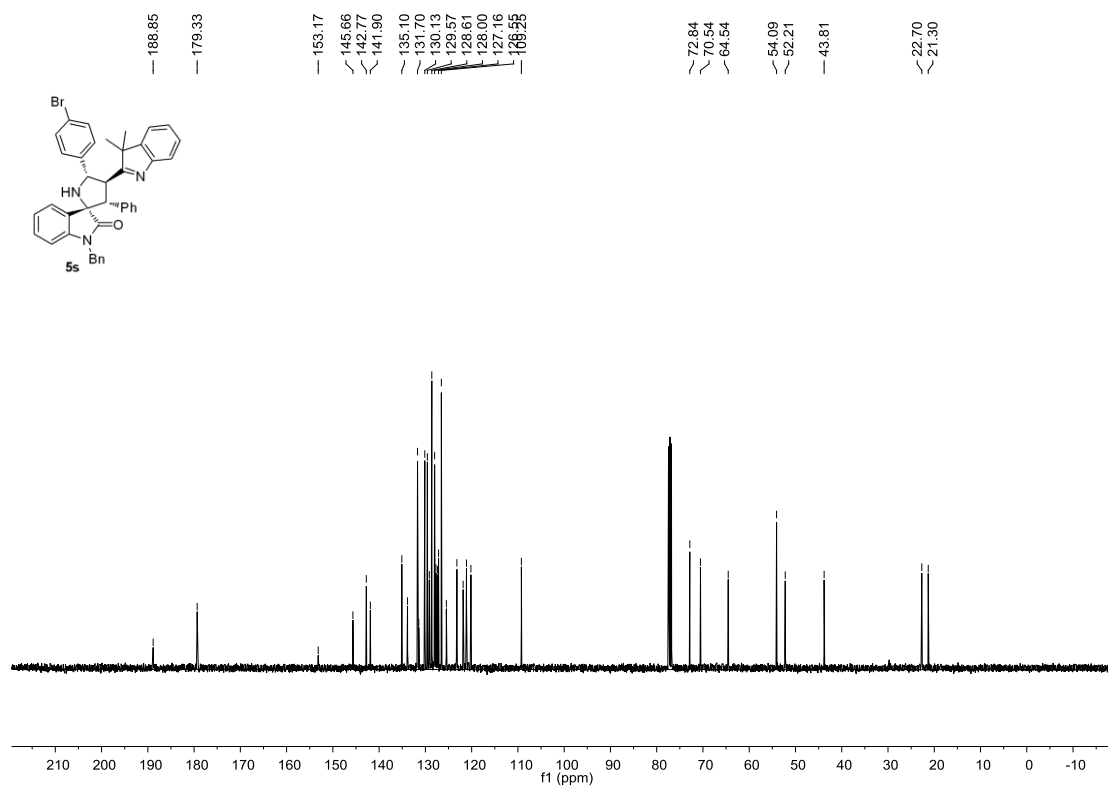
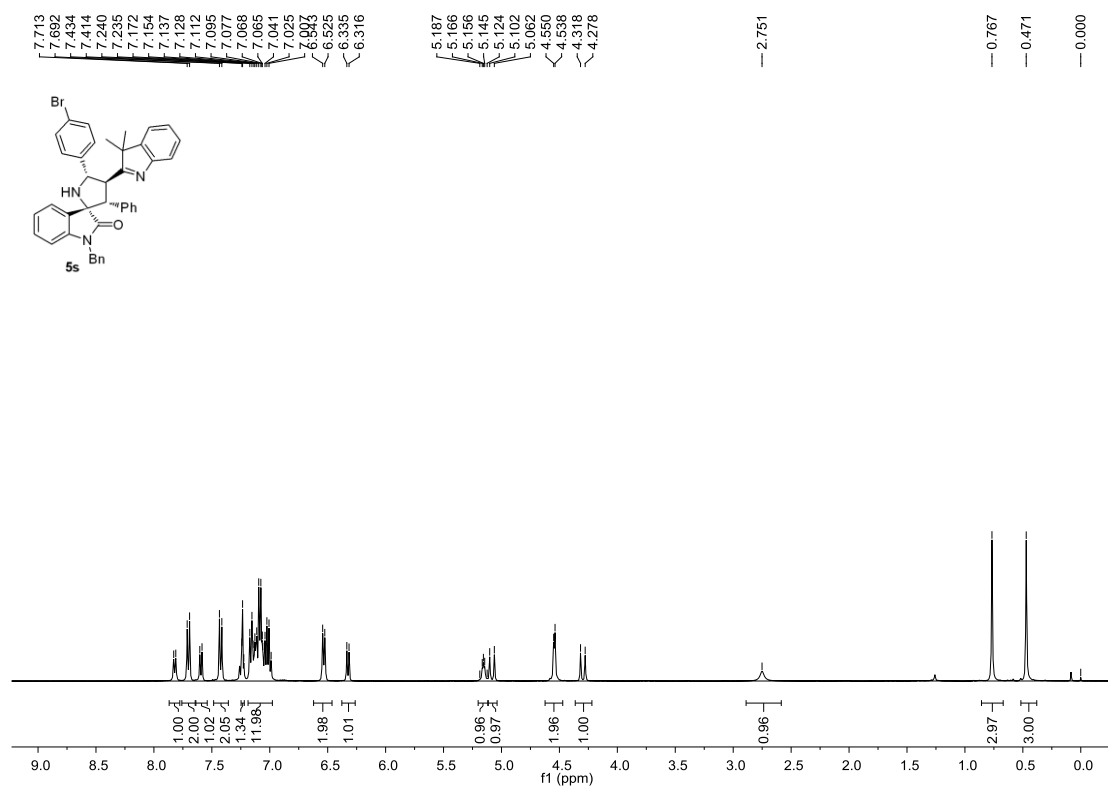


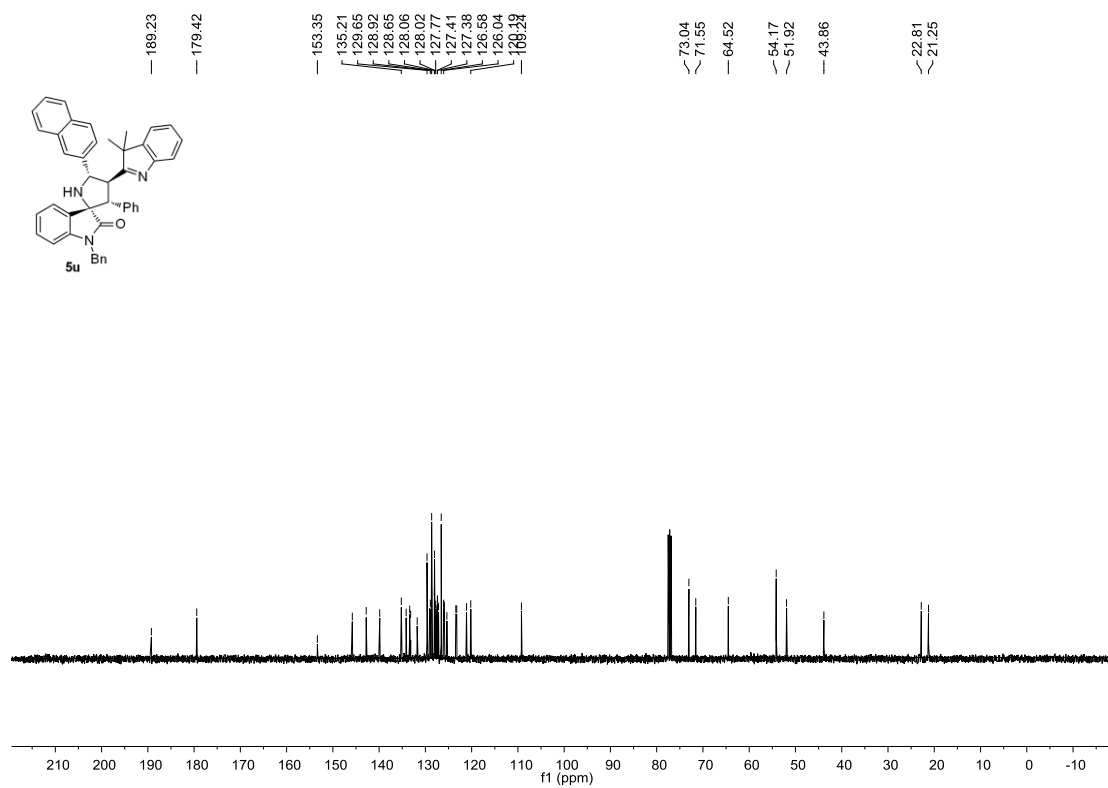
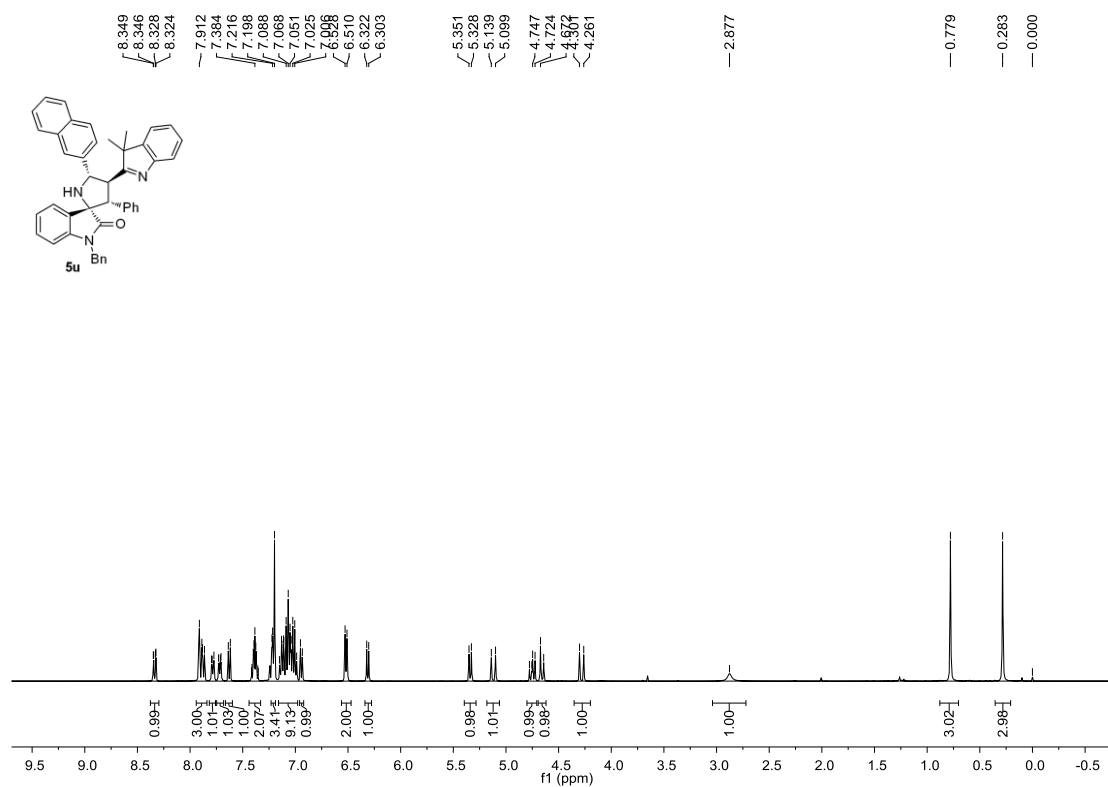


