

Supplementary Materials

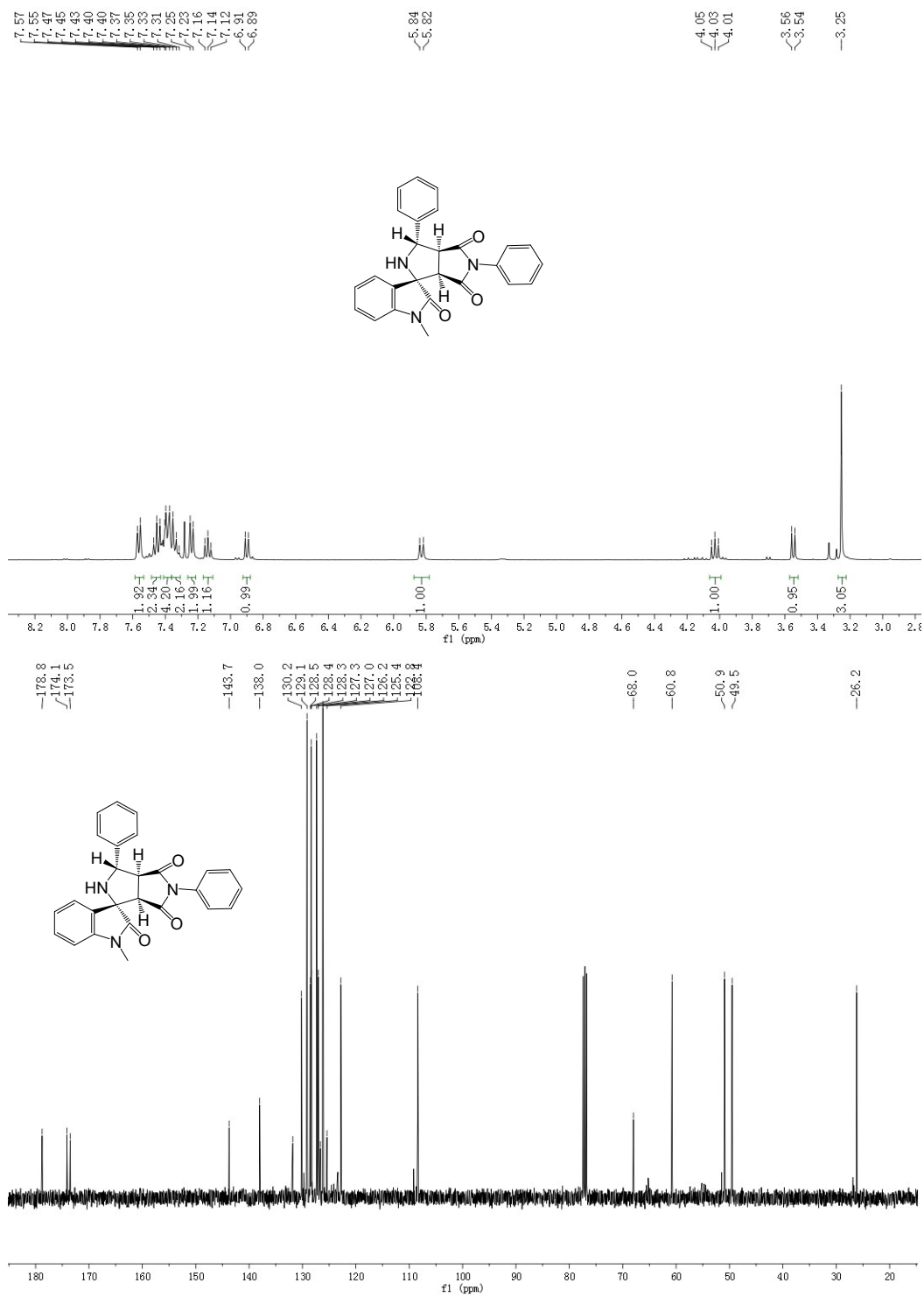
A facile one-pot construction of succimide-fused spiro[pyrrolidine-2,3'-oxindoles] via 1,3-dipolar cycloaddition involving 3-amino oxindoles and maleimides

Lunqiang Jin, Feng Liang*

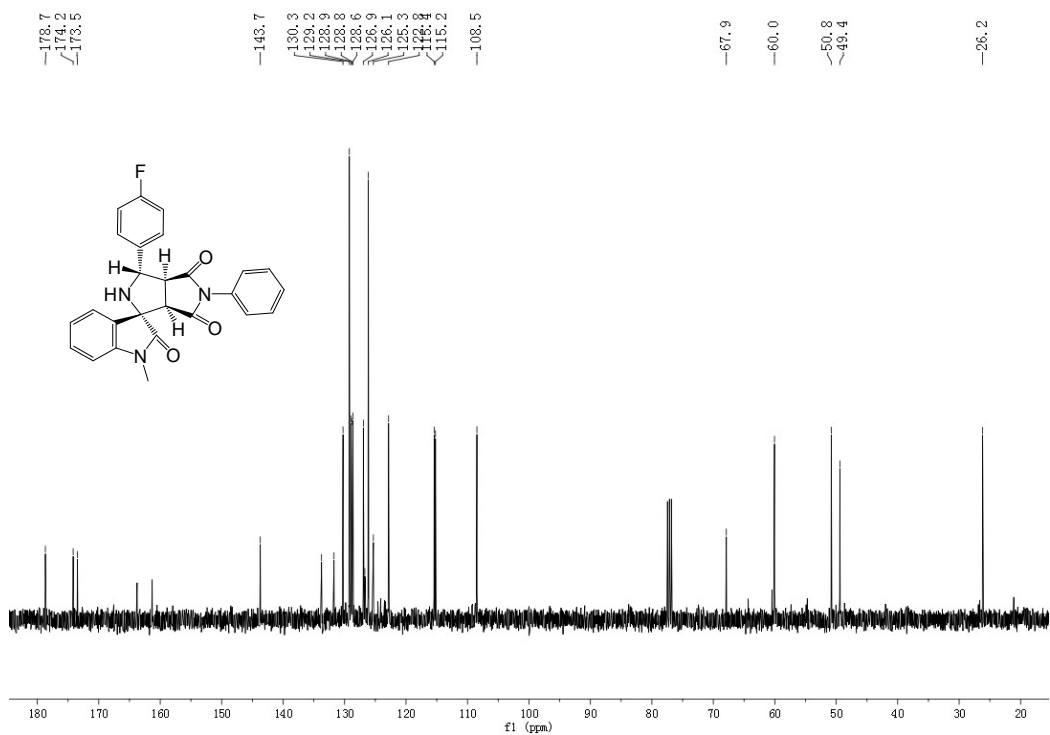
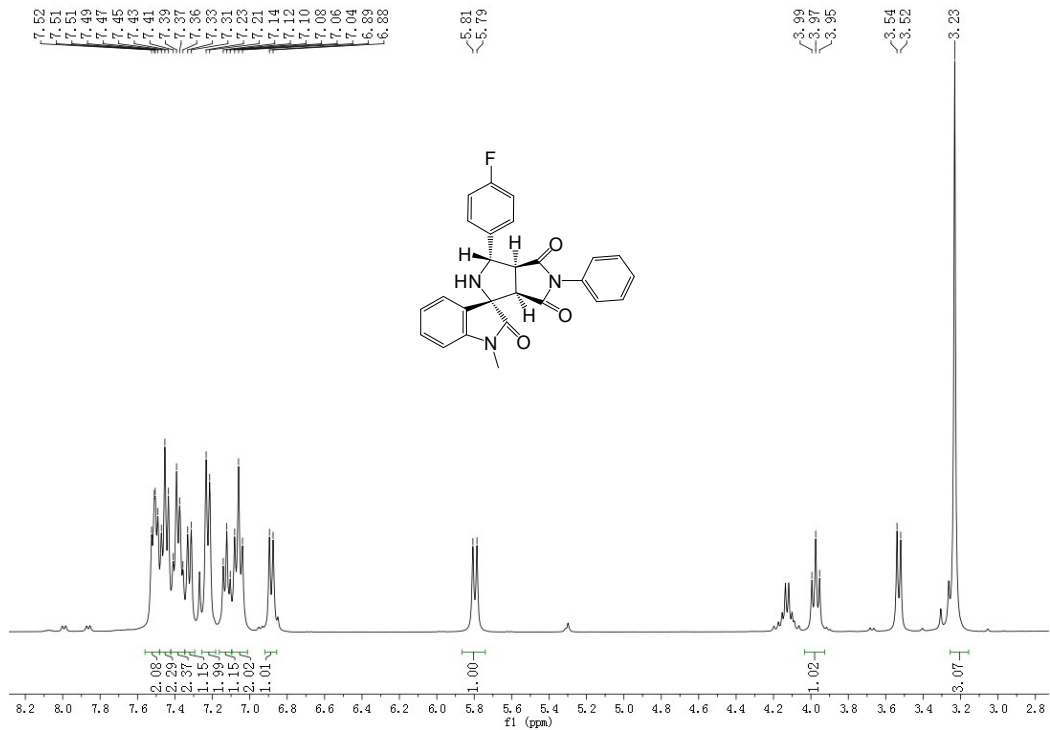
The State Key Laboratory of Refractories and Metallurgy, School of Chemistry &
Chemical Engineering, Wuhan University of Science and Technology, Wuhan 430081,
China

NMR Spectra

4a

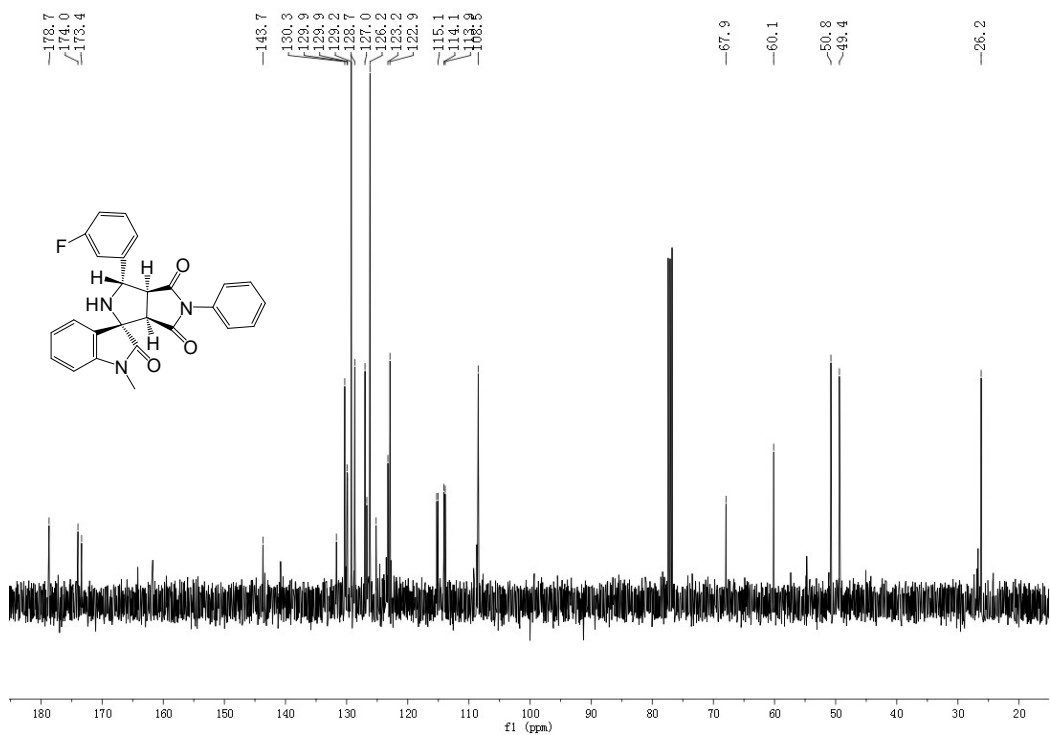


4b

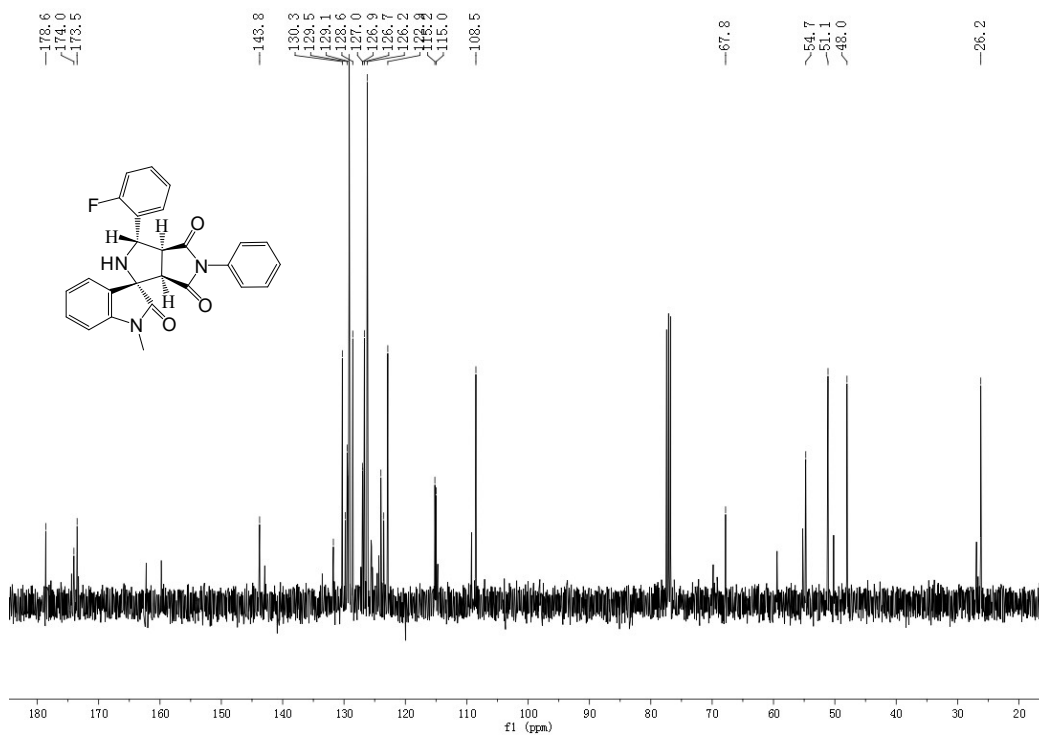
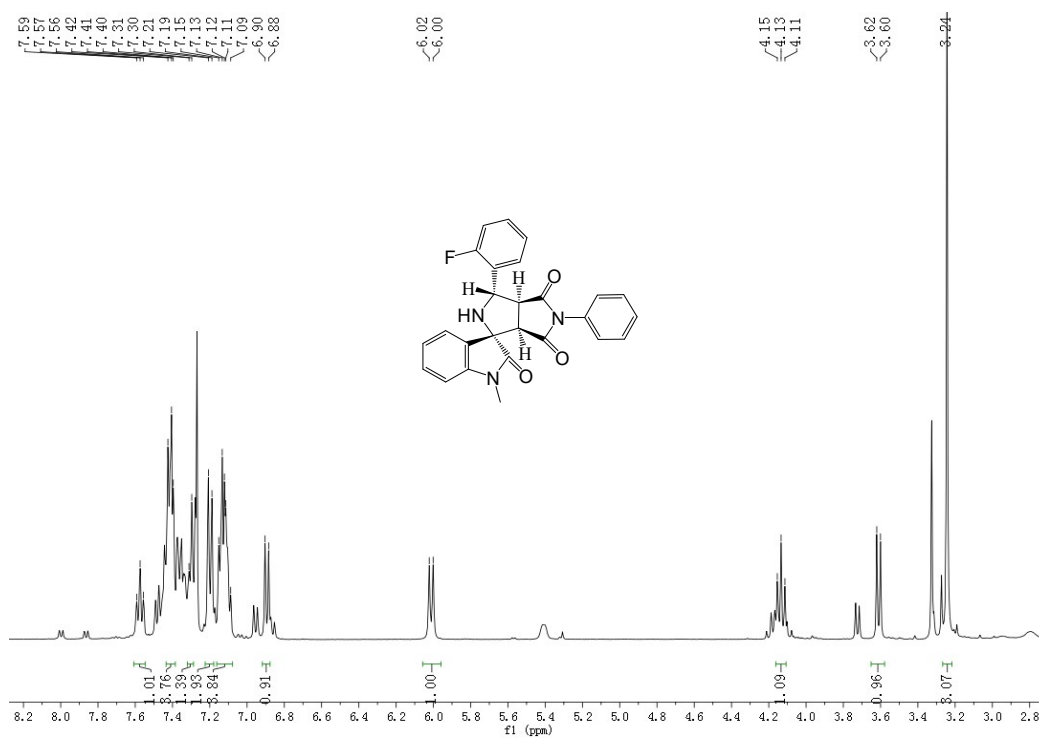