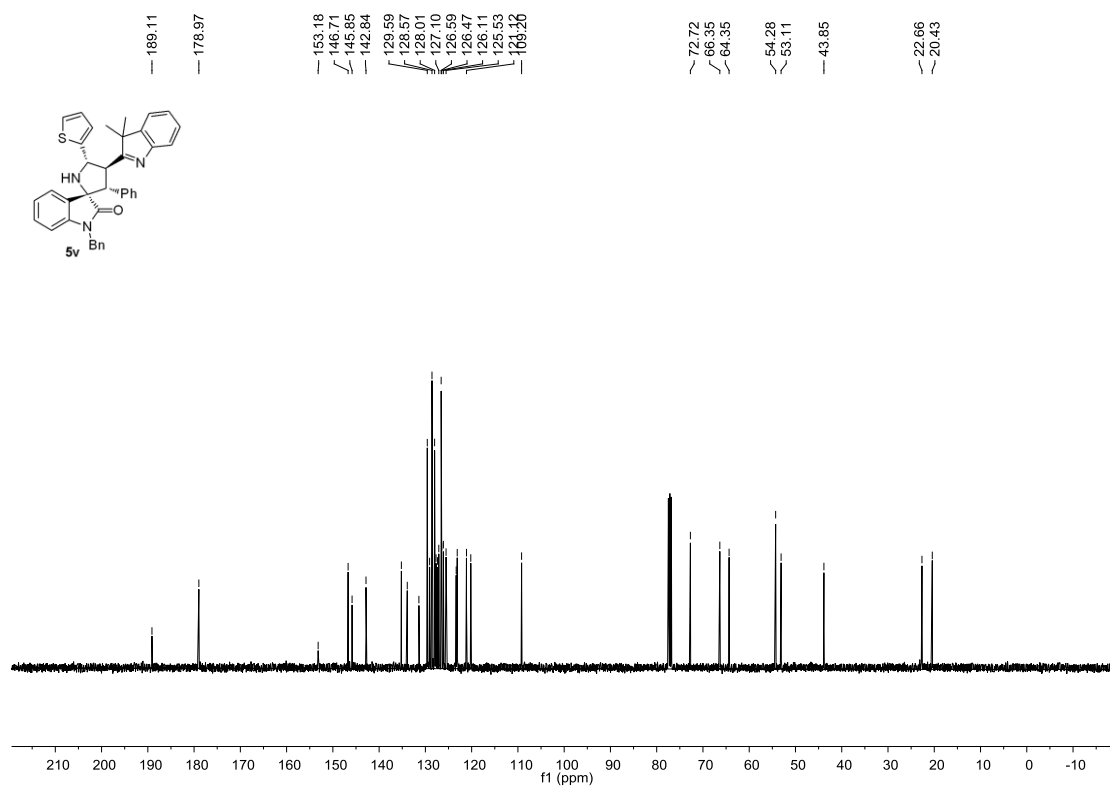
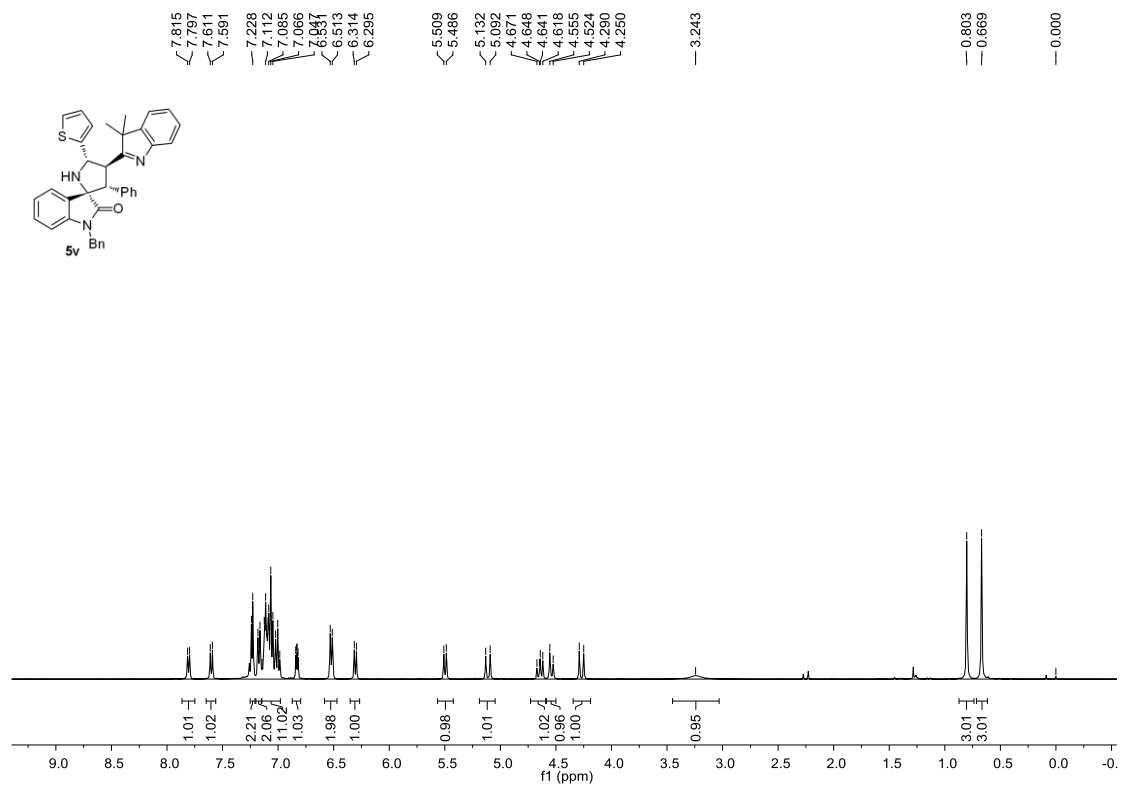


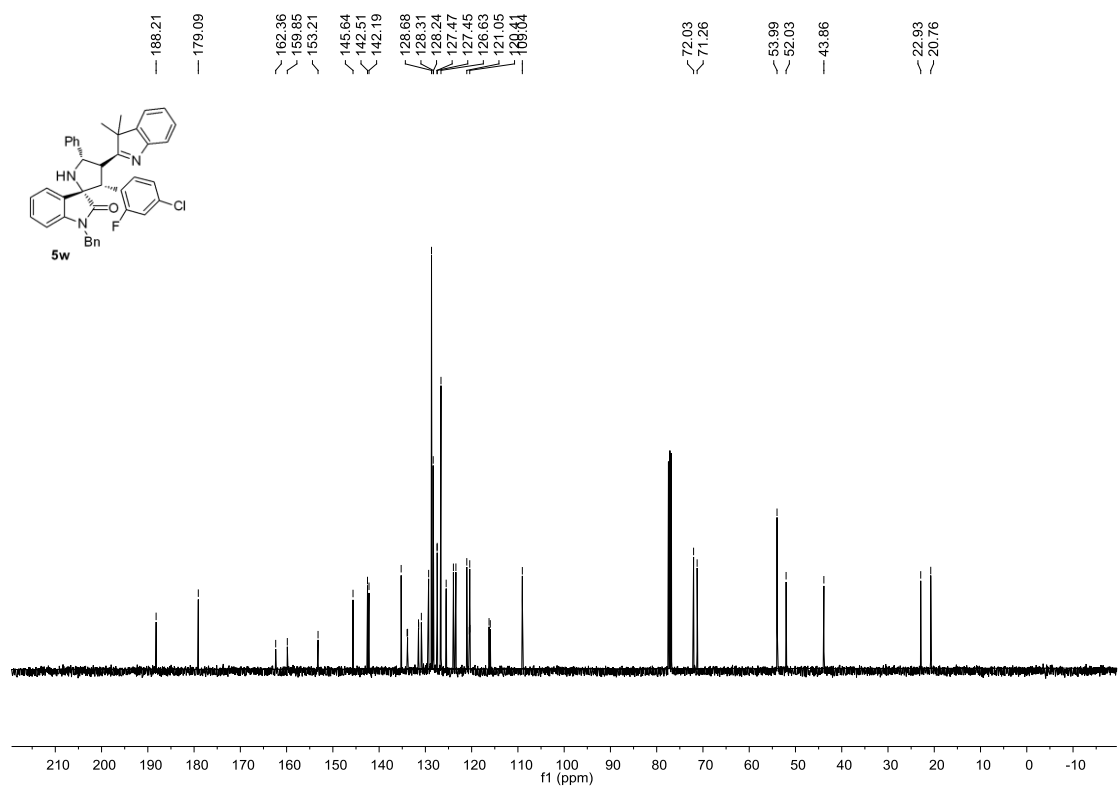
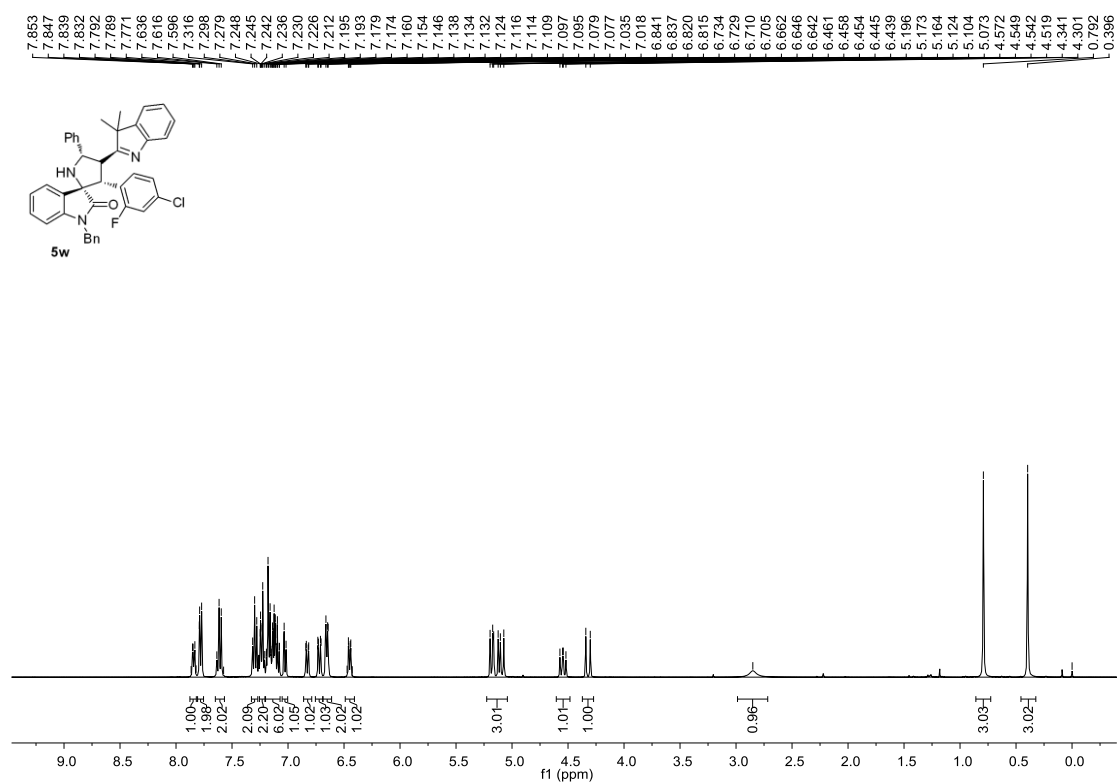