O=C1C(=O)N(c2ccccc2)[C@H]3[C@@H](C(=O)N3c4ccccc4F)[C@@H]1O

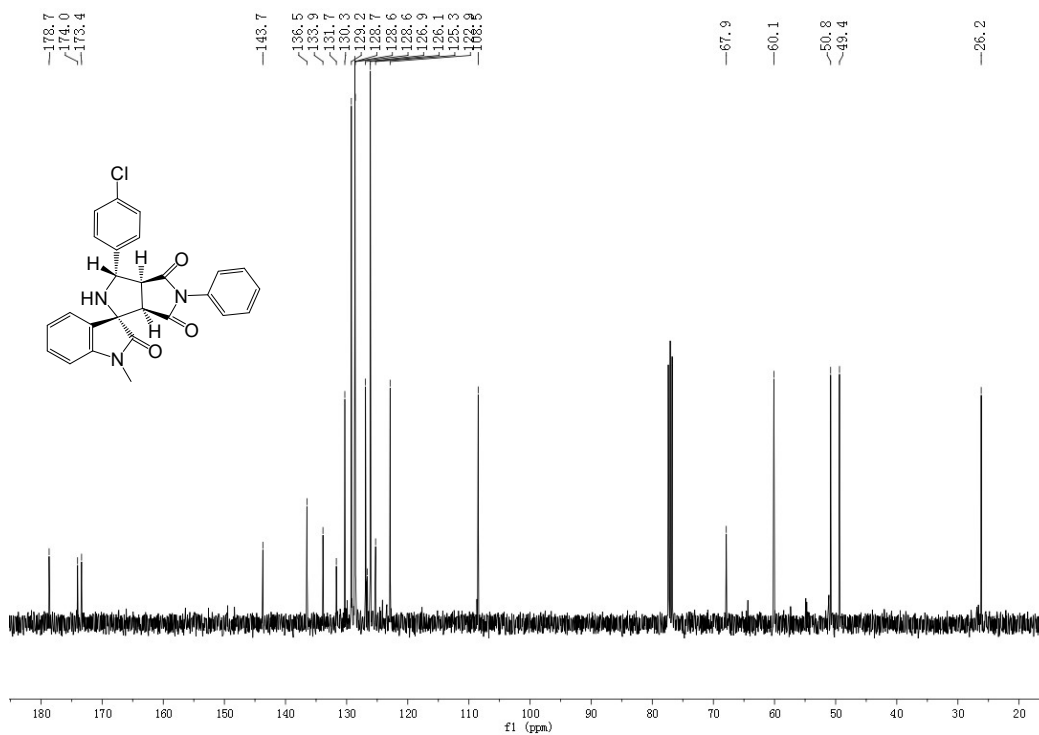
¹H NMR spectrum (CDCl₃) of (S)-1-((S)-2-oxo-2-phenyl-3-((S)-2-oxo-2-phenyl-3-oxo-1-phenyl-1H-indol-3-yl)propyl)pyrrolidine-2-one. The spectrum shows peaks from 1.6 to 7.5 ppm. Integration values are provided below the peaks: 1.00, 1.01, 0.97, 3.18. Chemical shift values are listed above the peaks: 7.47, 7.45, 7.43, 7.41, 7.39, 7.38, 7.36, 7.34, 7.32, 7.29, 7.23, 7.21, 7.15, 7.11, 7.10, 7.00, 6.99, 6.98, 6.90, 4.03, 4.01, 3.99, 3.55, 3.53, 3.24.



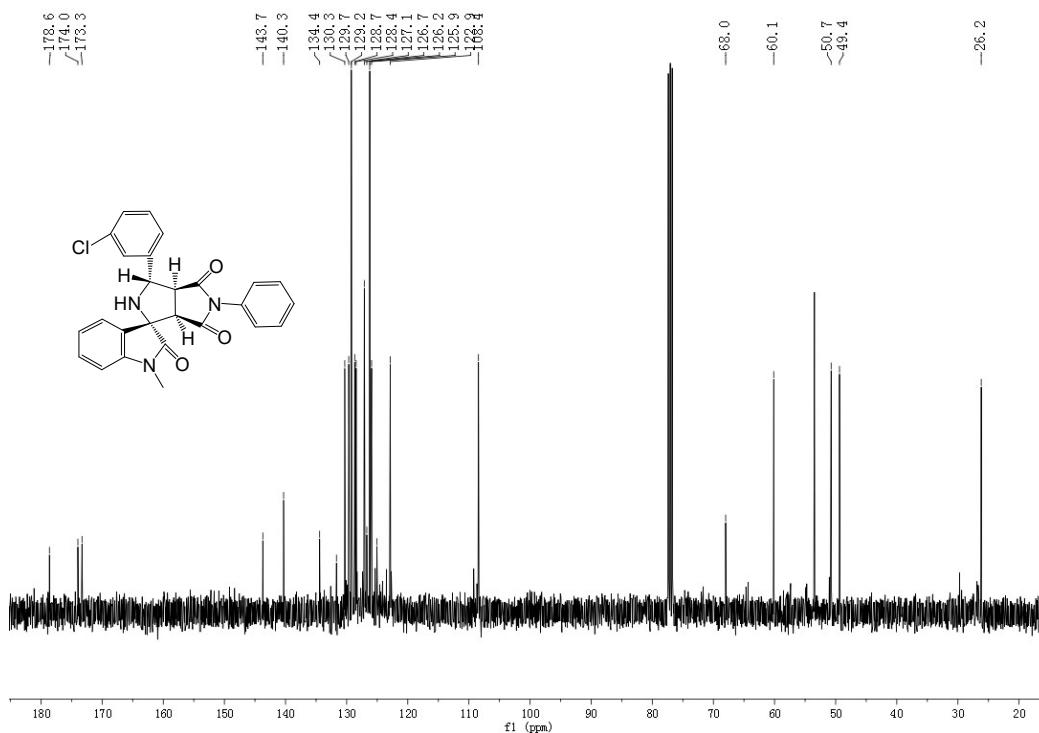
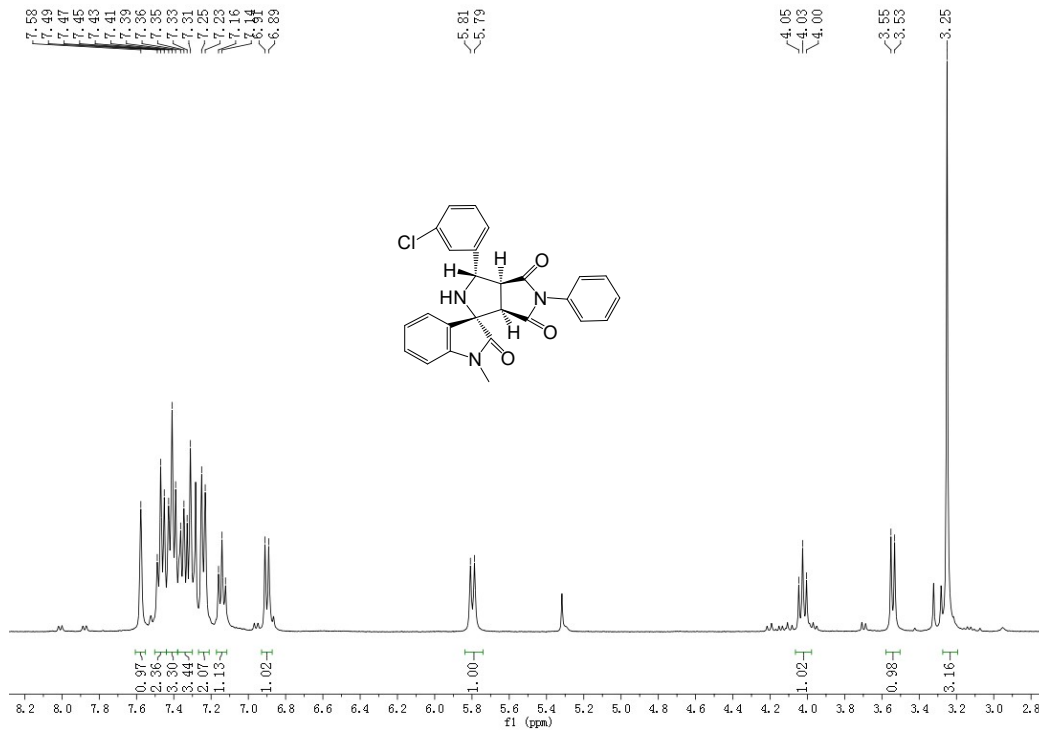
4d



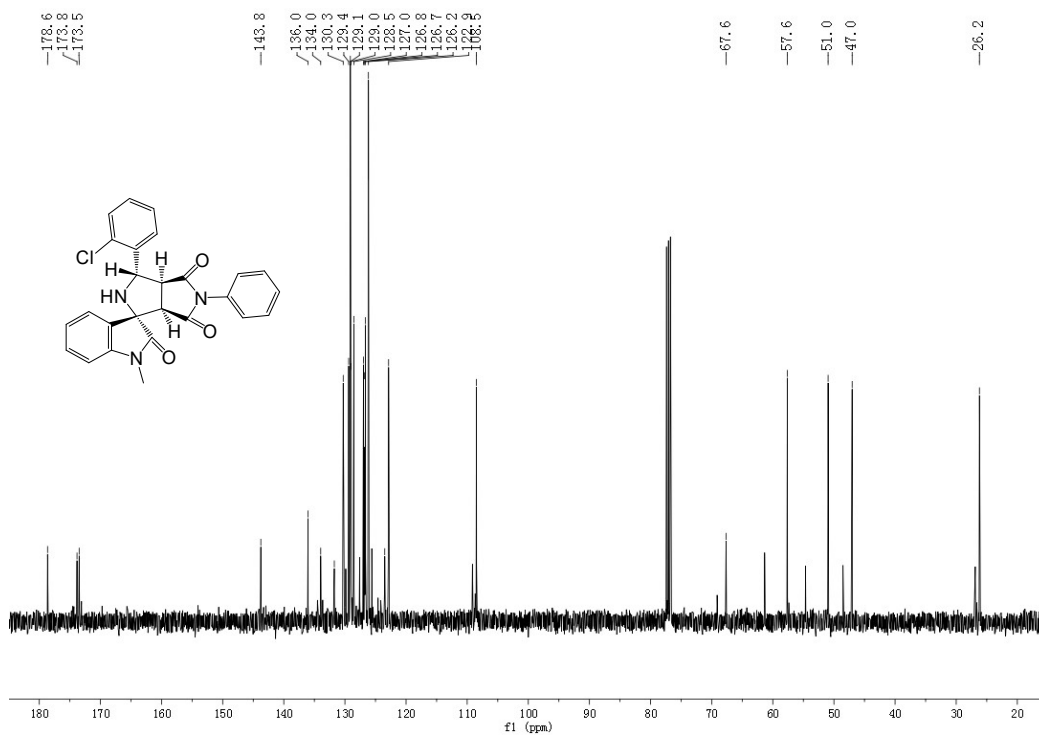
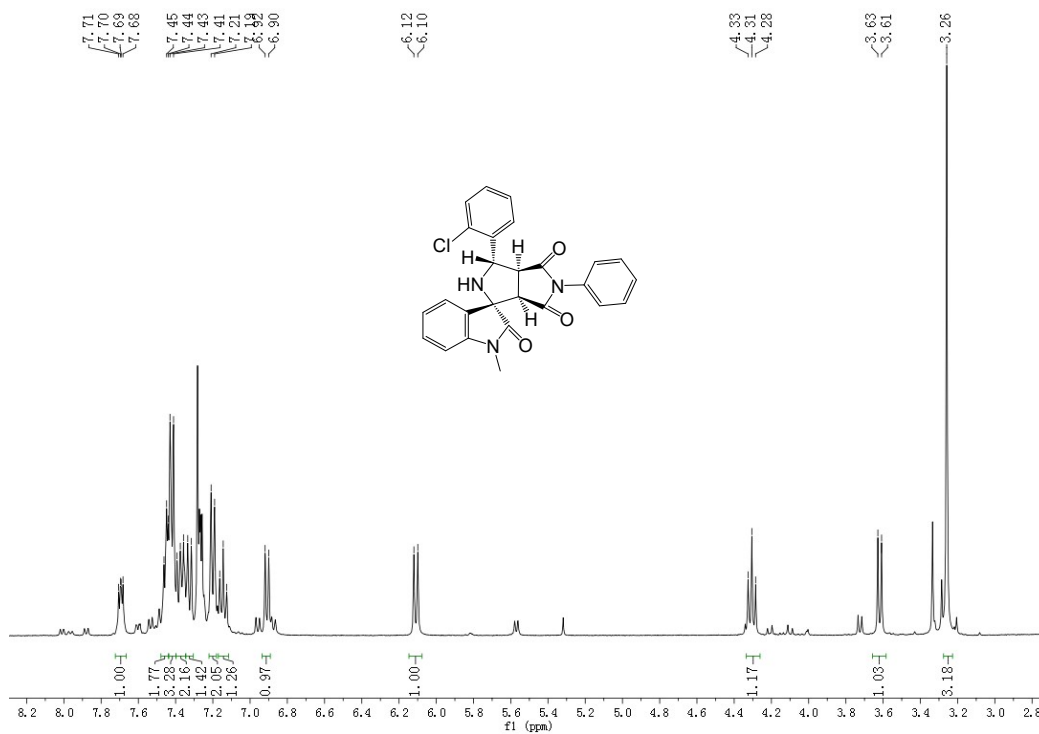
Chemical structure of compound 10 is shown above the spectrum. The structure is a complex molecule featuring a central carbon atom bonded to a chlorine atom, a phenyl ring, a benzimidazole ring, and a benzimidazole ring. The benzimidazole rings are substituted with a phenyl group and a benzimidazole ring. The spectrum shows peaks from 3.2 to 7.5 ppm. Integration values are provided below the baseline: 4.87, 4.21, 2.16, 1.33, 1.11, 1.00, 1.03, 1.10, and 3.11.