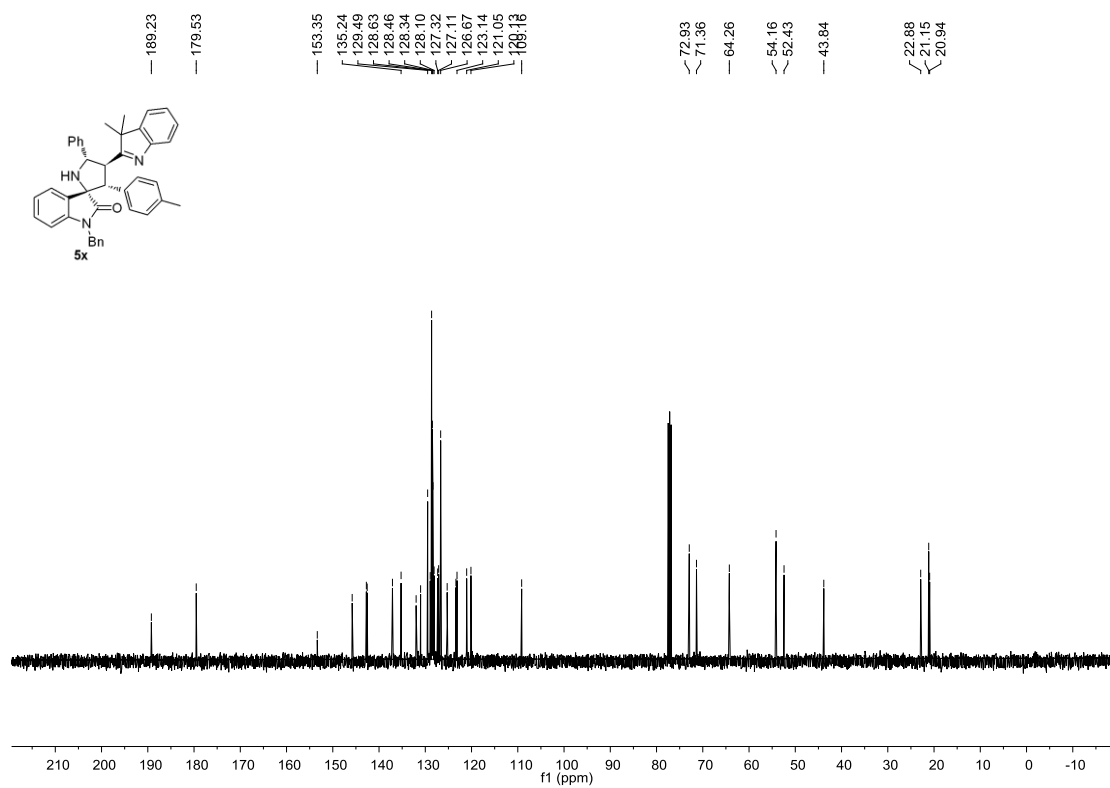
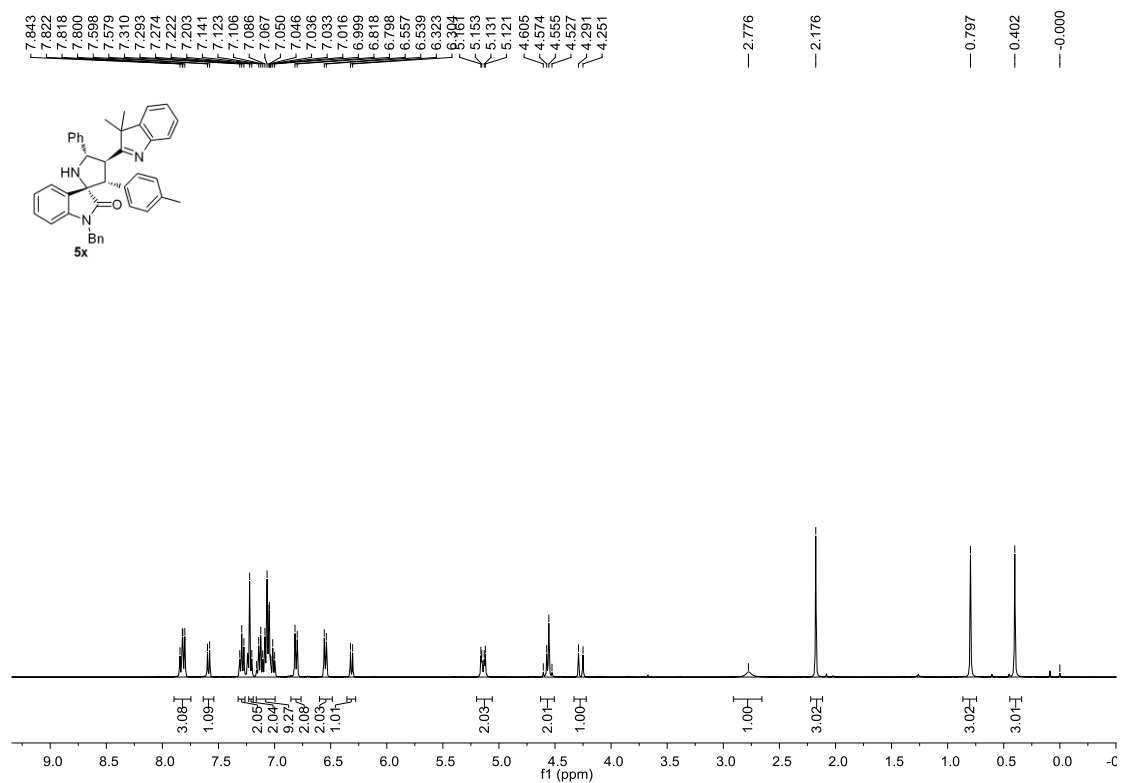


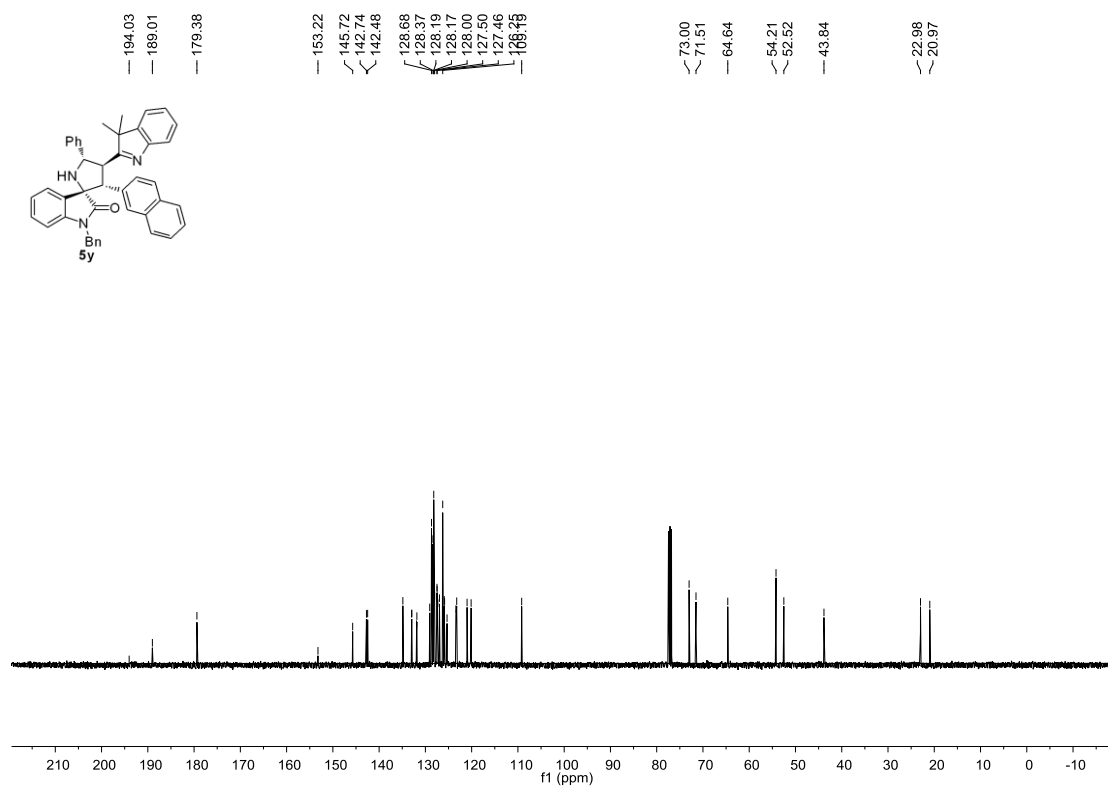
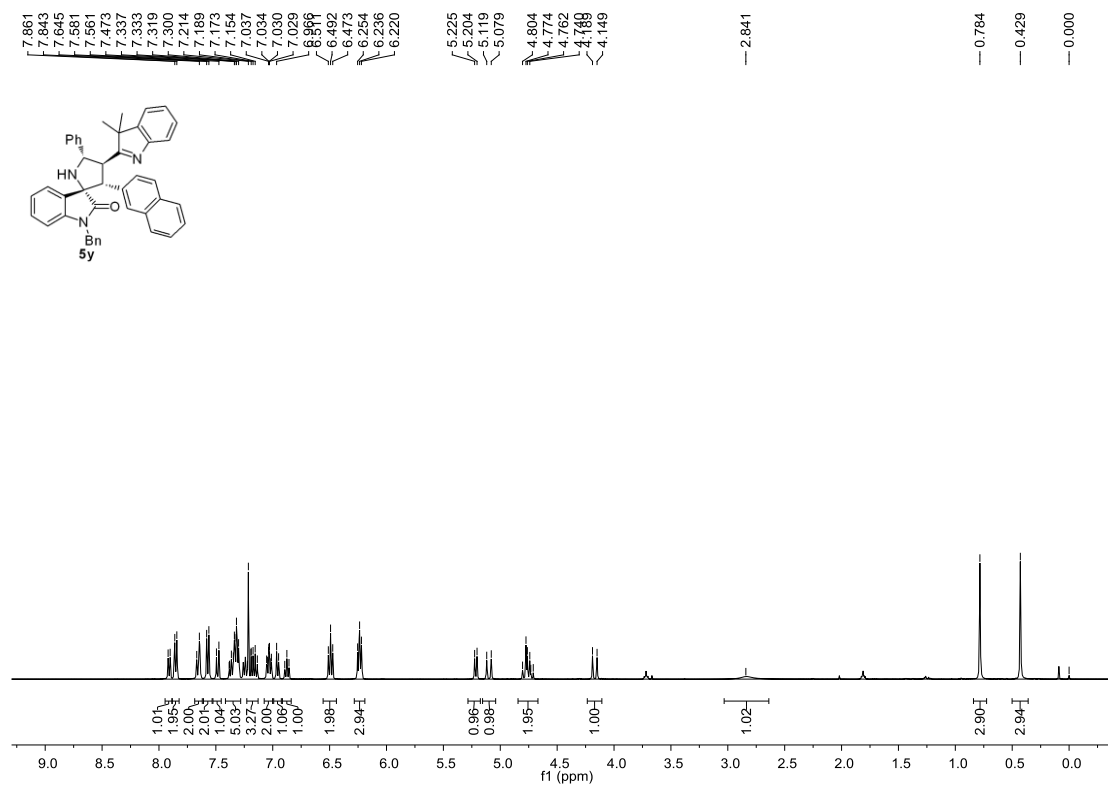


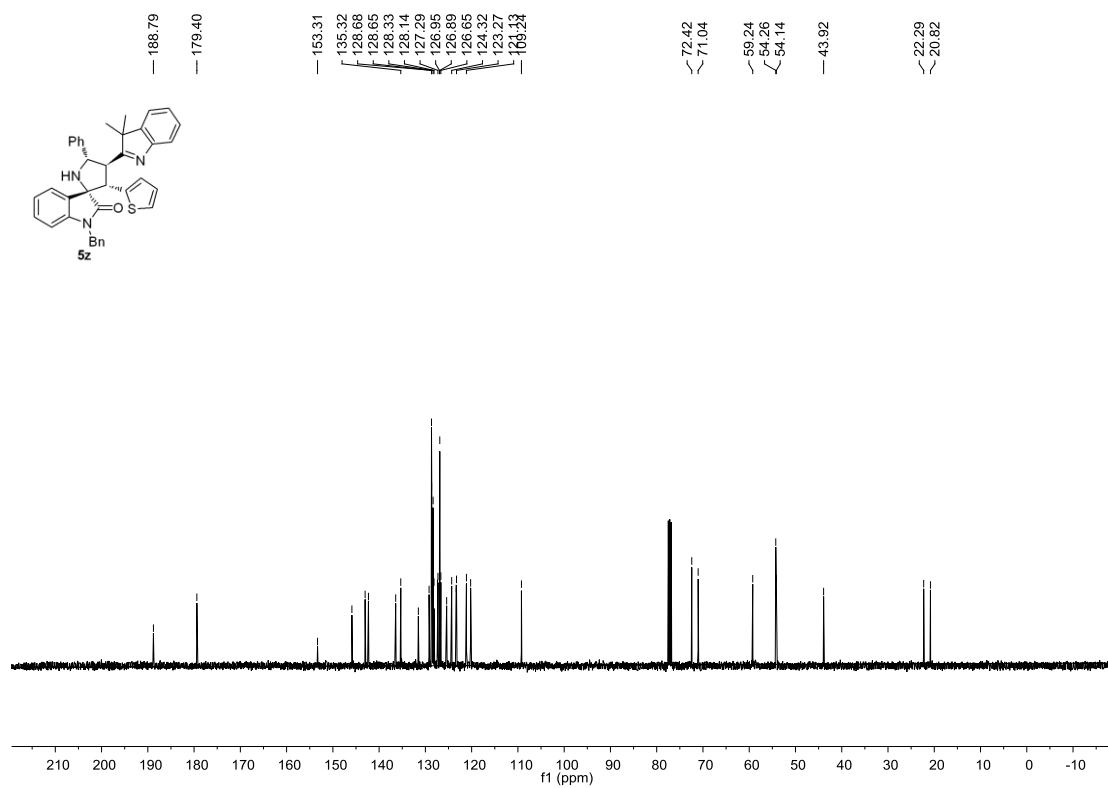
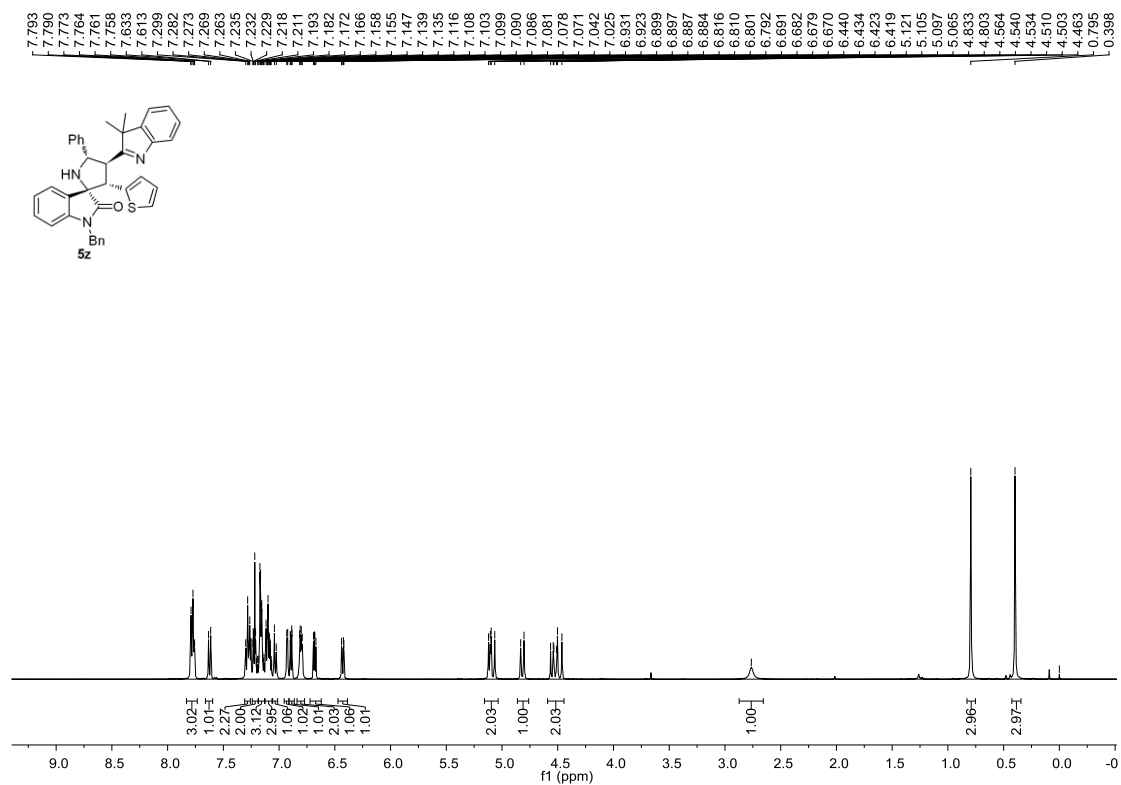


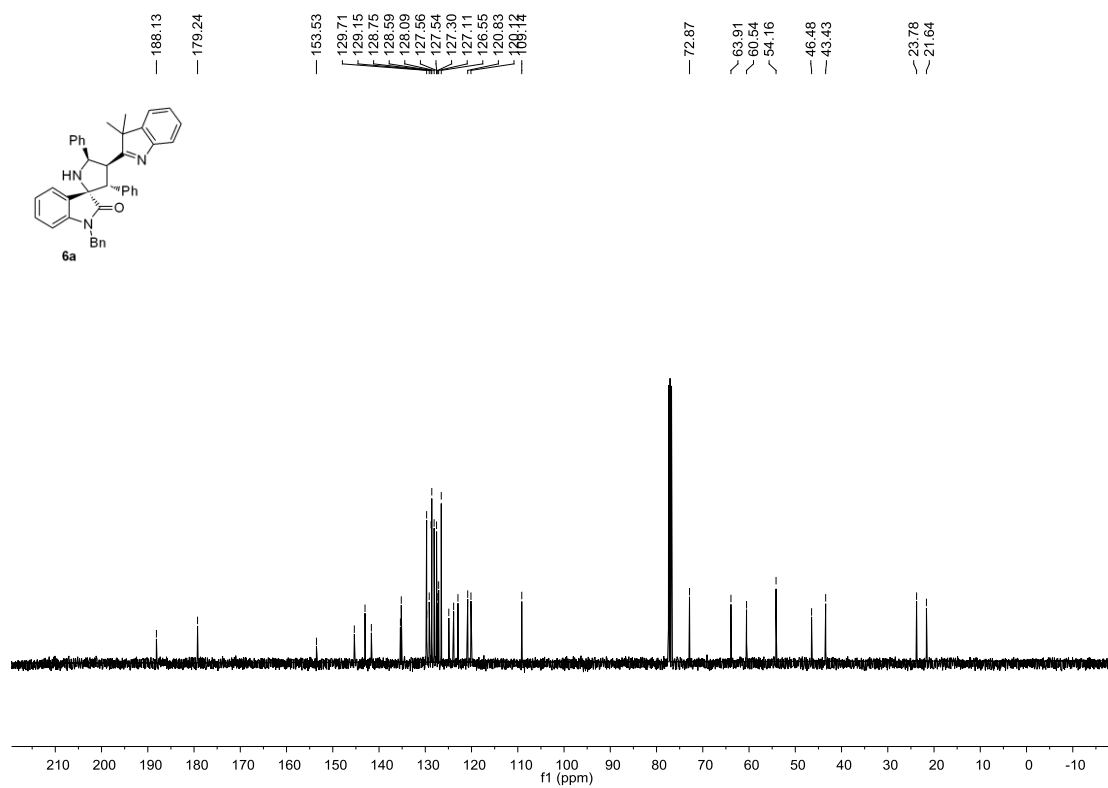
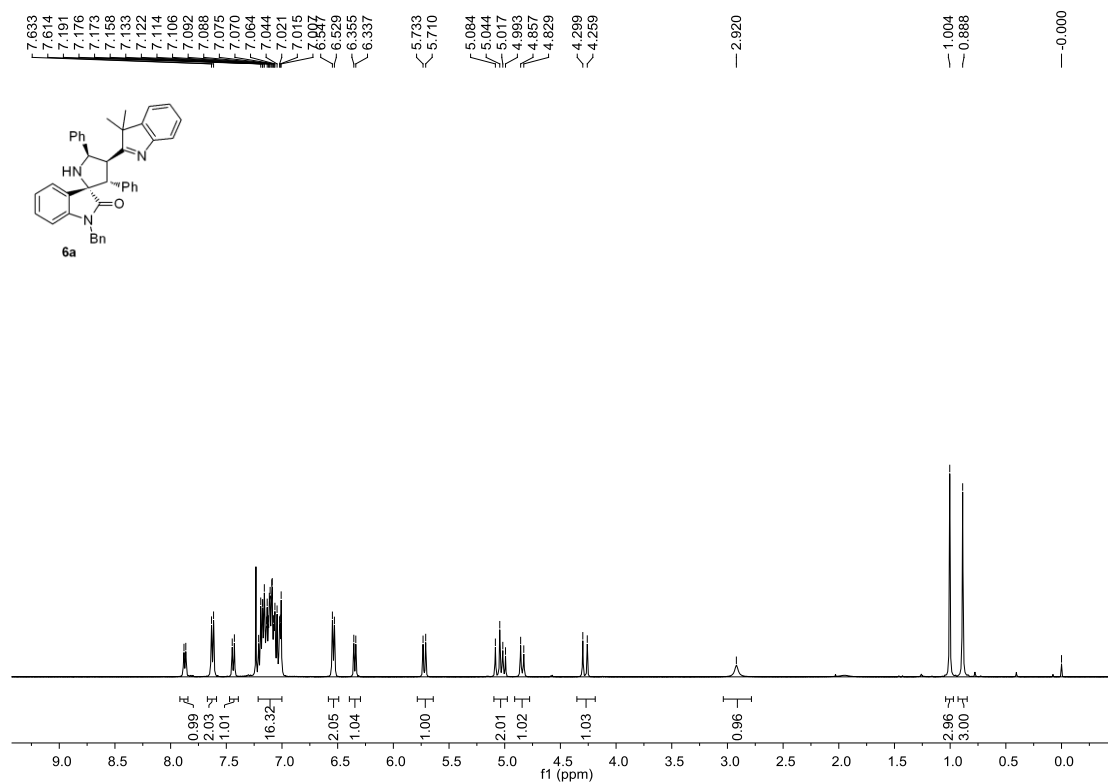


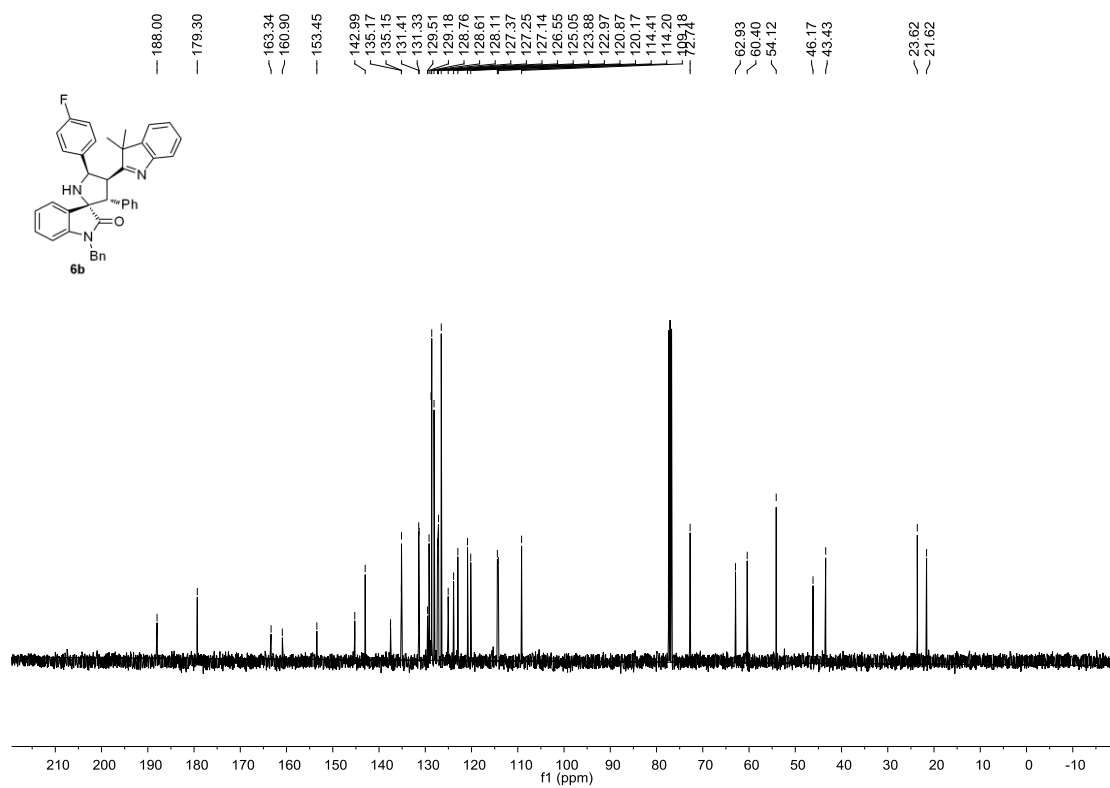
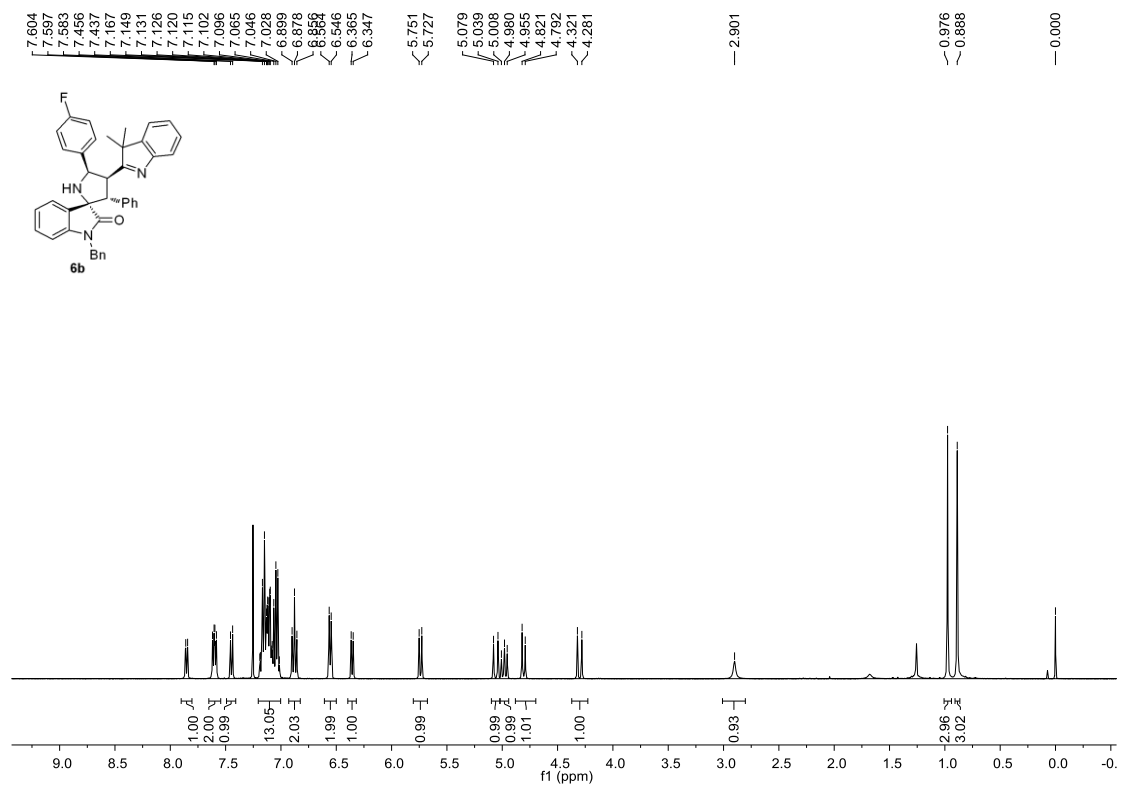