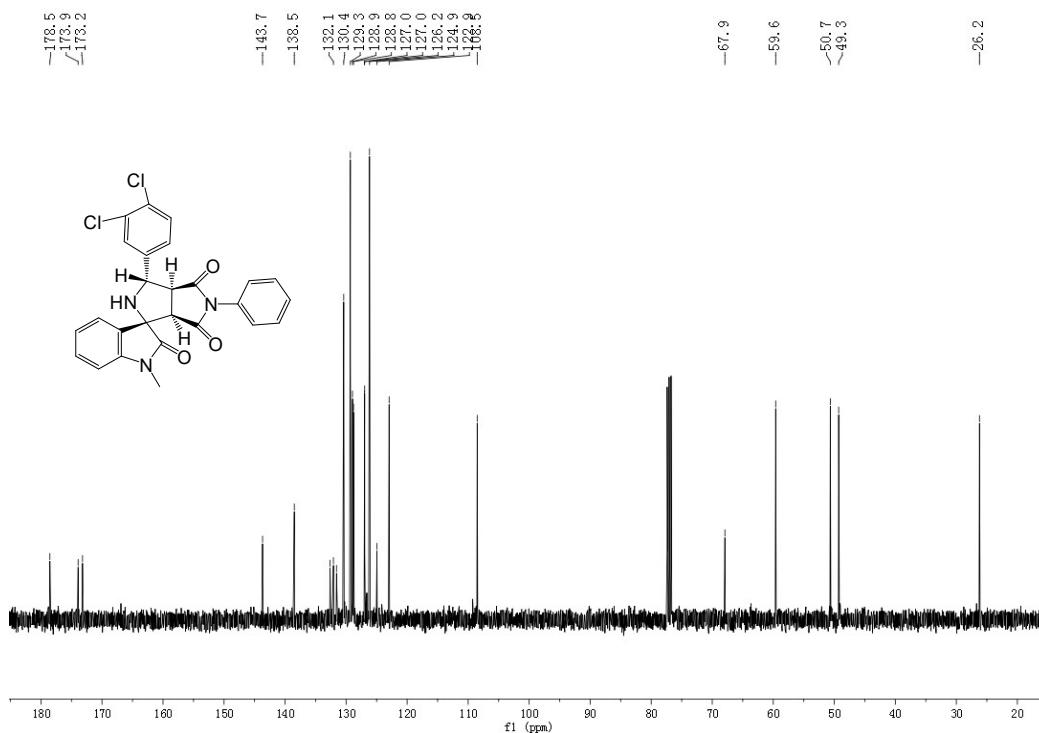
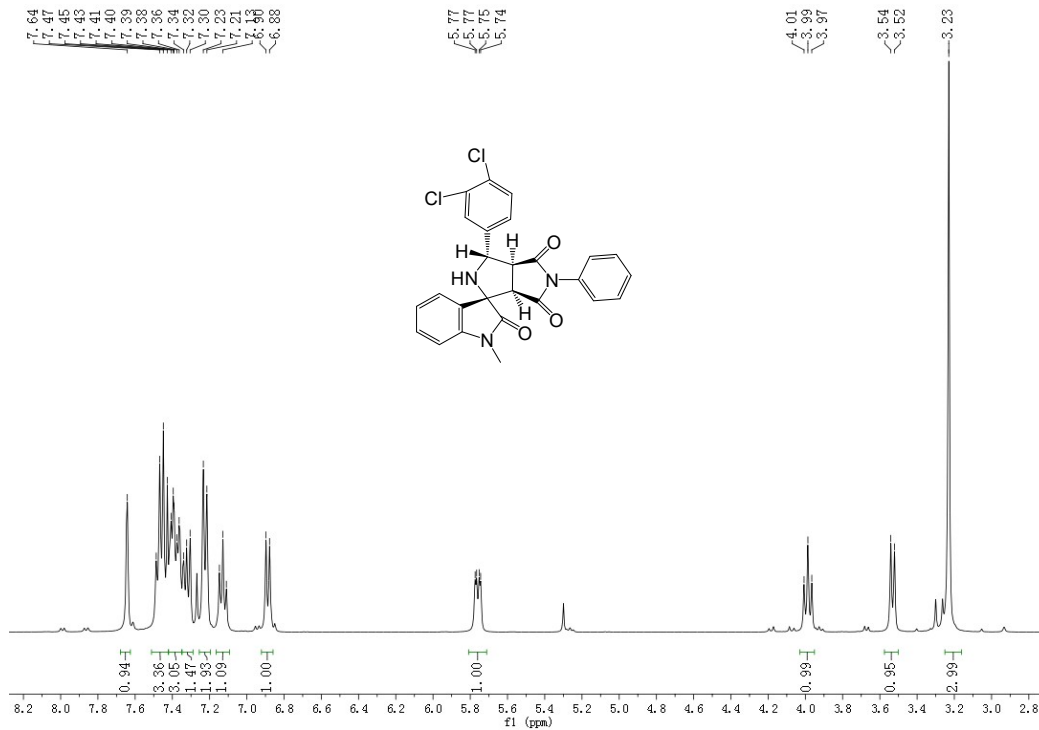
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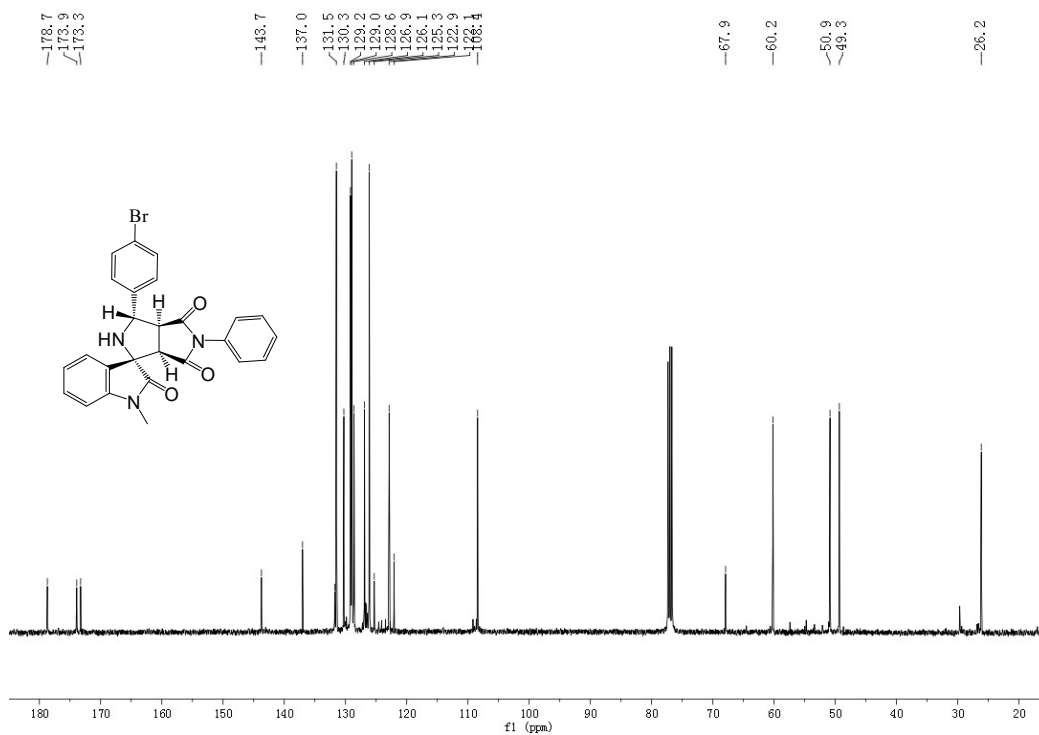
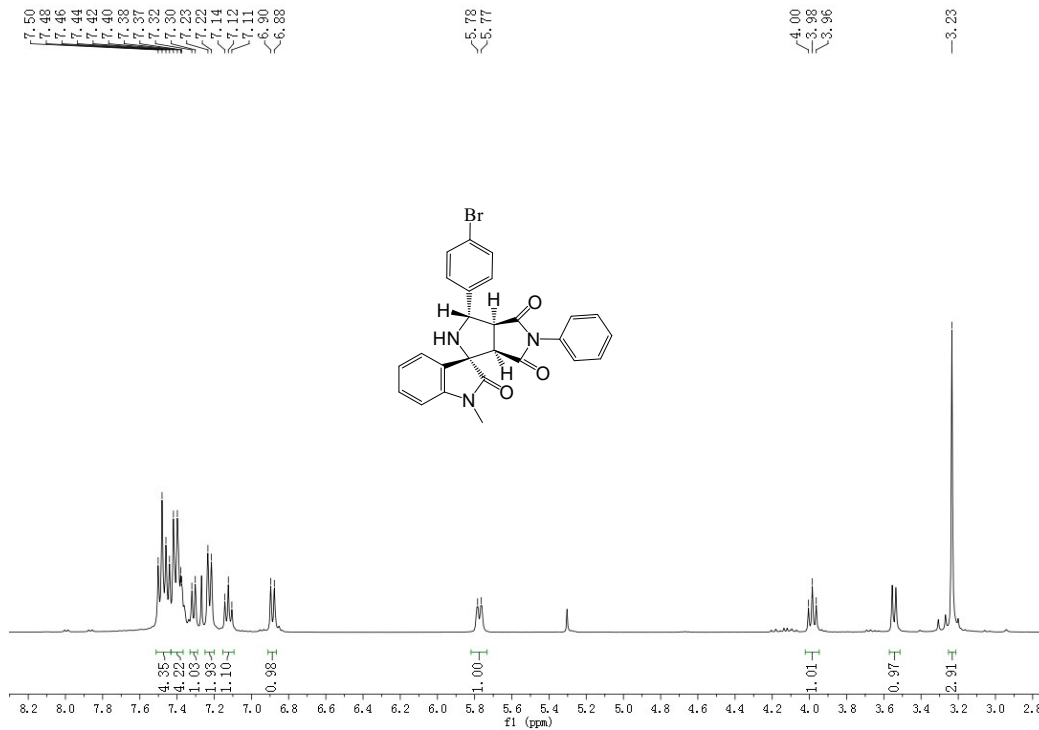
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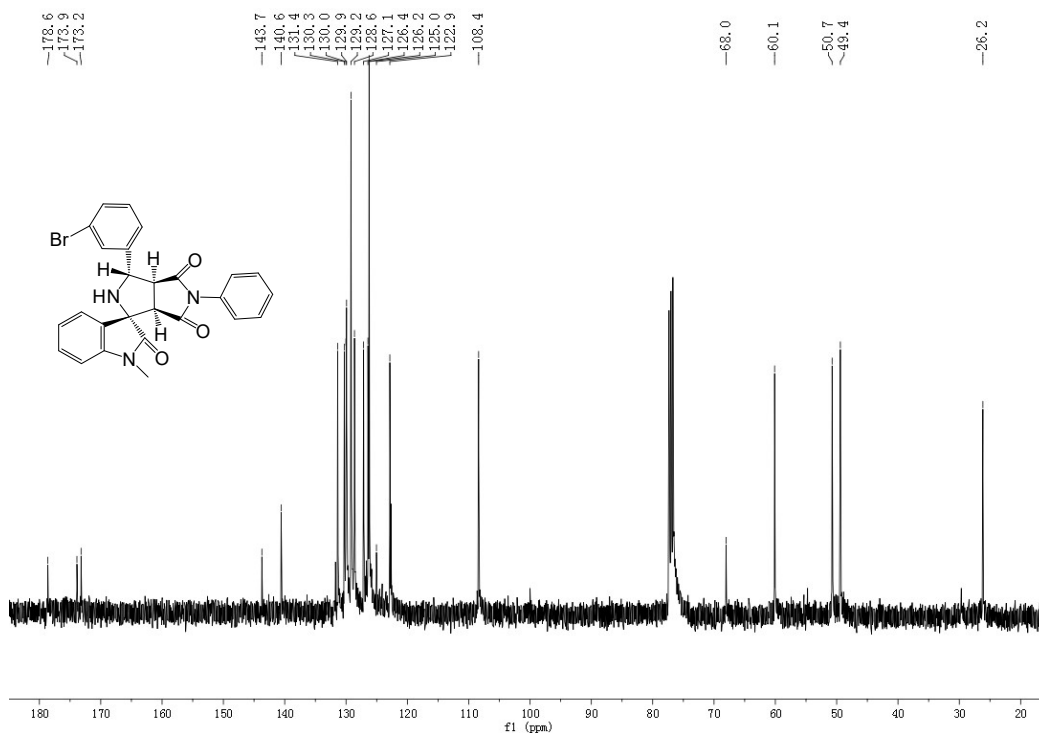
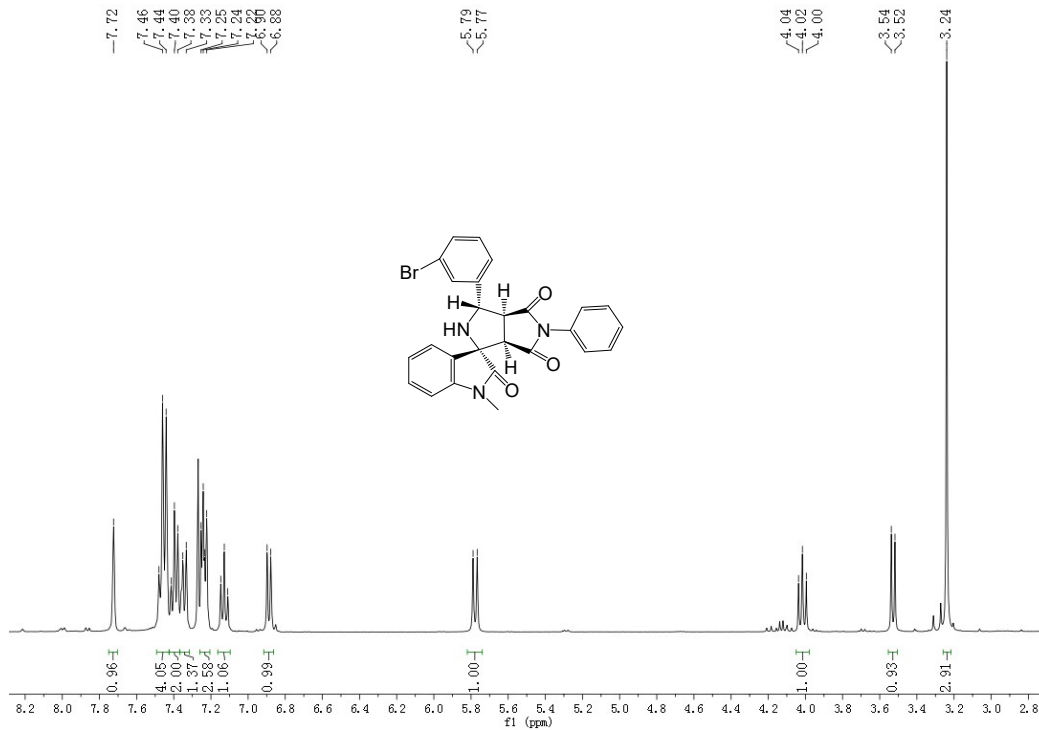
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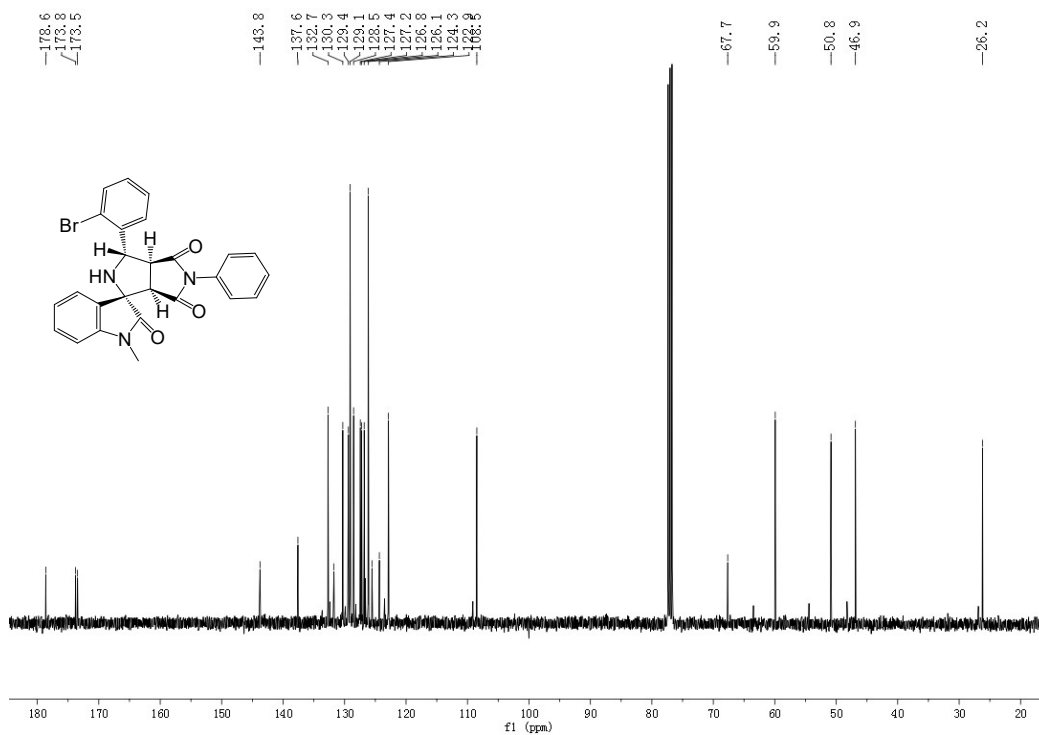
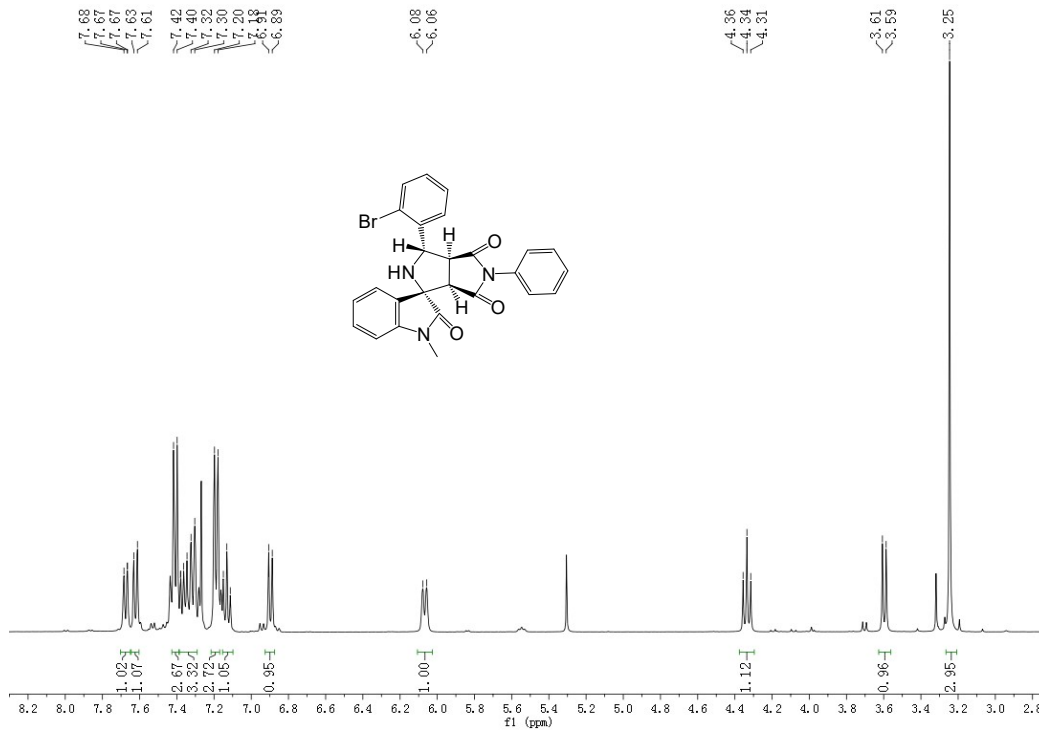
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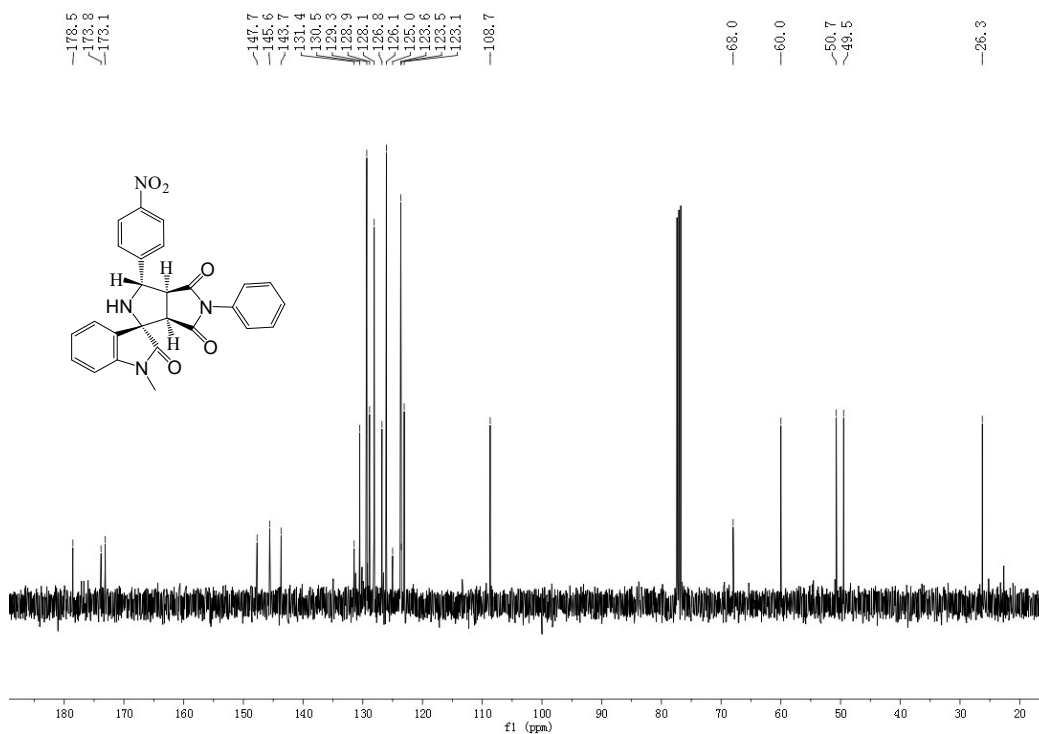
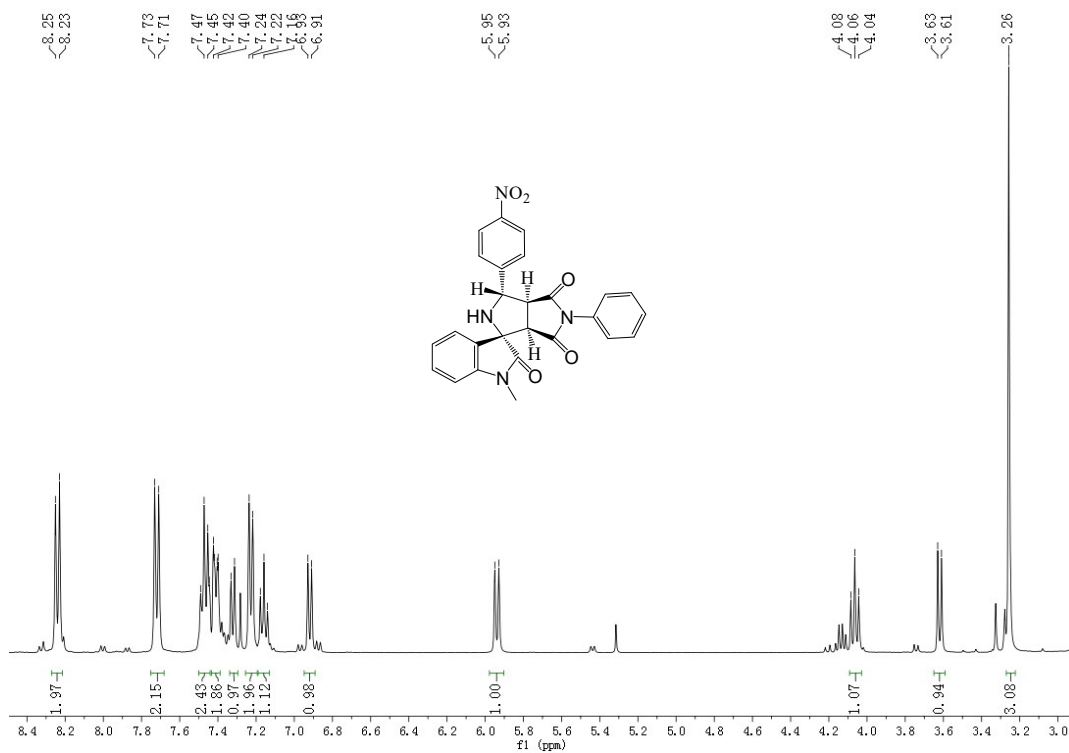
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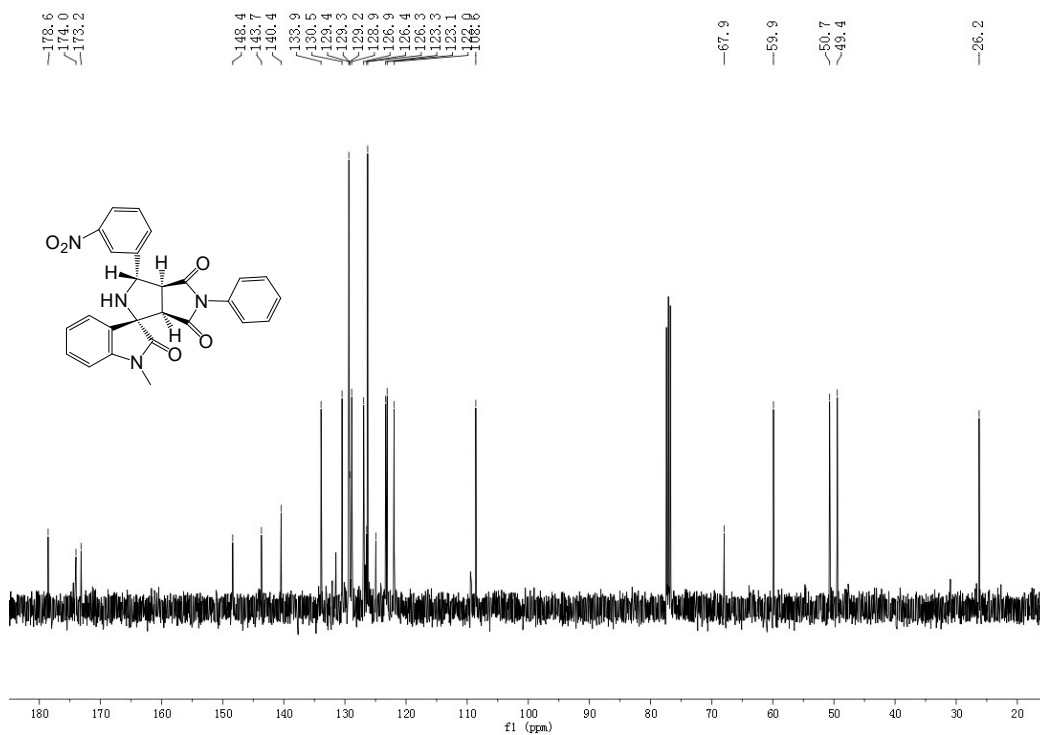
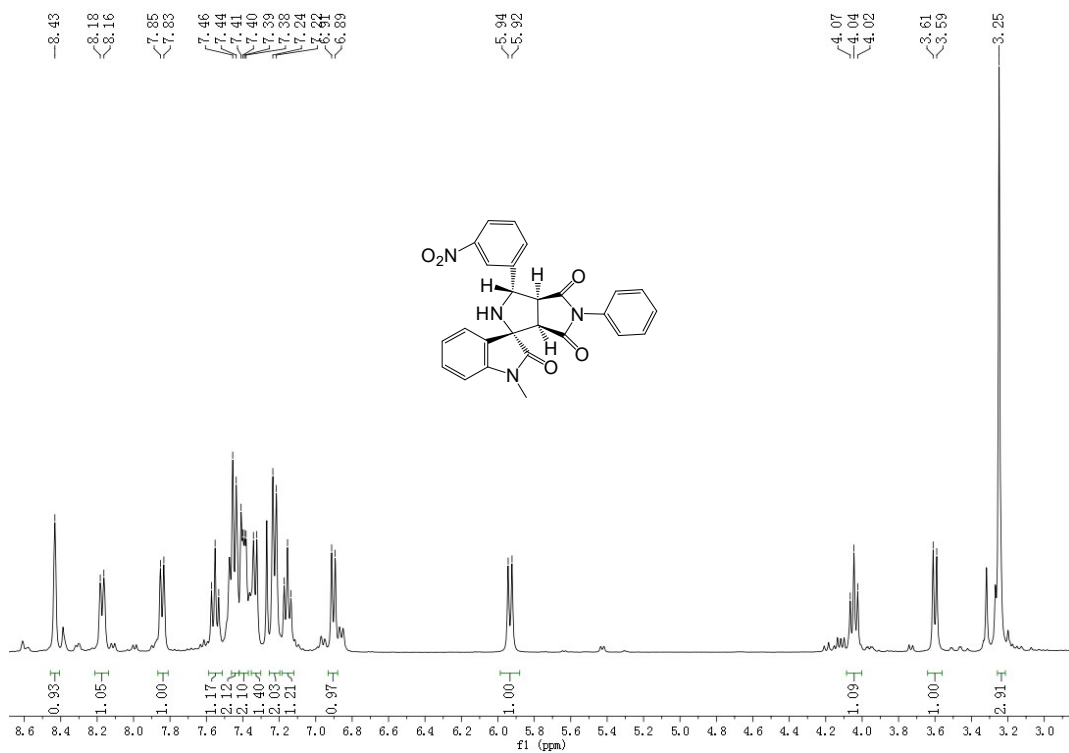
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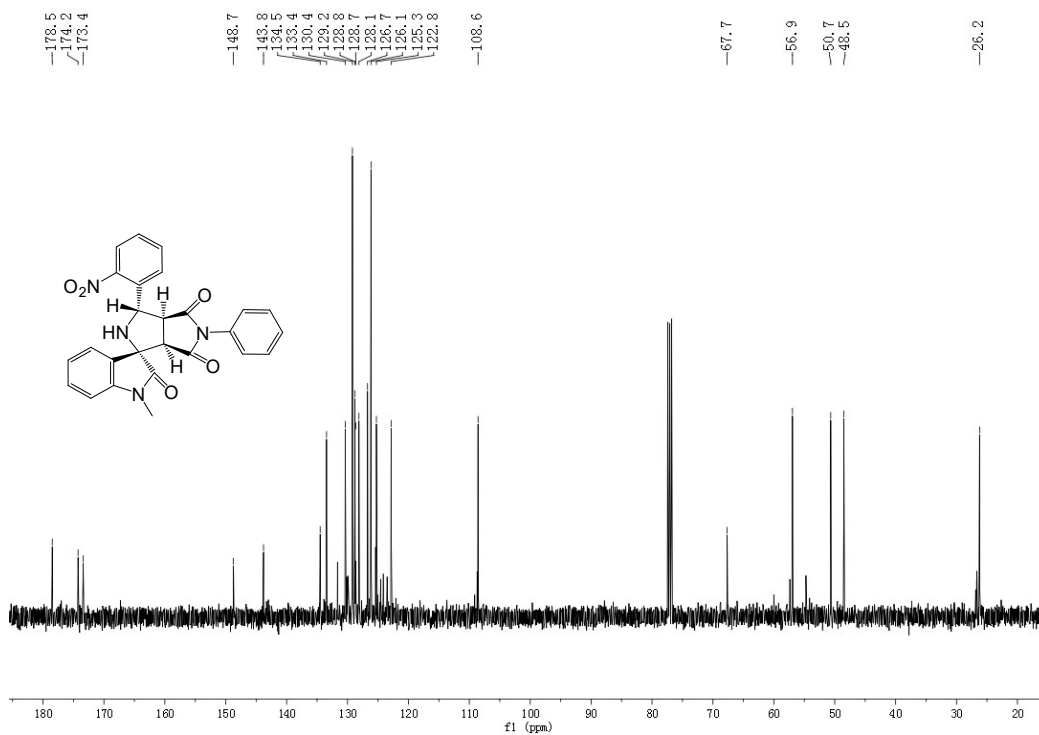
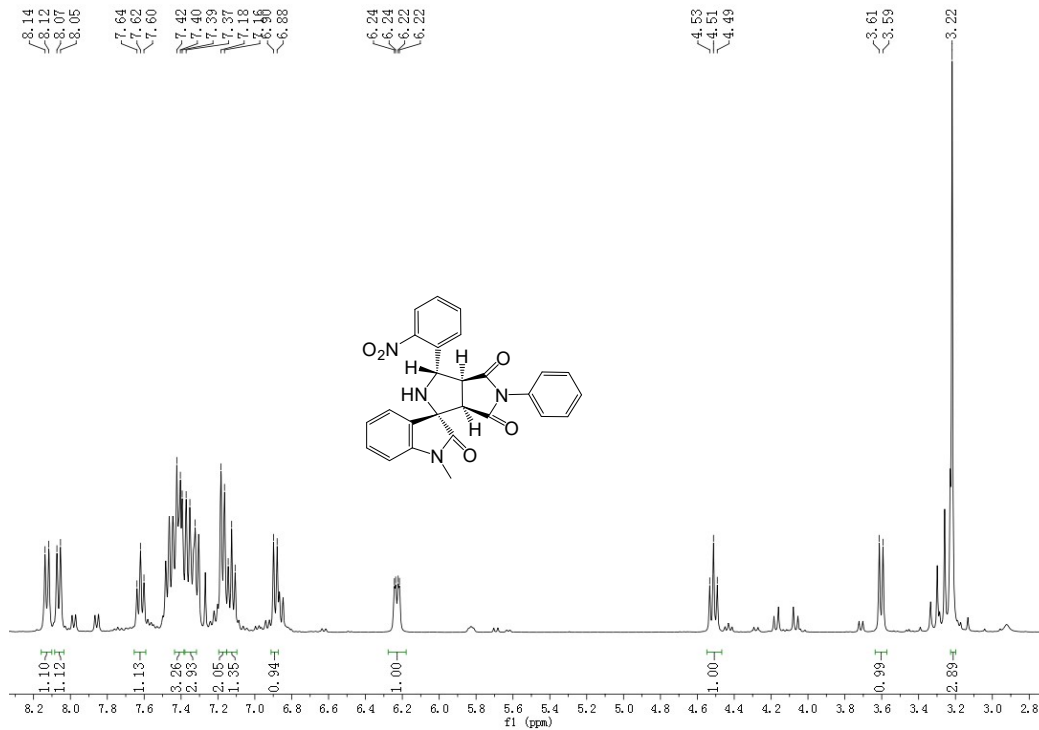
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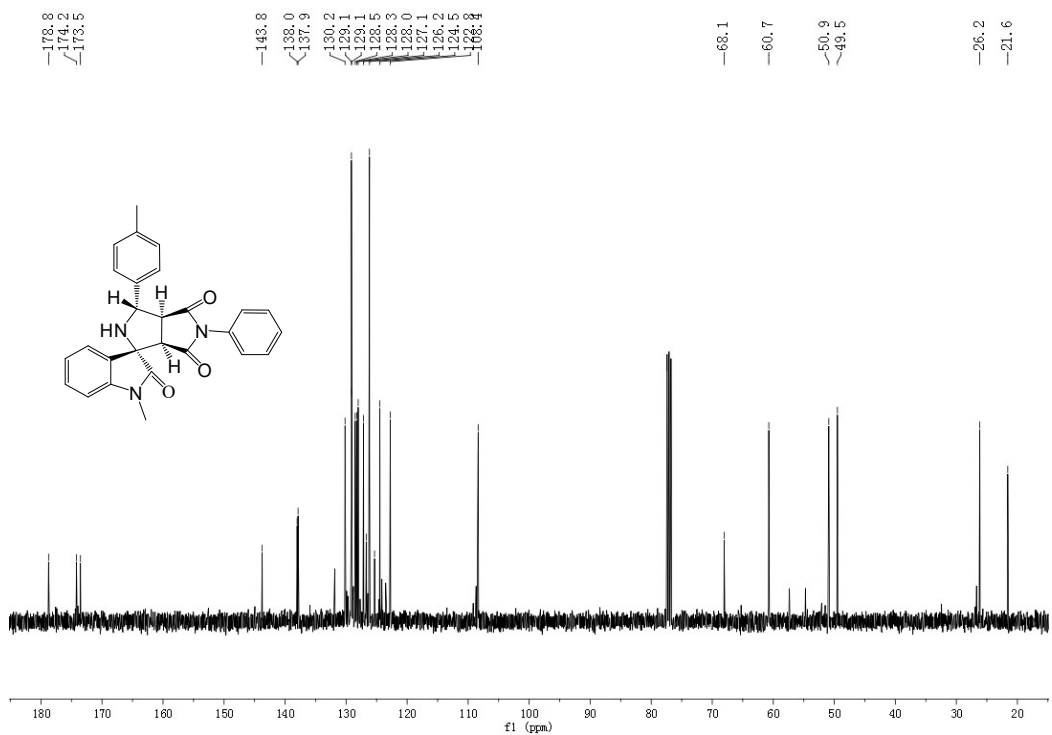
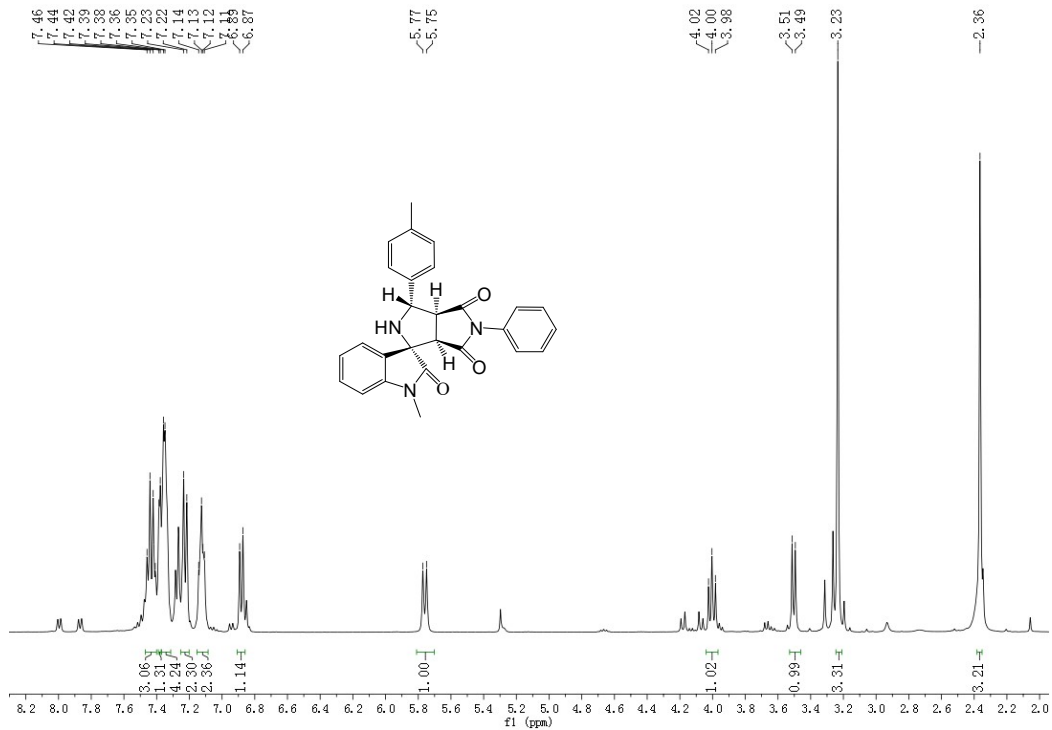
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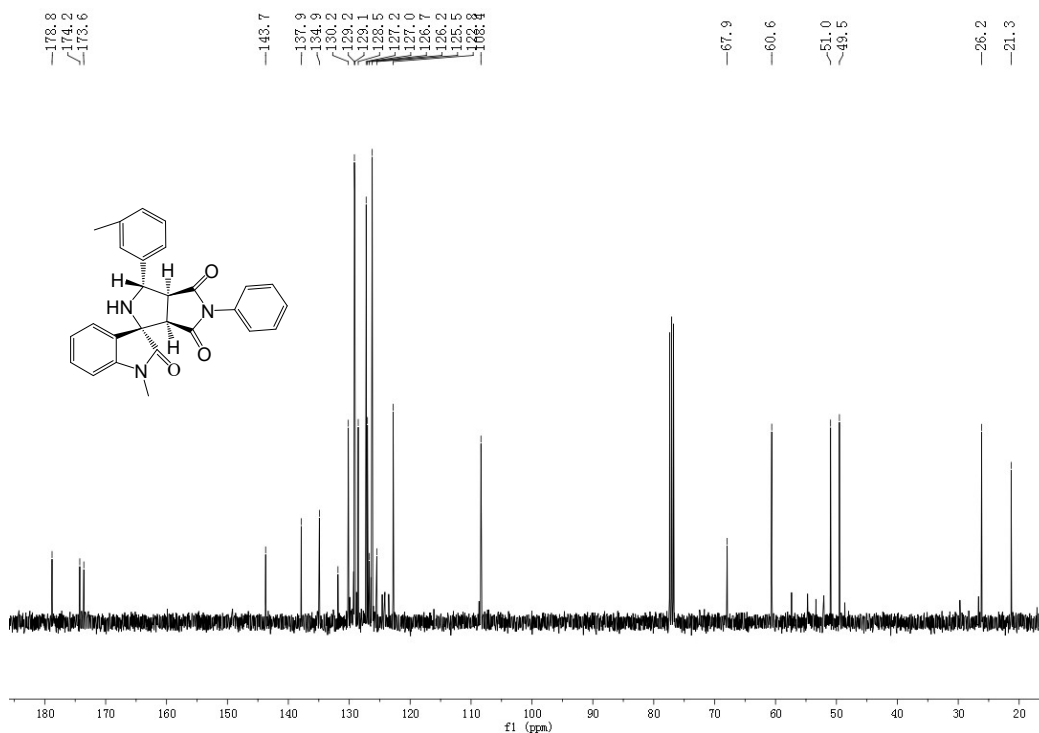
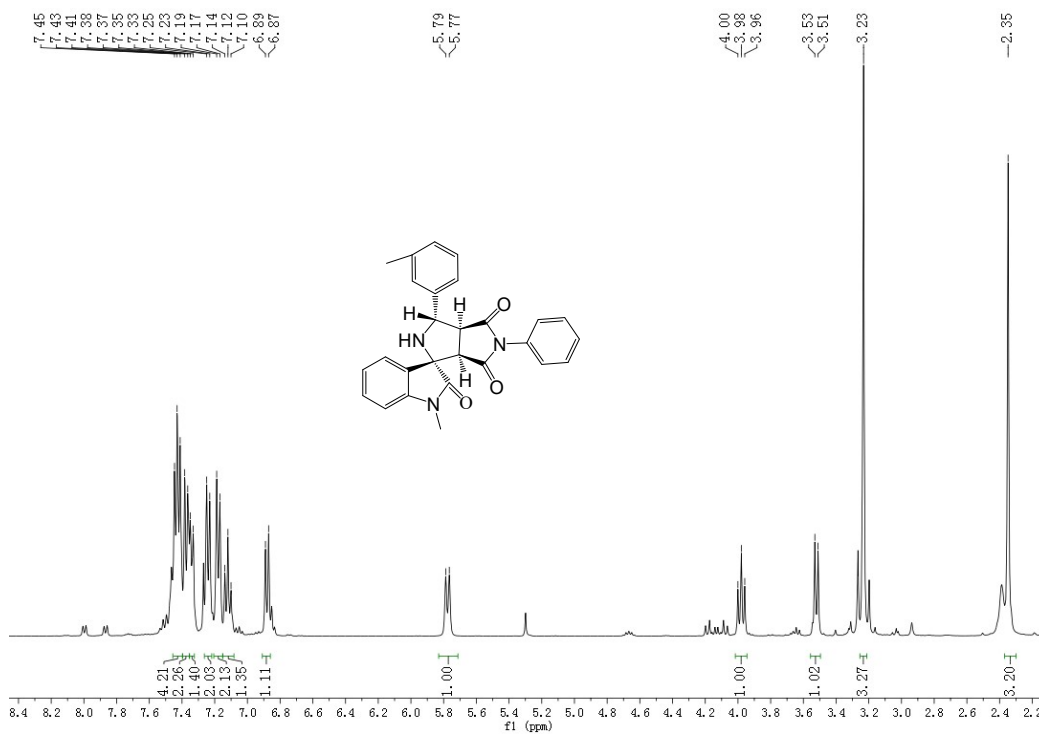
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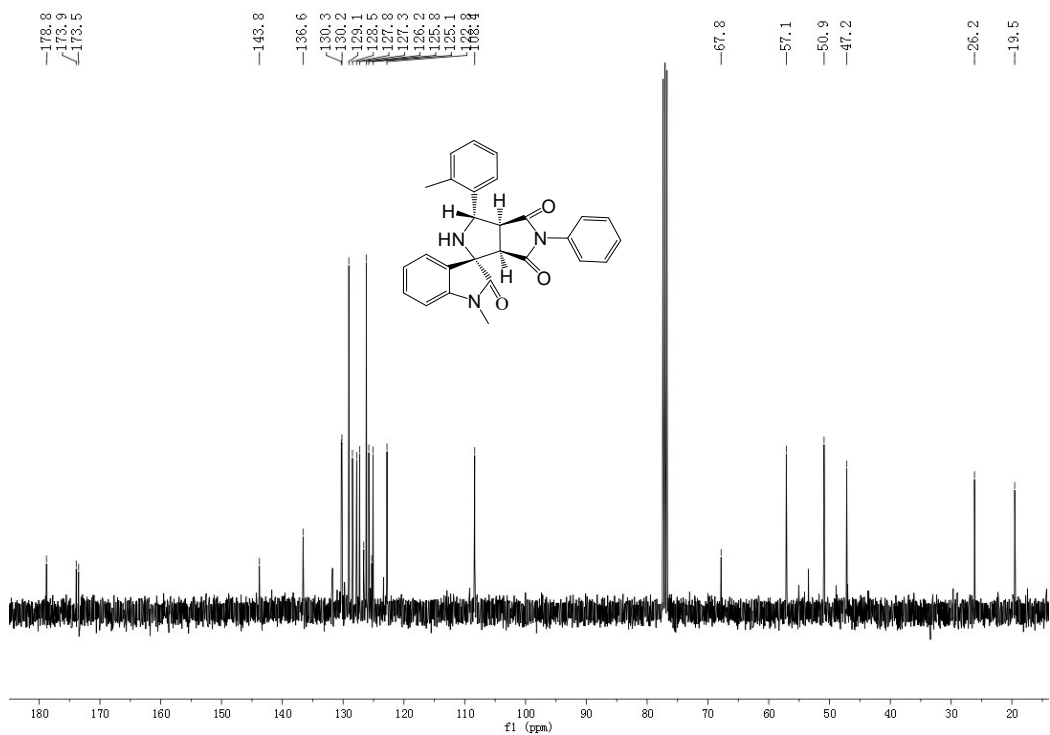
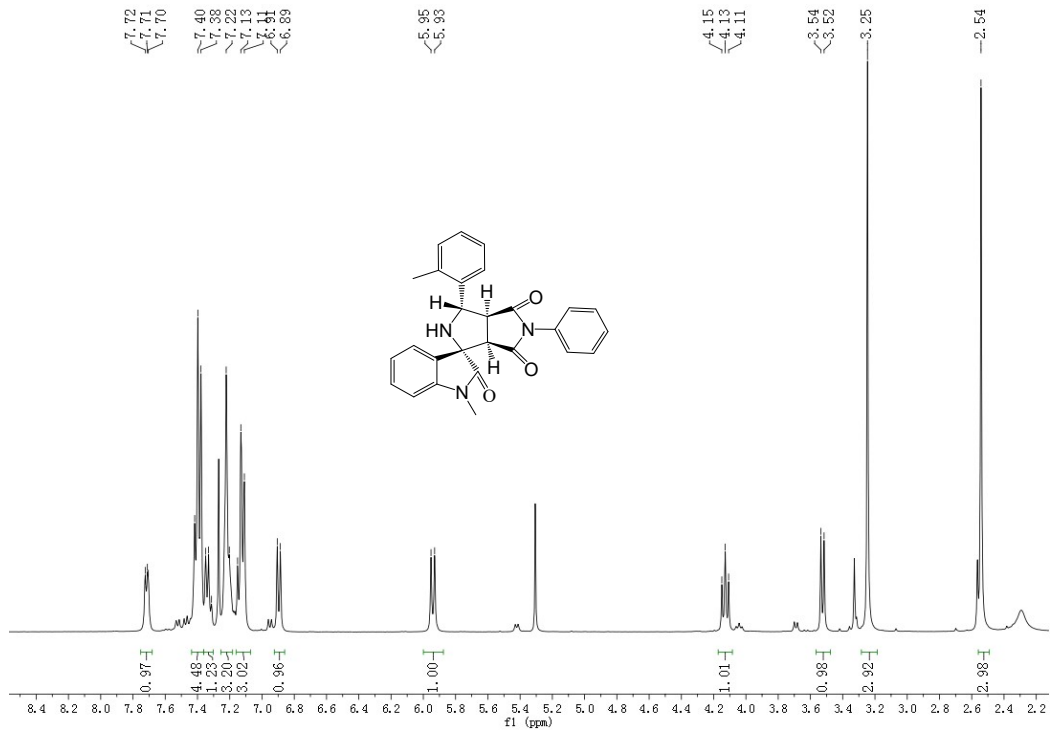