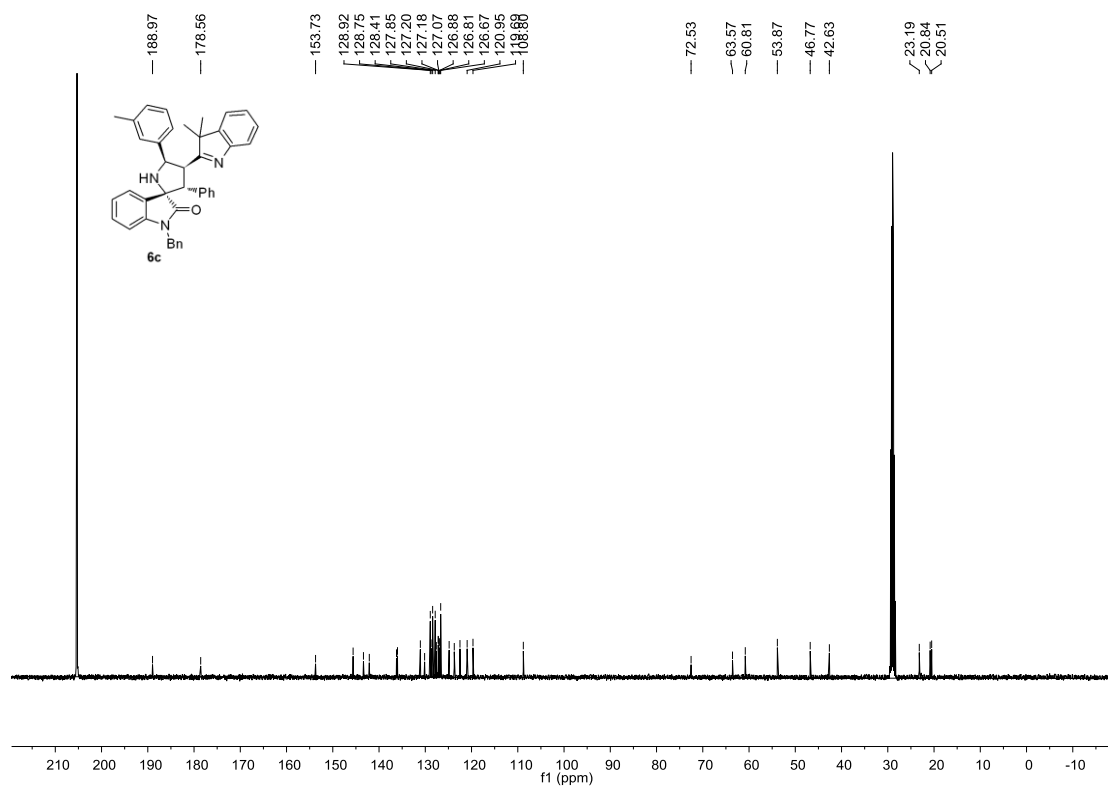
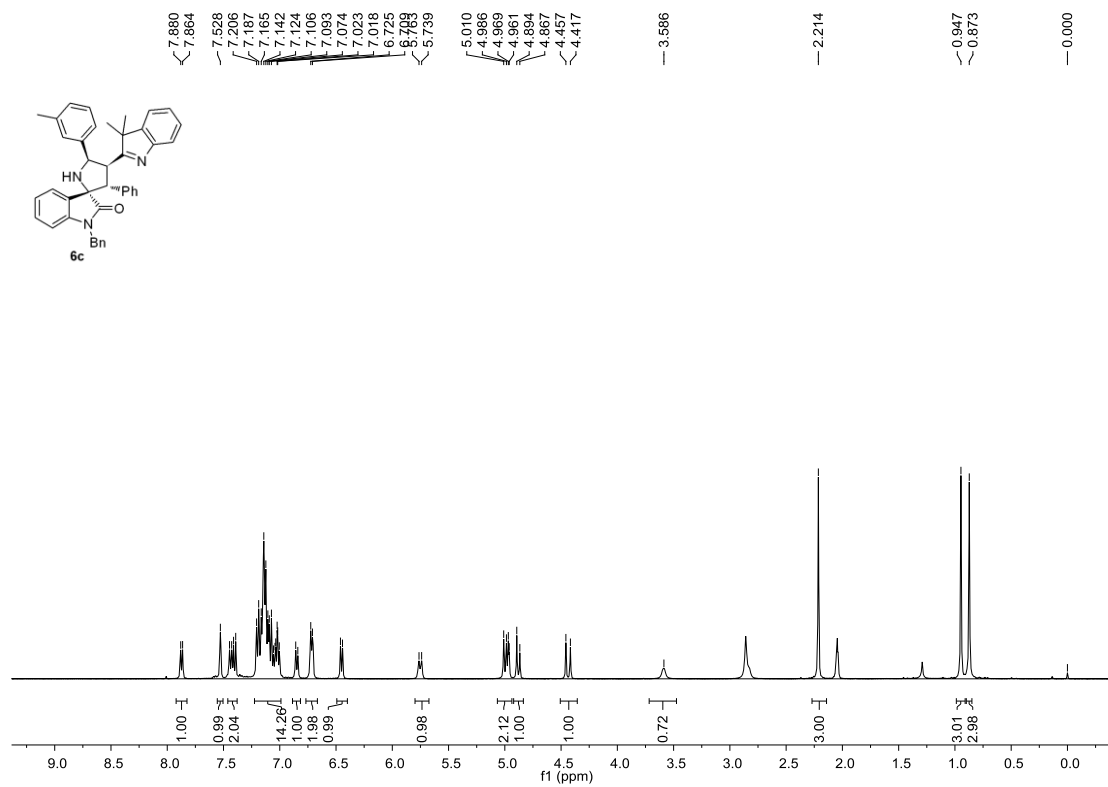


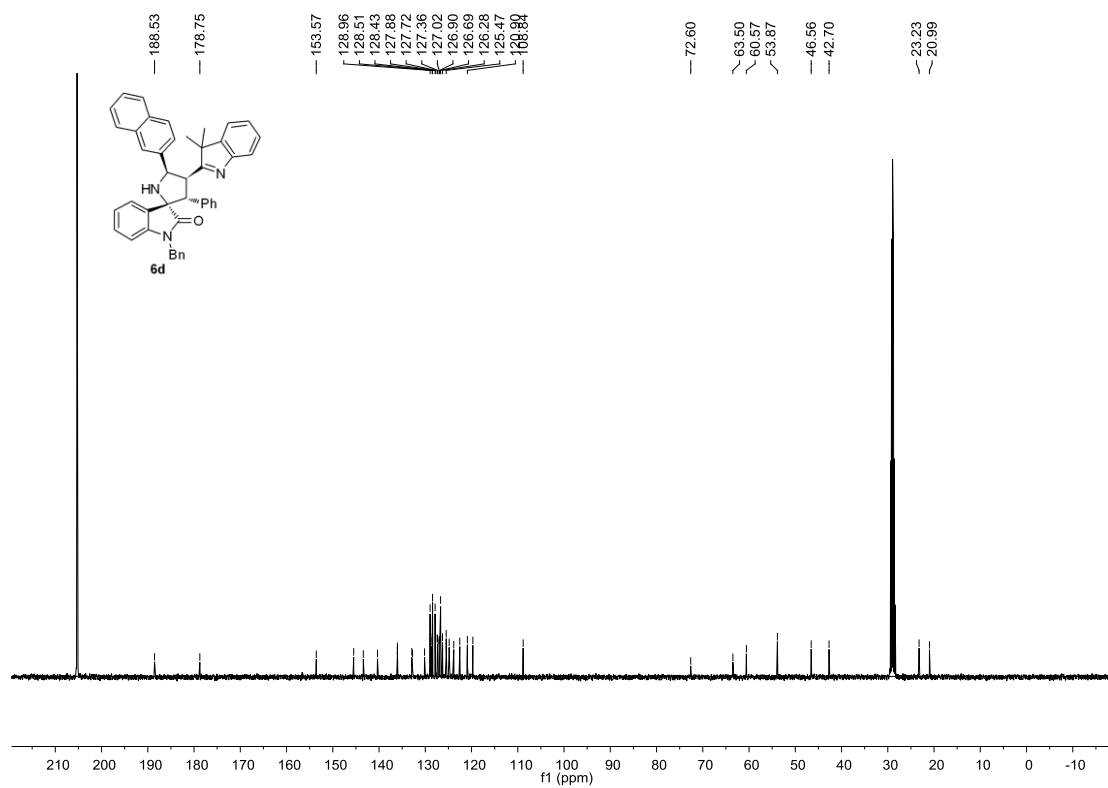
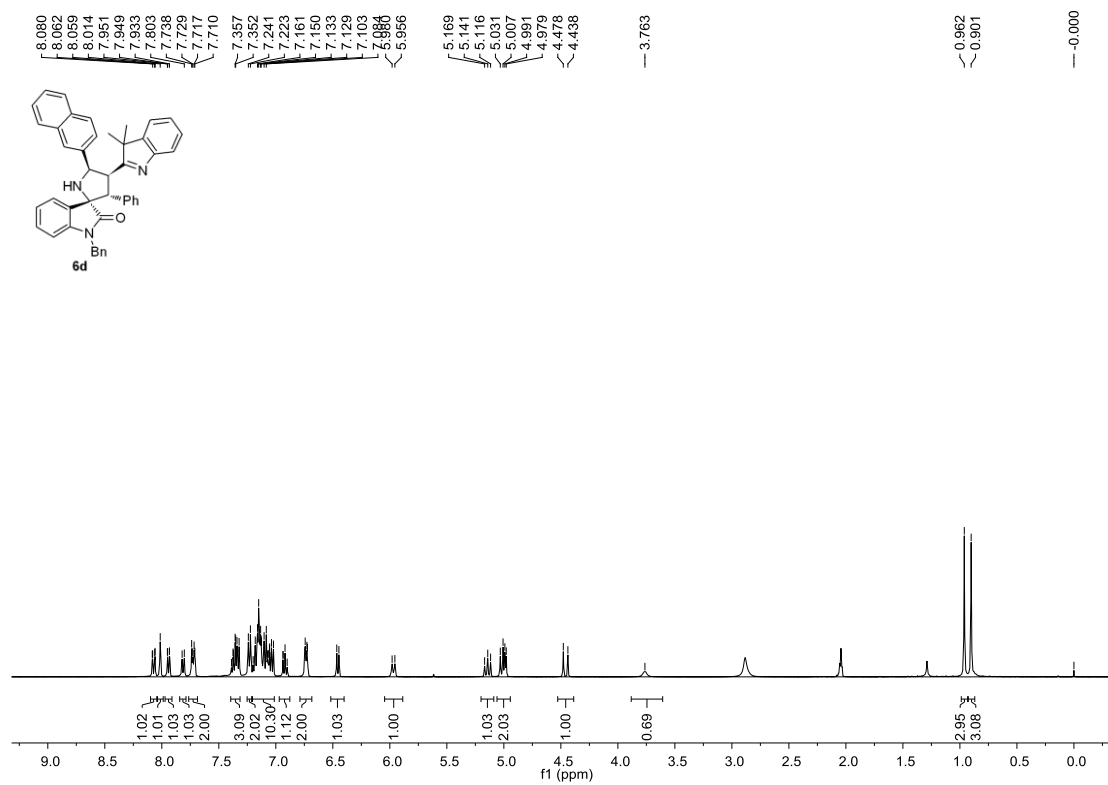