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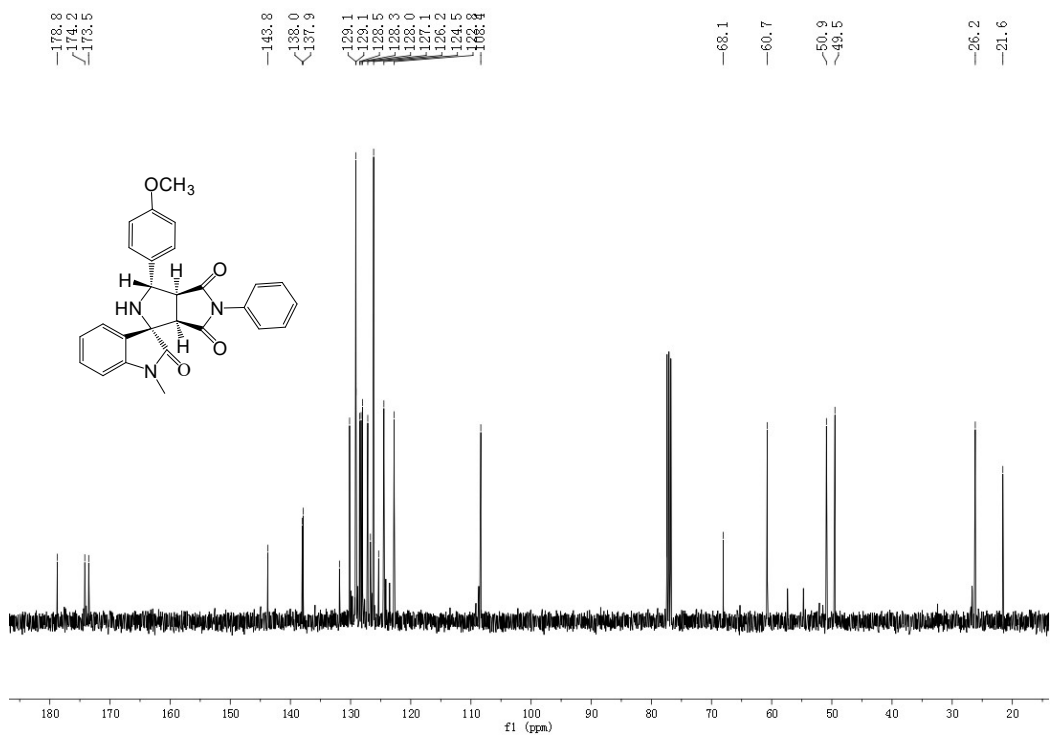
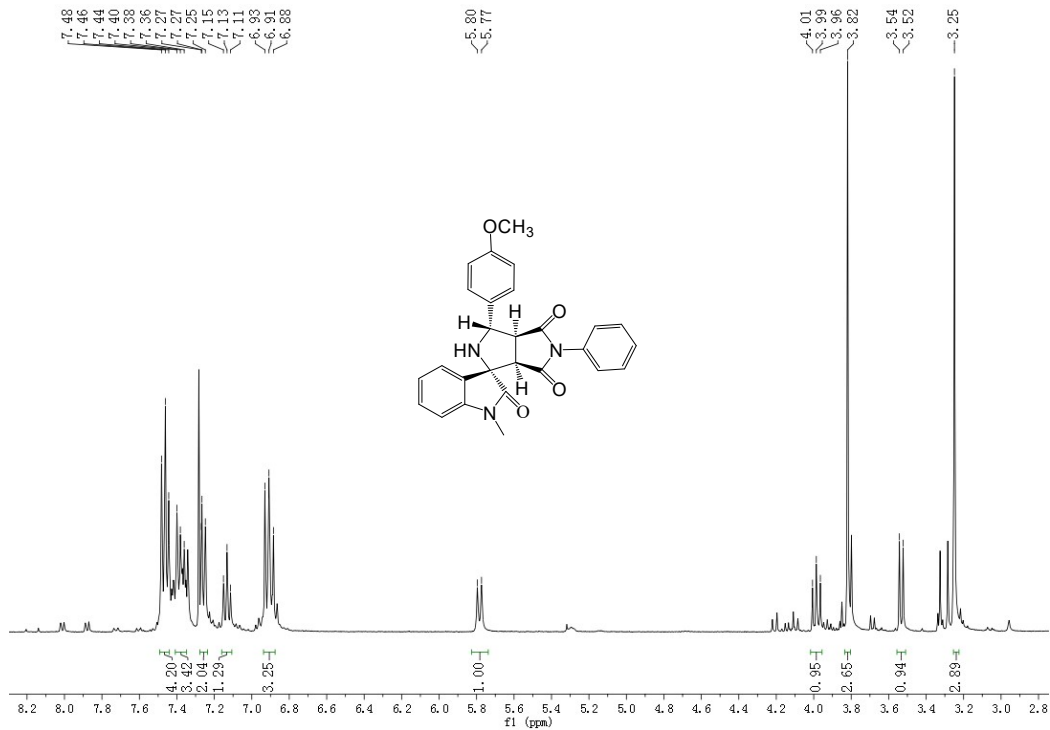
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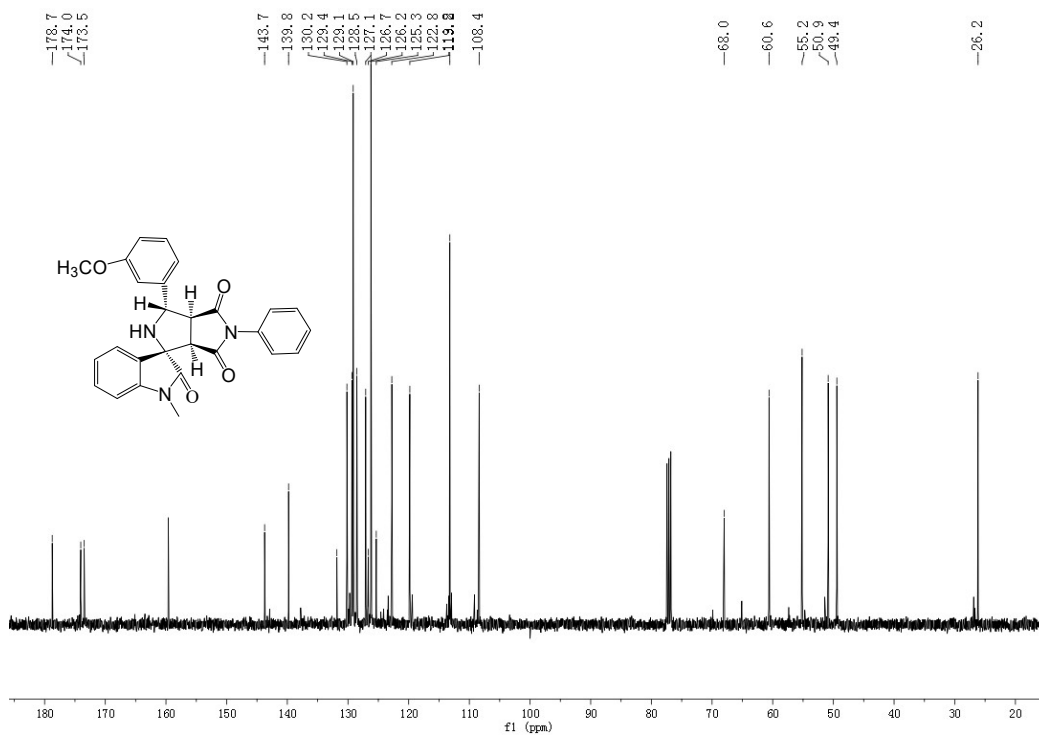
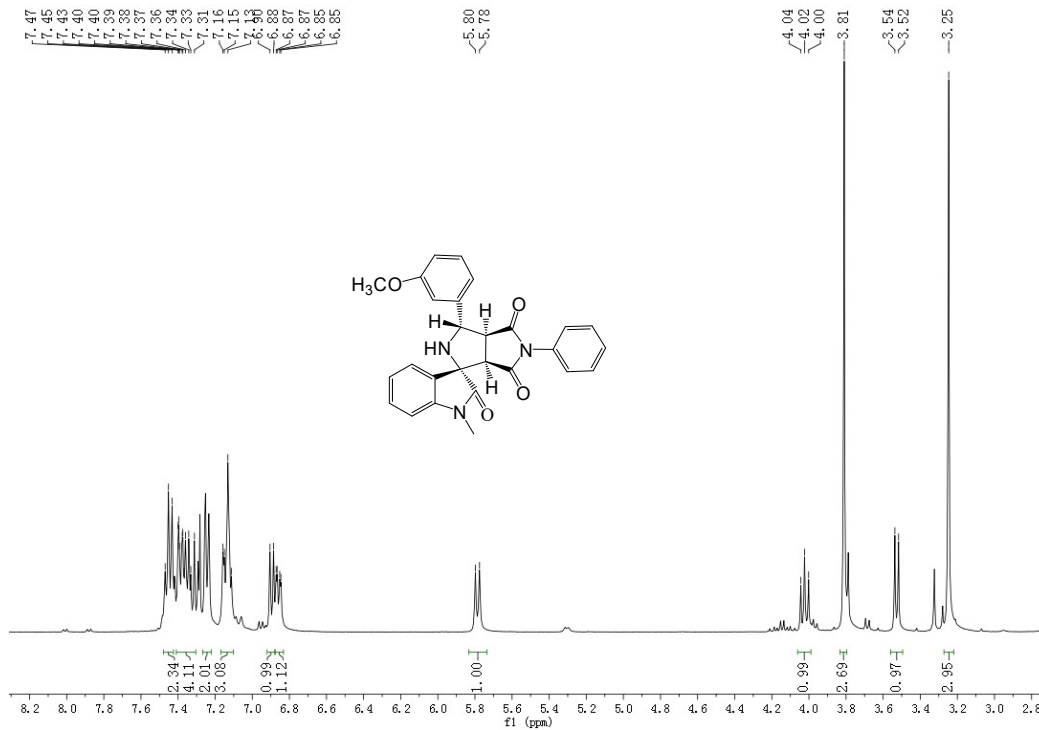
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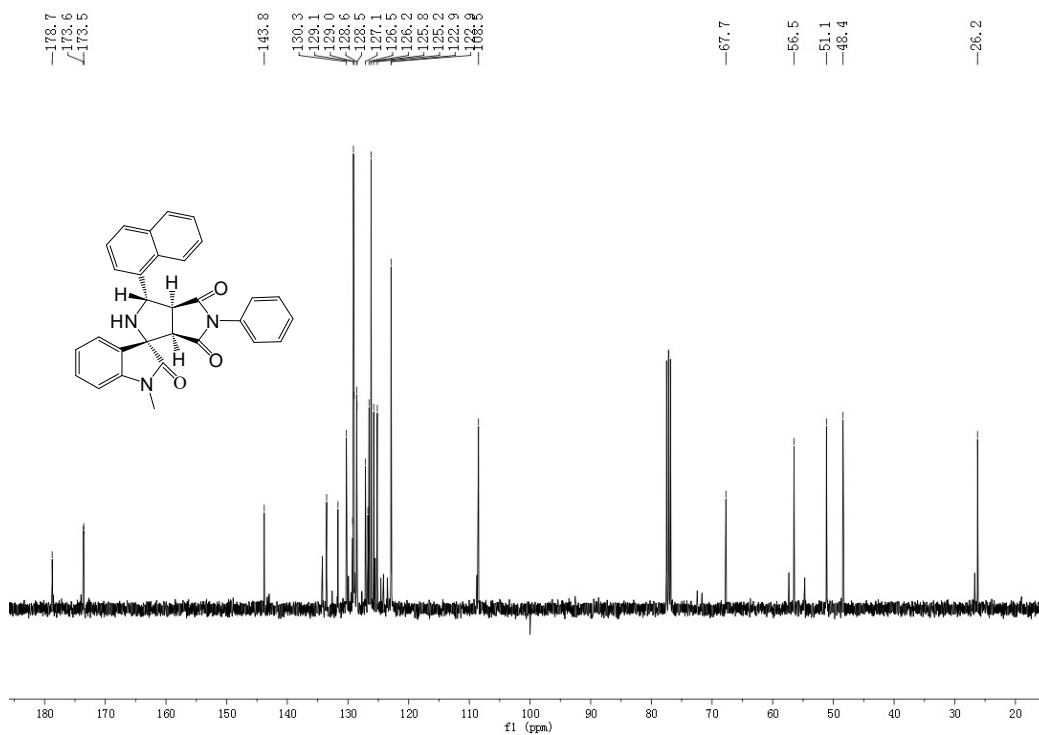
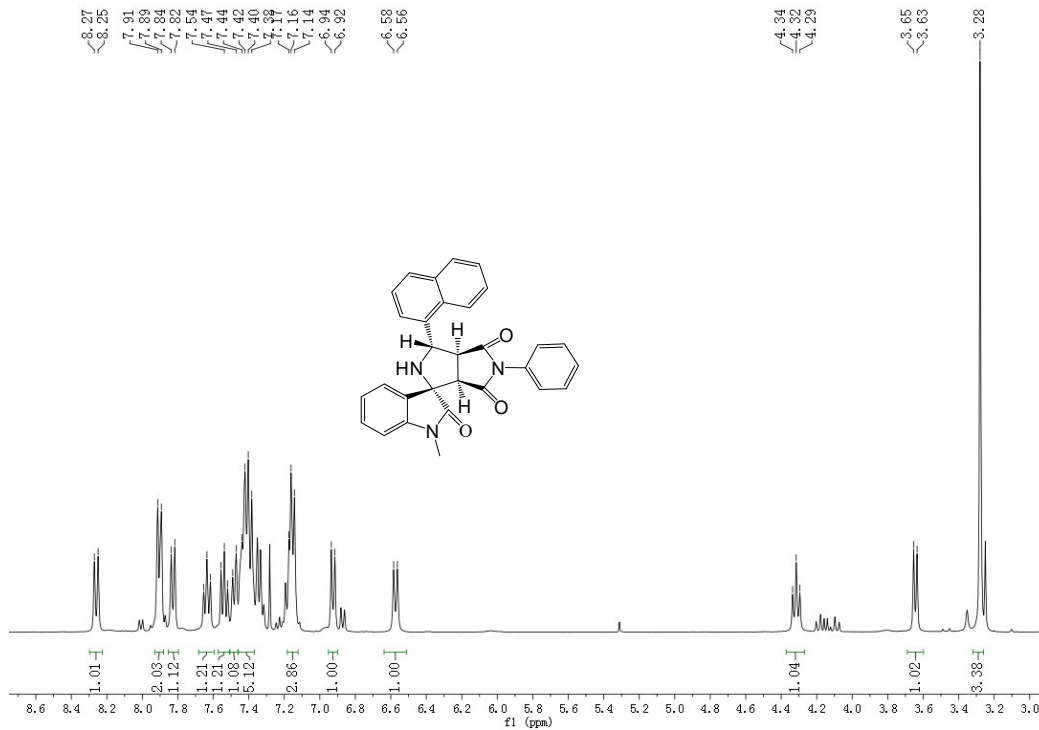
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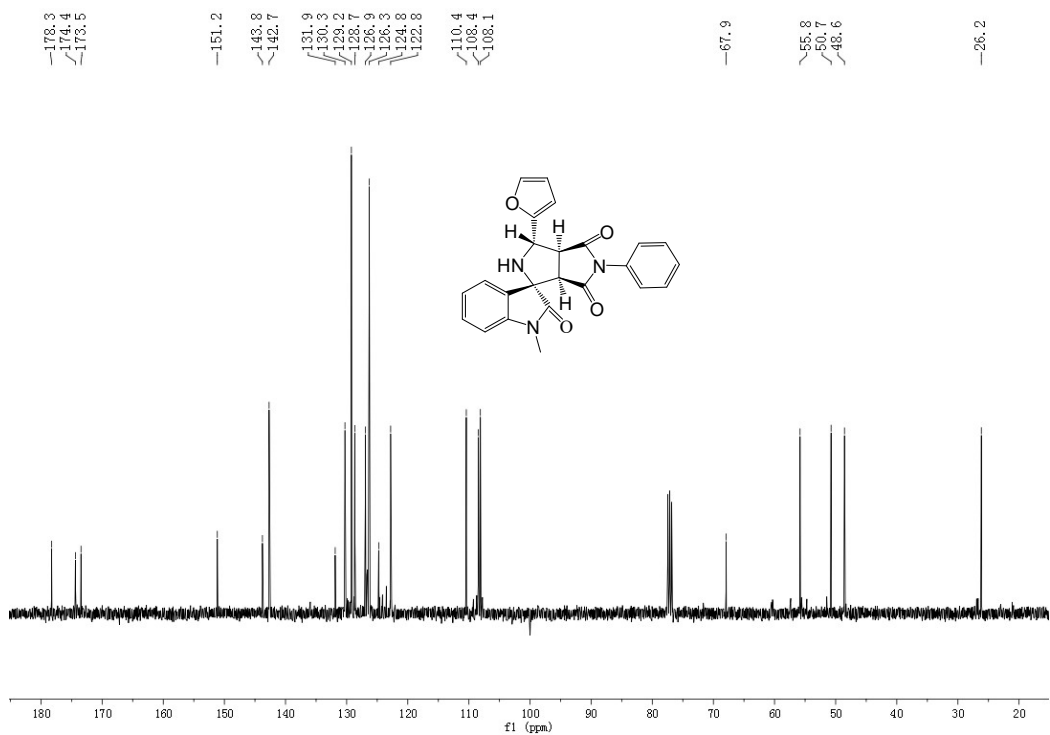
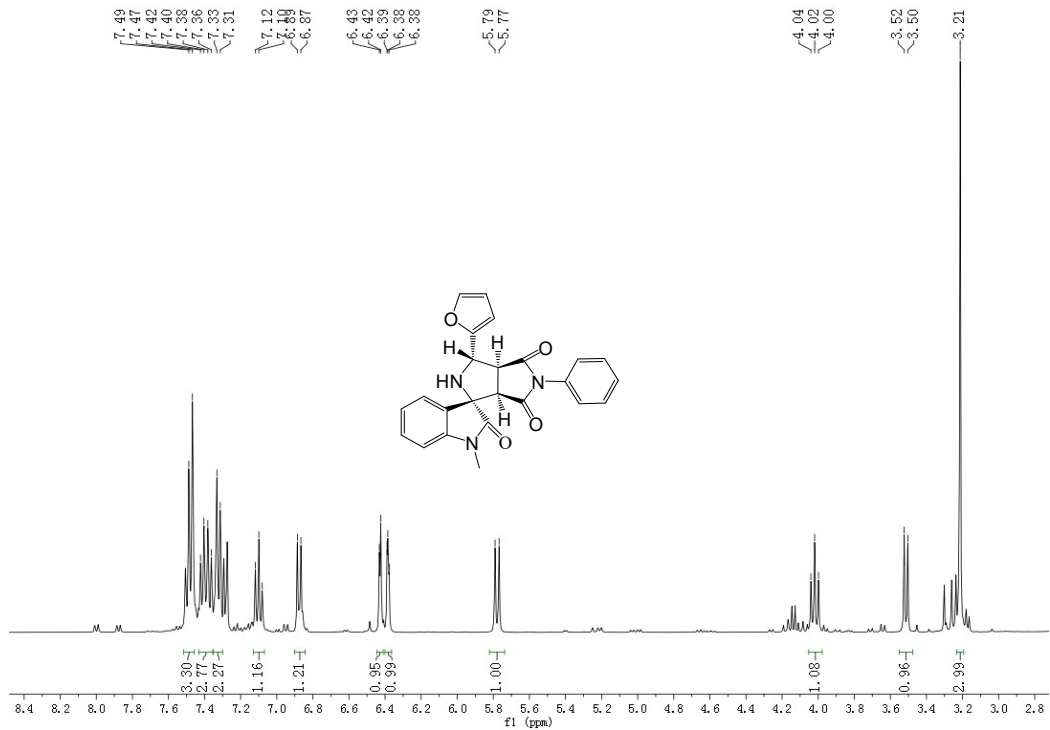
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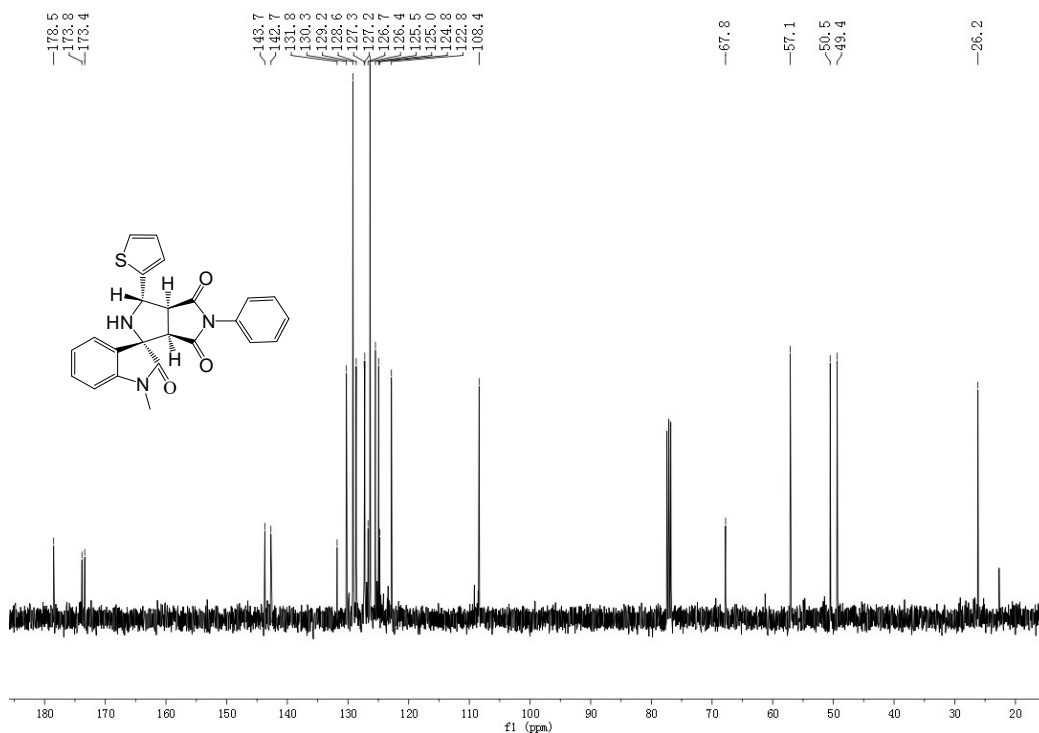


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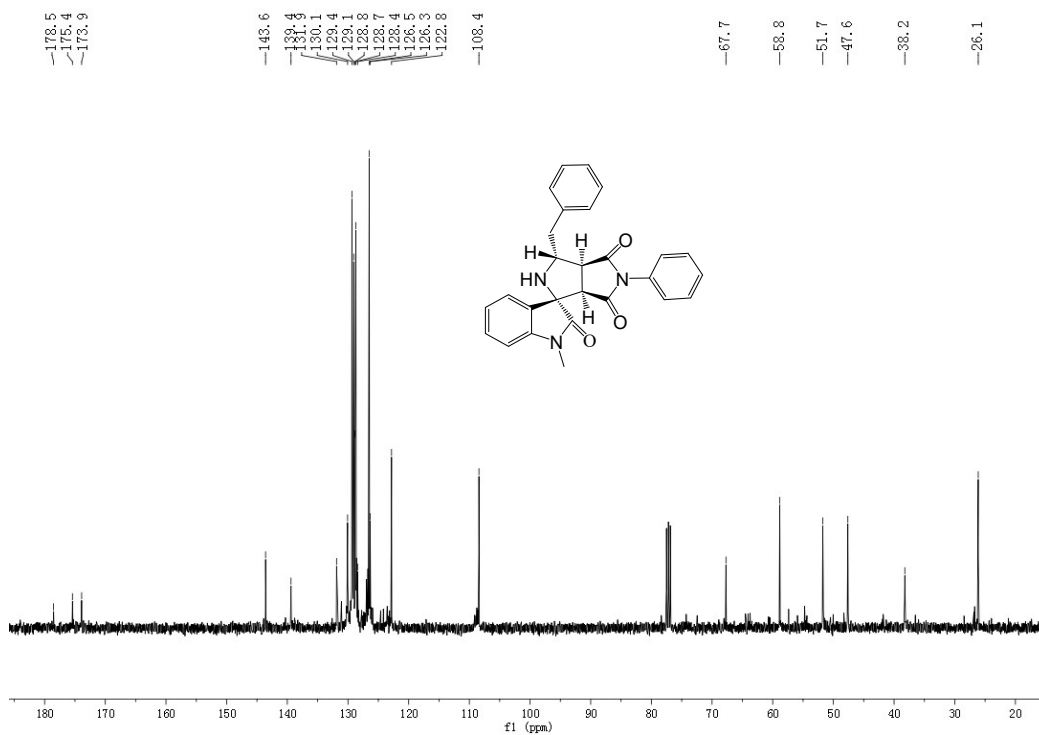
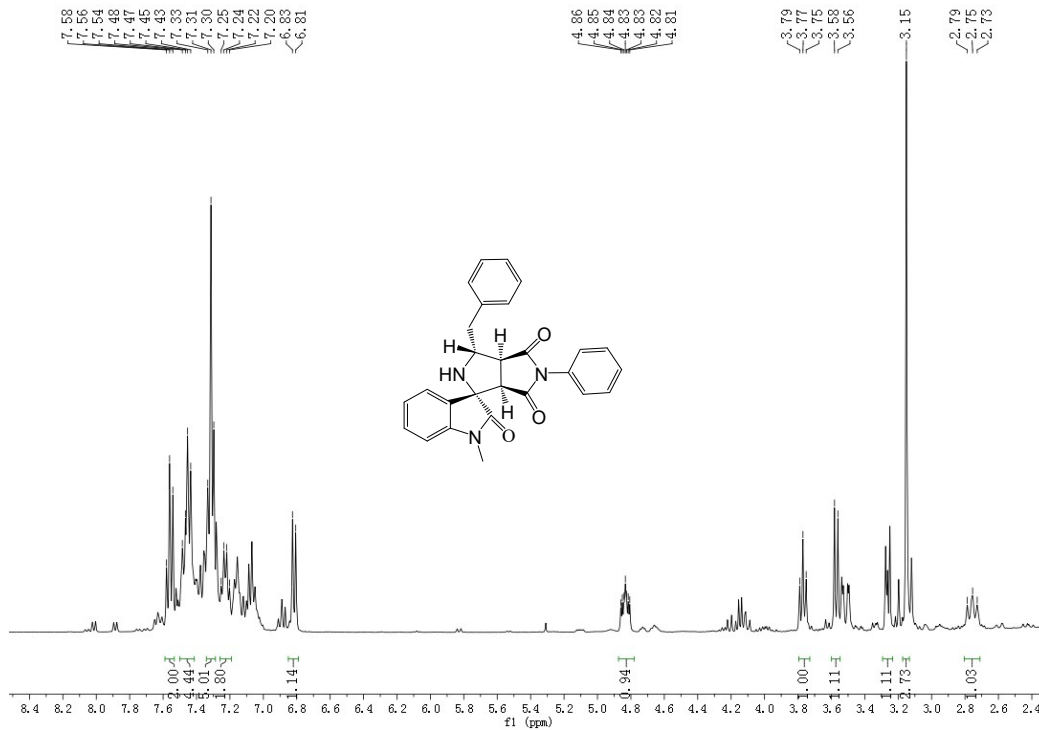