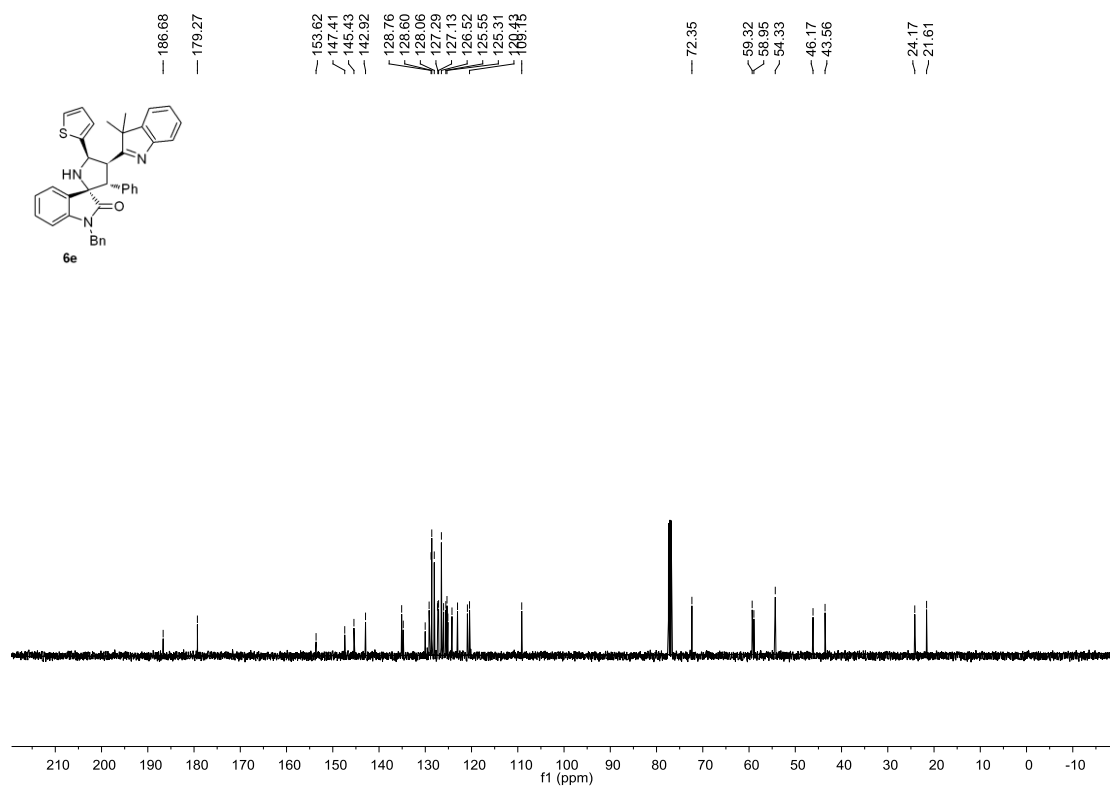
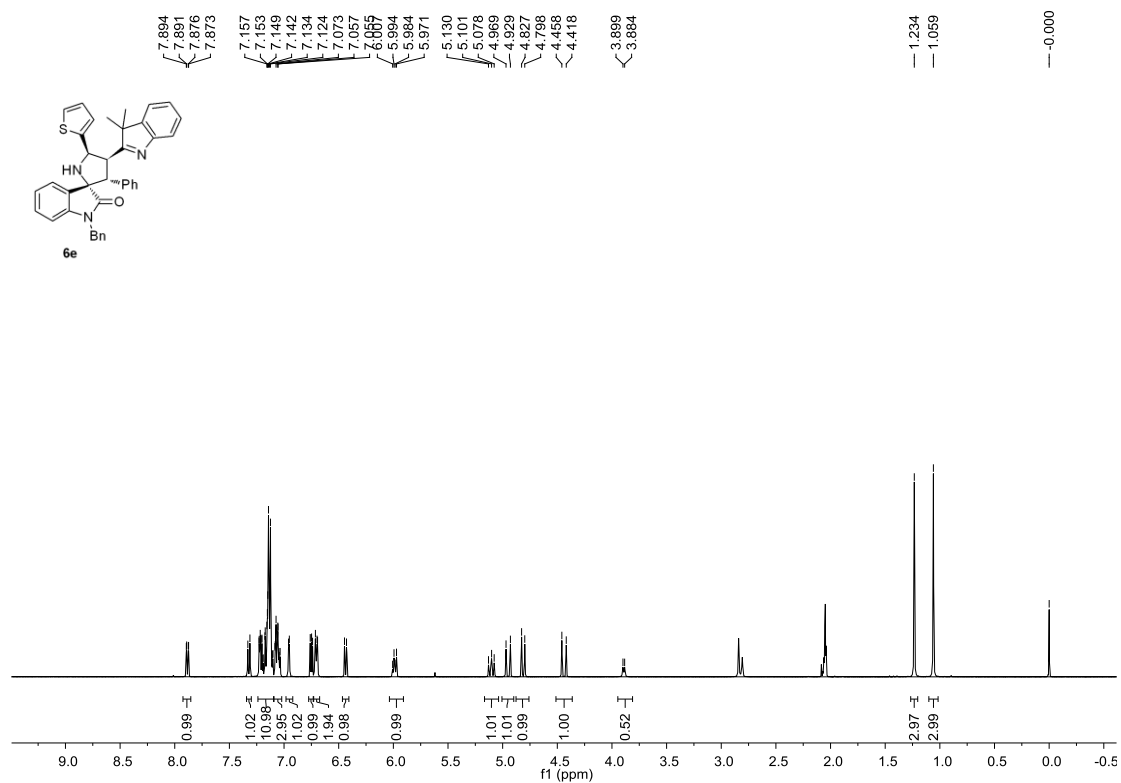


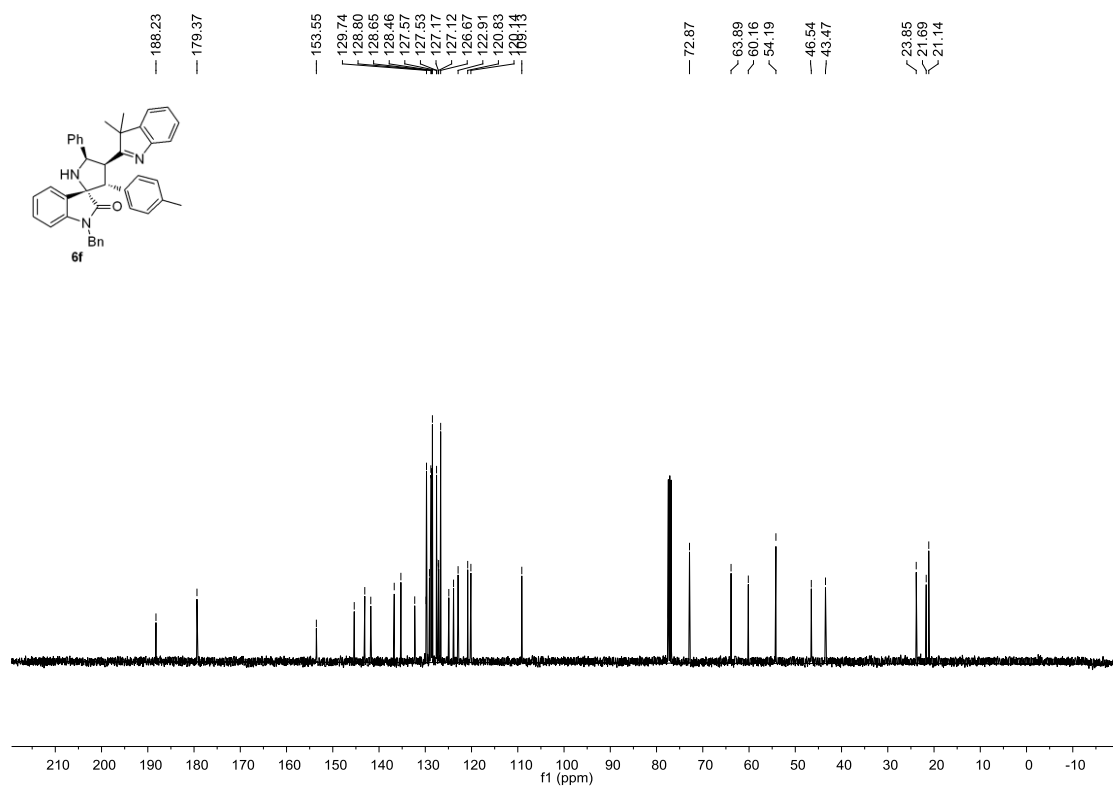
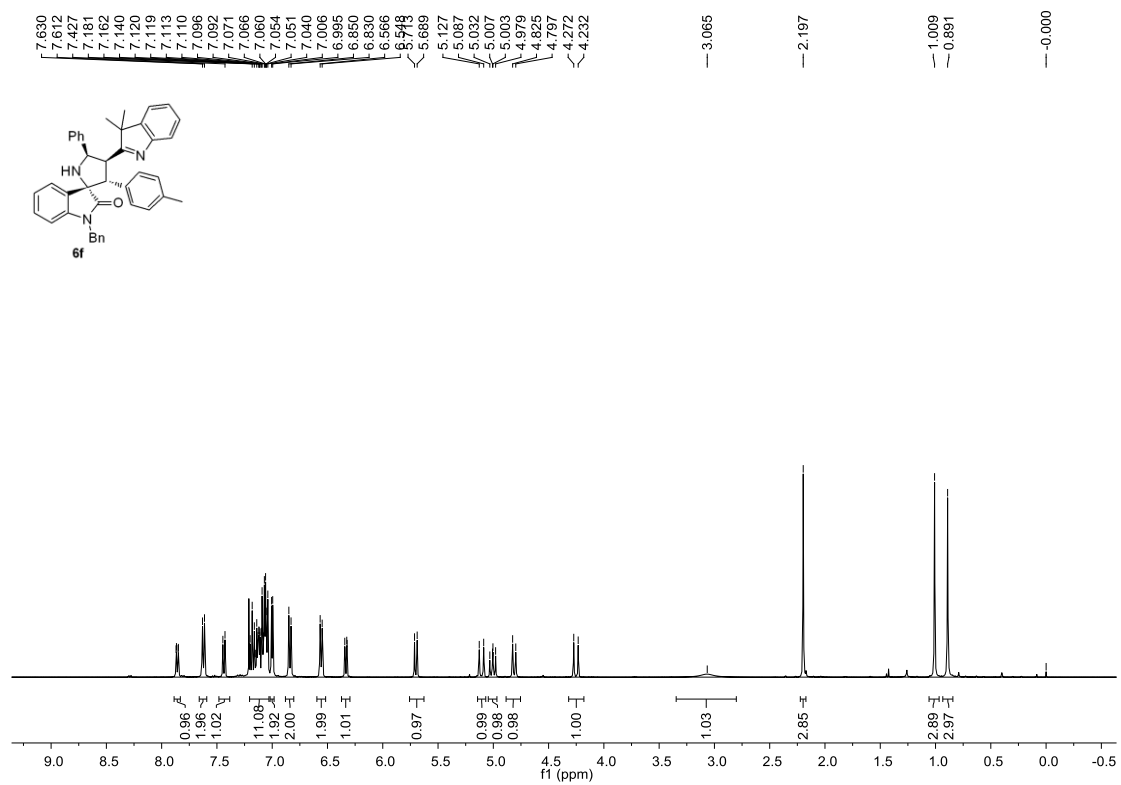