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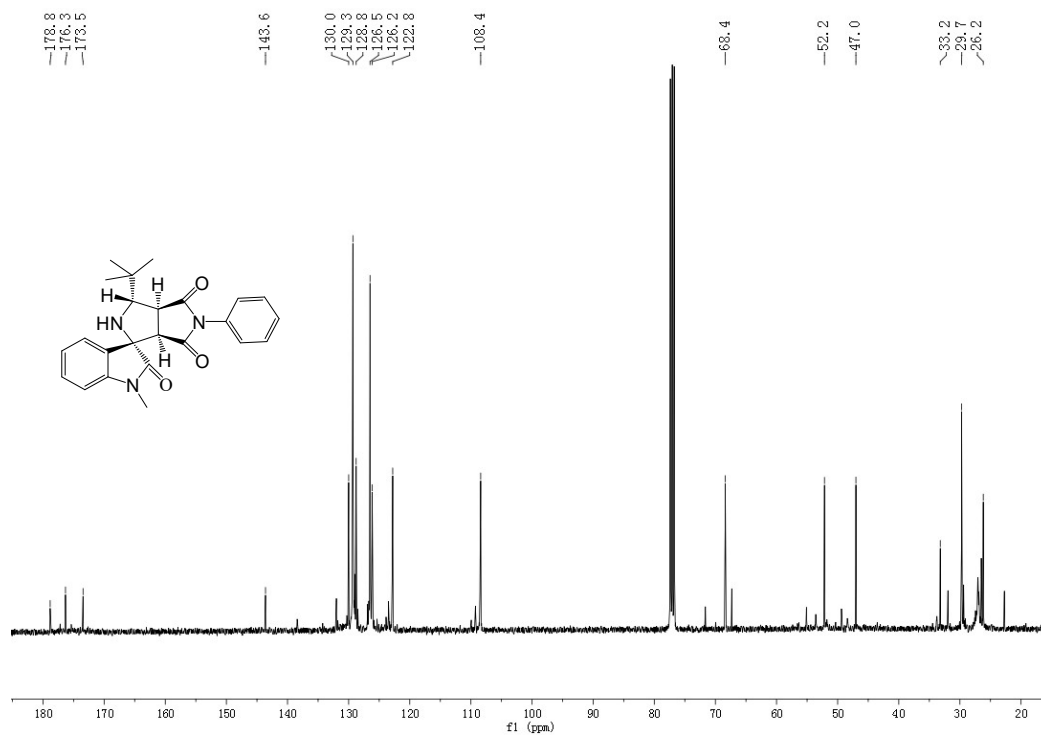
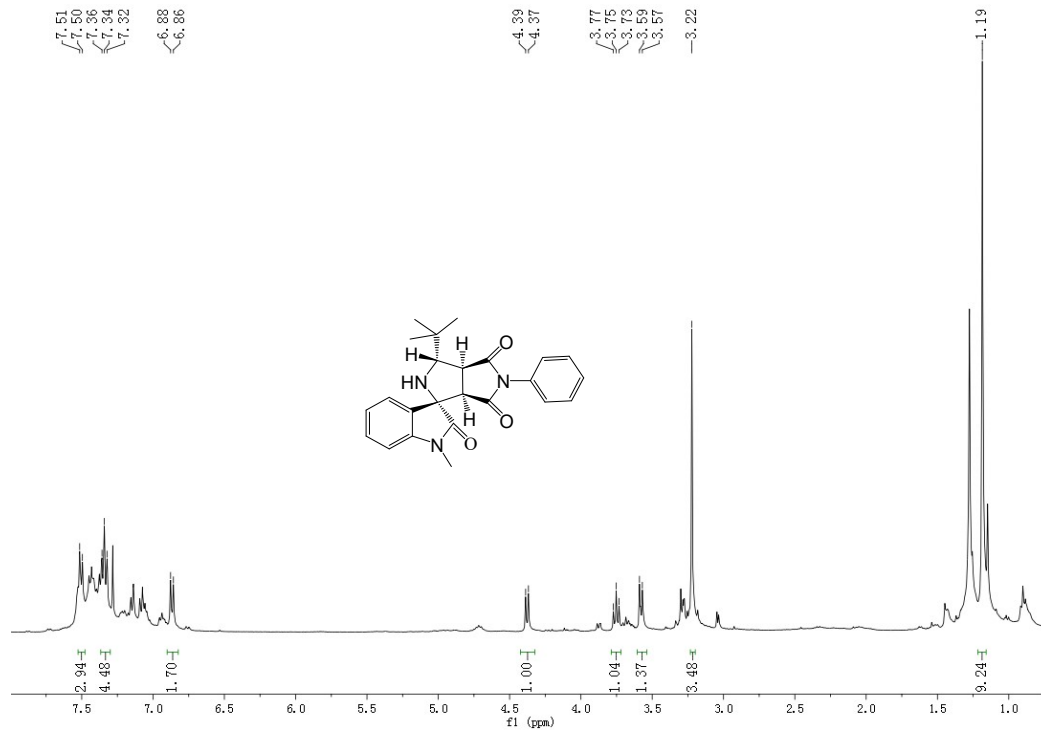