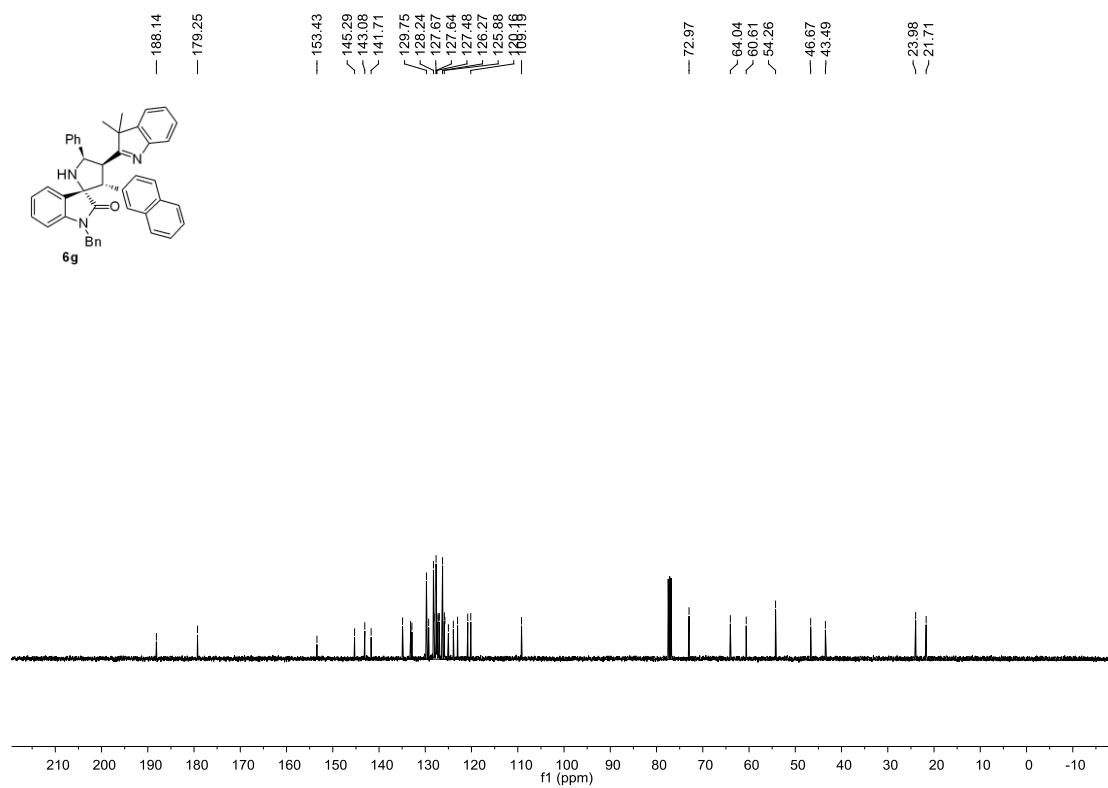
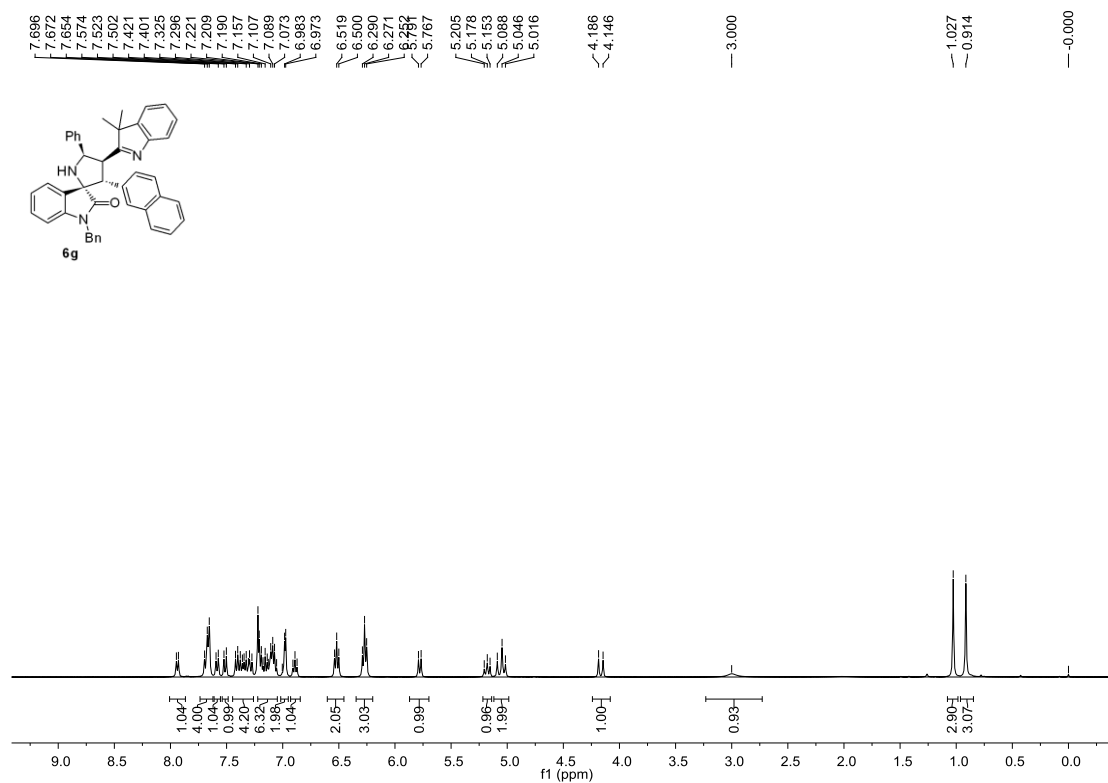


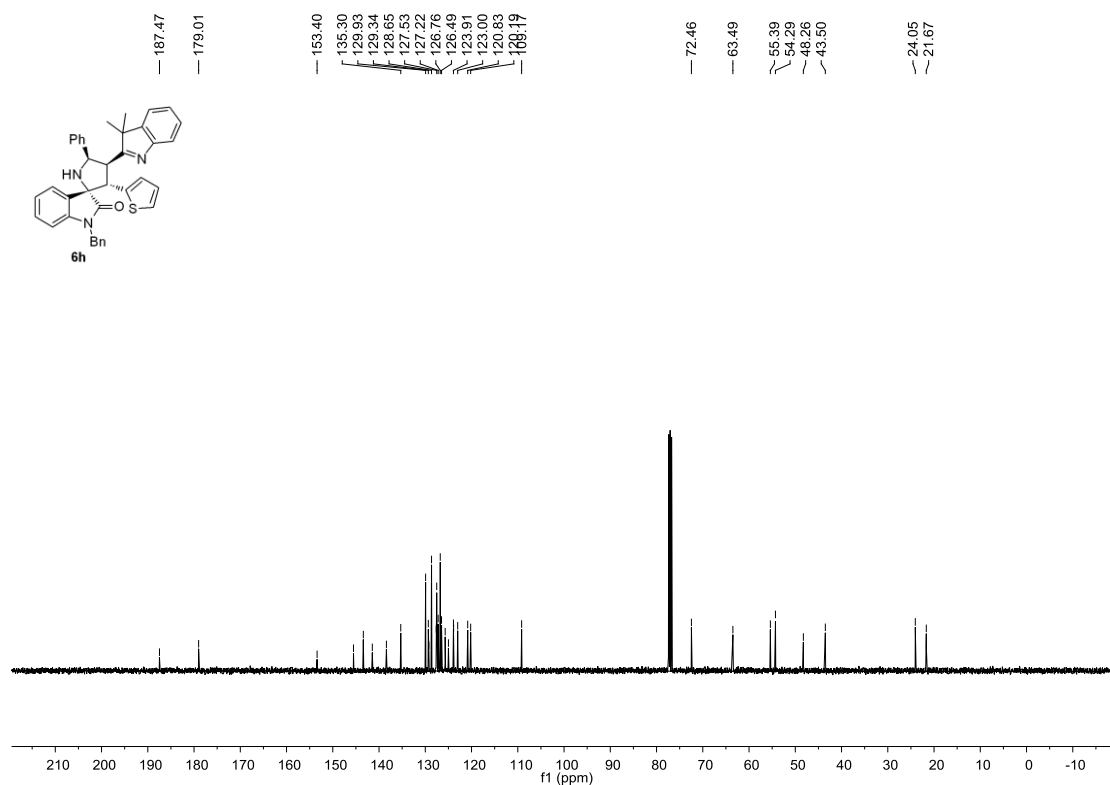
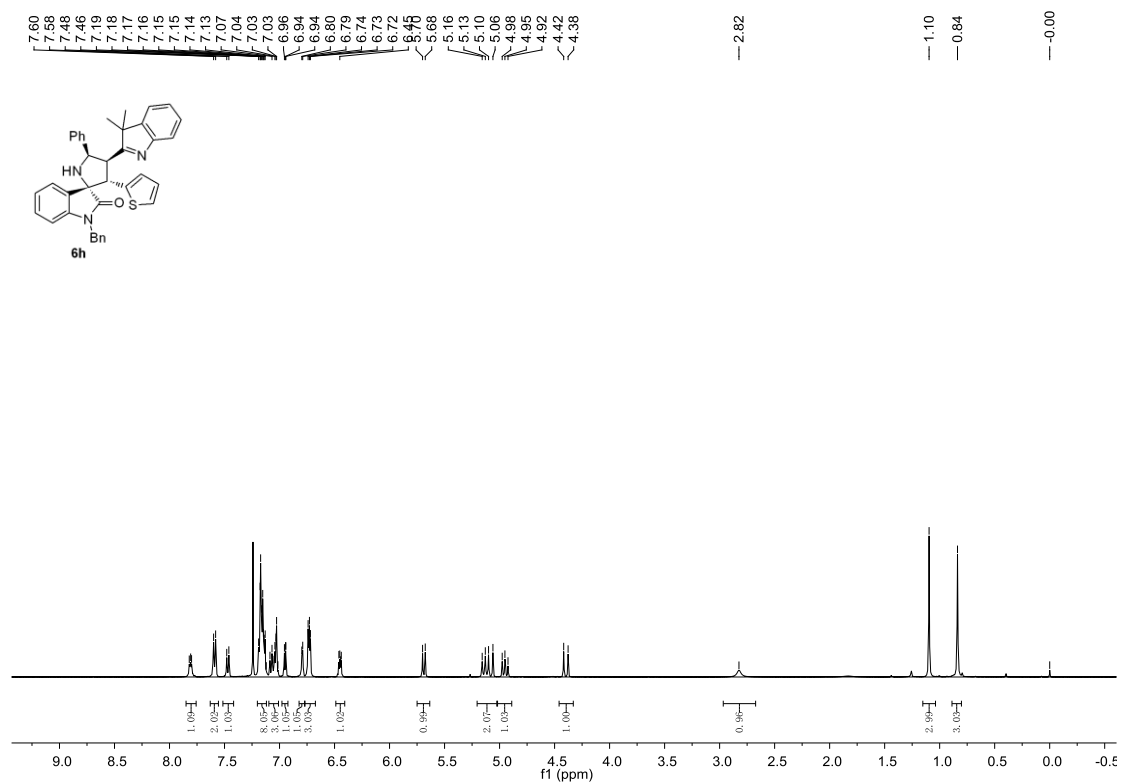


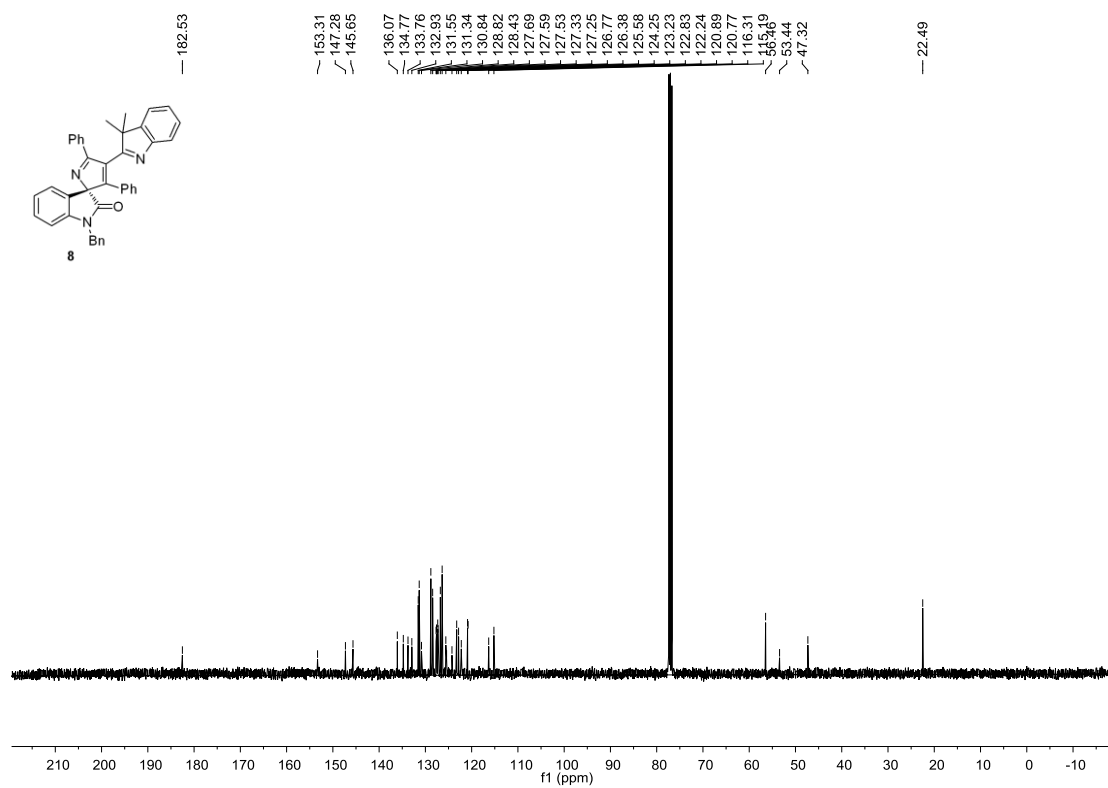
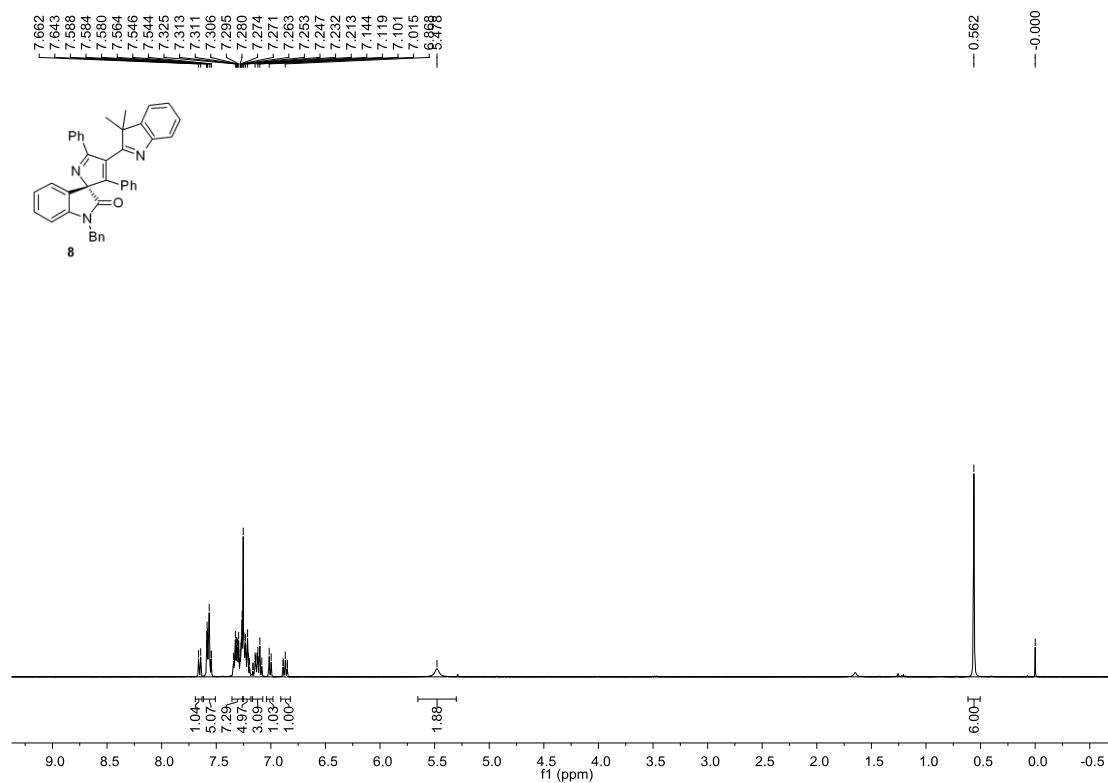


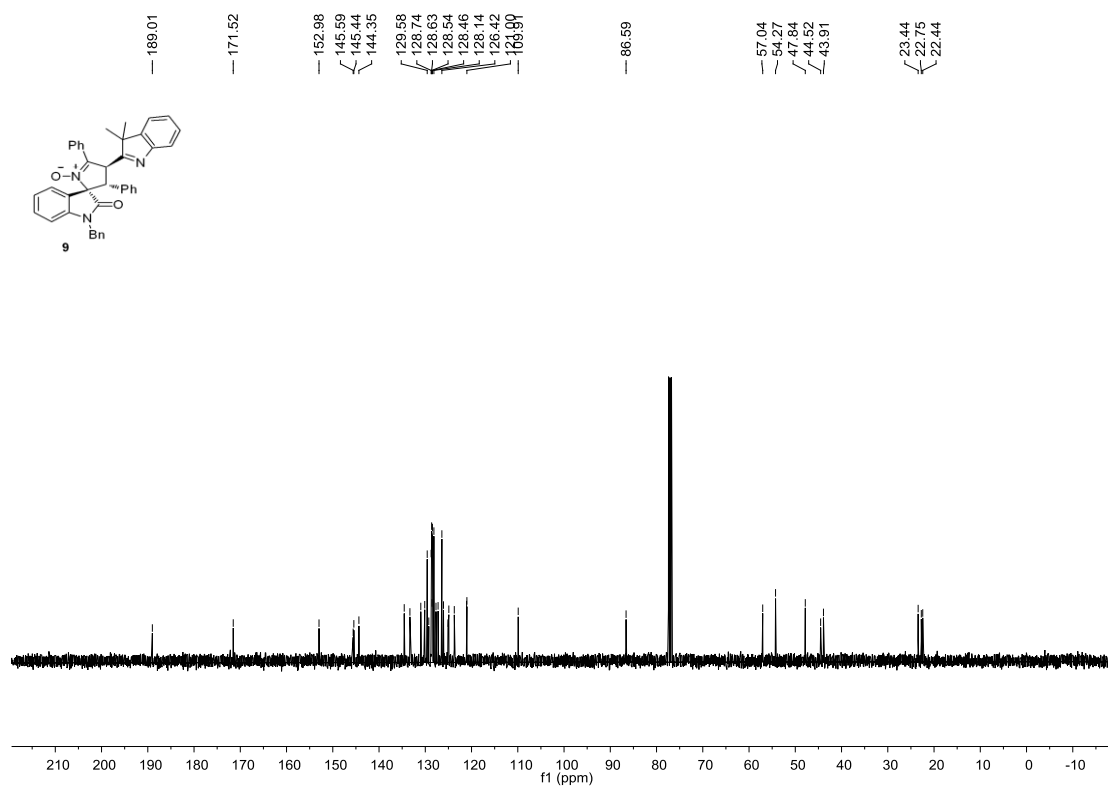
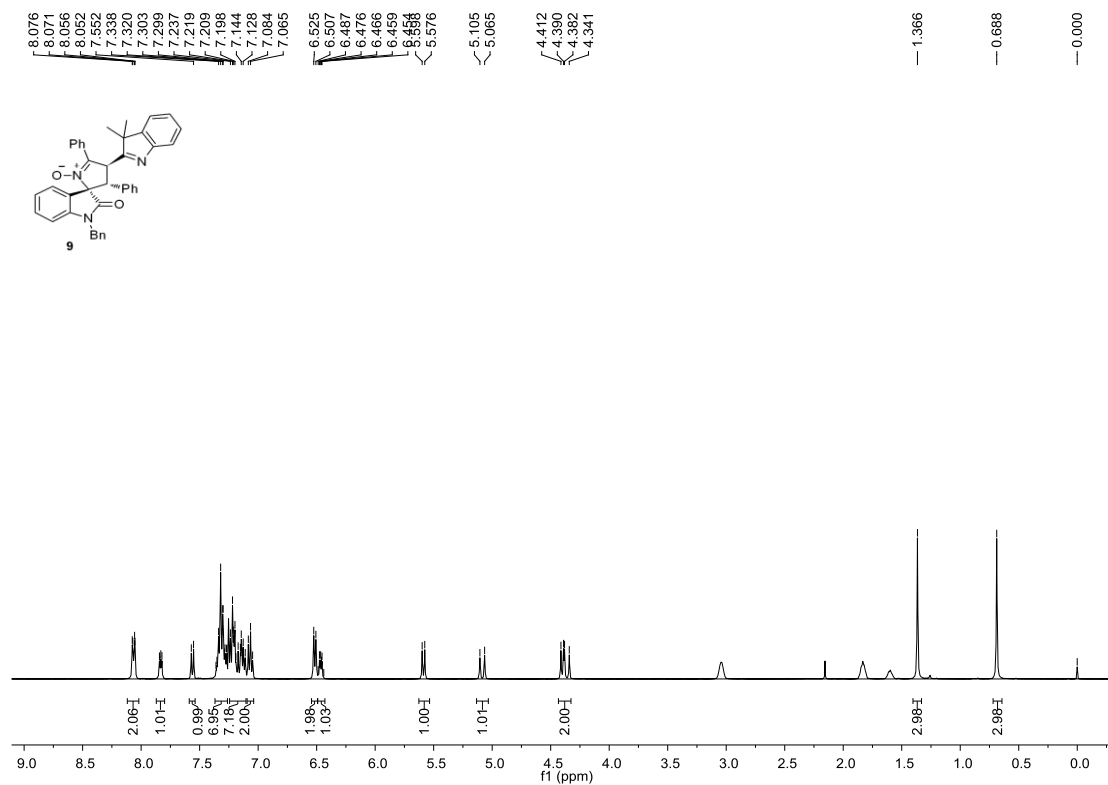








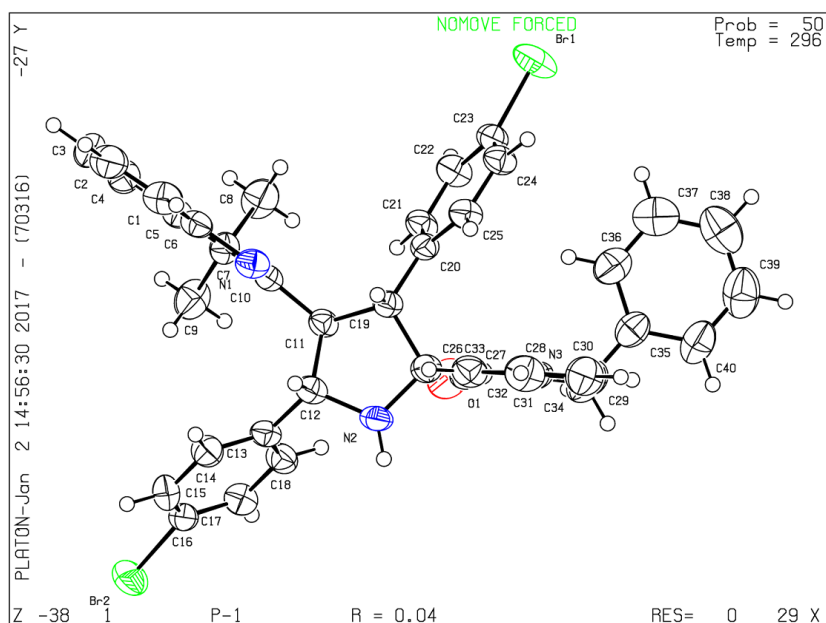




2. Crystal data of 5e and 6a

For 5e:

Empirical formula C₄₀ H₃₄ Br₂ N₃ O
Formula weight 732.52
Temperature (K) 296(2)
Crystal system triclinic
Space group P-1
a (Å) 11.2235(2)
b (Å) 12.7822(3)
c (Å) 13.1058(3)
 $\alpha(^{\circ})$ 62.5490(10)
 $\beta(^{\circ})$ 83.9160(10)
 $\gamma(^{\circ})$ 87.4370(10)
Volume (Å³) 1659.06(6)
Z 2
D_{calc} (g cm³) = 1.466
 μ (mm⁻¹) = 2.480
F(000) = 746
Theta range for data collection 1.76 to 25.00
Index ranges -13 ≤ h ≤ 13, -13 ≤ k ≤ 15, -15 ≤ l ≤ 15
Reflections collected 22697
Independent reflections 5857 [R(int) = 0.0352]
Data/restraints/parameters 5857/0/415
GOF (on F²) 0.930
Final R indexes [I ≥ 2σ(I)] R₁ = 0.0391, wR₂ = 0.1211
Final R indexes [all data] R₁ = 0.0568, wR₂ = 0.1311
Largest diff. peak and hole (e Å⁻³) 0.764/-0.635



For 6a:

Empirical formula C_{14.67} H_{14.33} N O
Formula weight 220.60
Temperature (K) 296(2)
Crystal system triclinic
Space group P-1
a (Å) 9.9514(2)
b (Å) 12.1514(3)
c (Å) 16.6468(3)
 $\alpha(^{\circ})$ 109.6120(10)
 $\beta(^{\circ})$ 93.8450(10)
 $\gamma(^{\circ})$ 106.0740(10)
Volume (Å³) 1793.39(7)
Z 6
D_{calc} (g cm⁻³) = 1.226
 μ (mm⁻¹) = 0.077
F (000) = 704
Theta range for data collection 1.85 to 24.99
Index ranges -11 ≤ h ≤ 11, -14 ≤ k ≤ 14, -19 ≤ l ≤ 19
Reflections collected 18484
Independent reflections 6299 [R(int) = 0.0199]
Data/restraints/parameters 6299/1/451
GOF (on F²) 1.082
Final R indexes [I ≥ 2σ(I)] R₁ = 0.0504, wR₂ = 0.1354
Final R indexes [all data] R₁ = 0.0590, wR₂ = 0.1425
Largest diff. peak and hole (e Å⁻³) 0.584/-0.609