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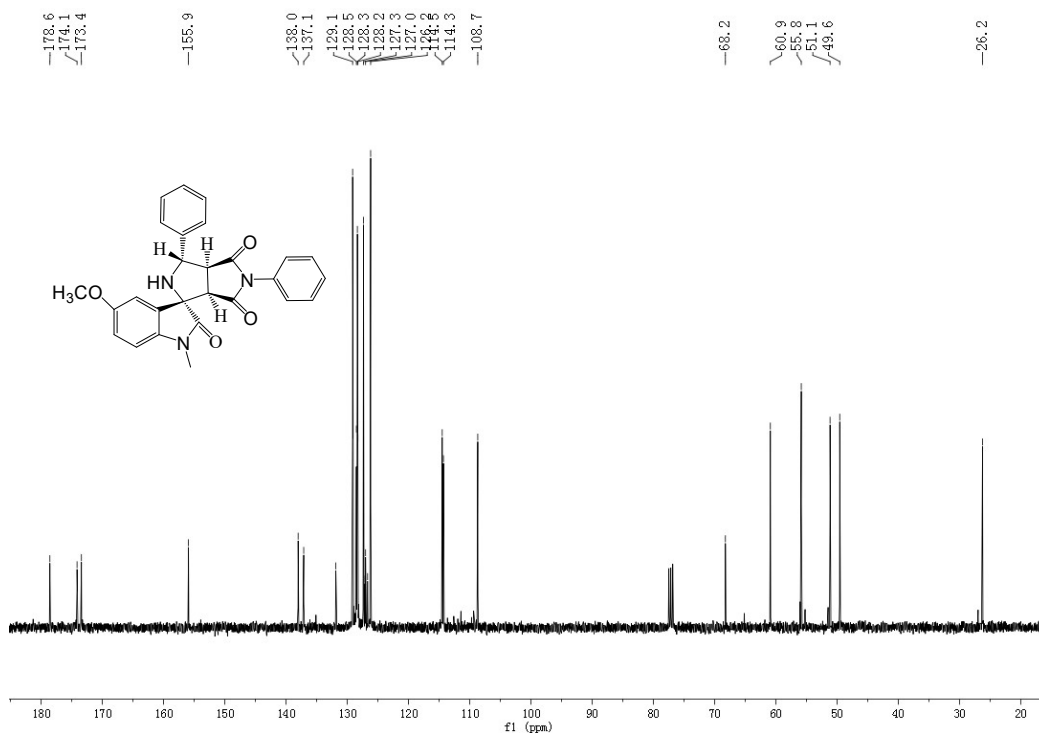
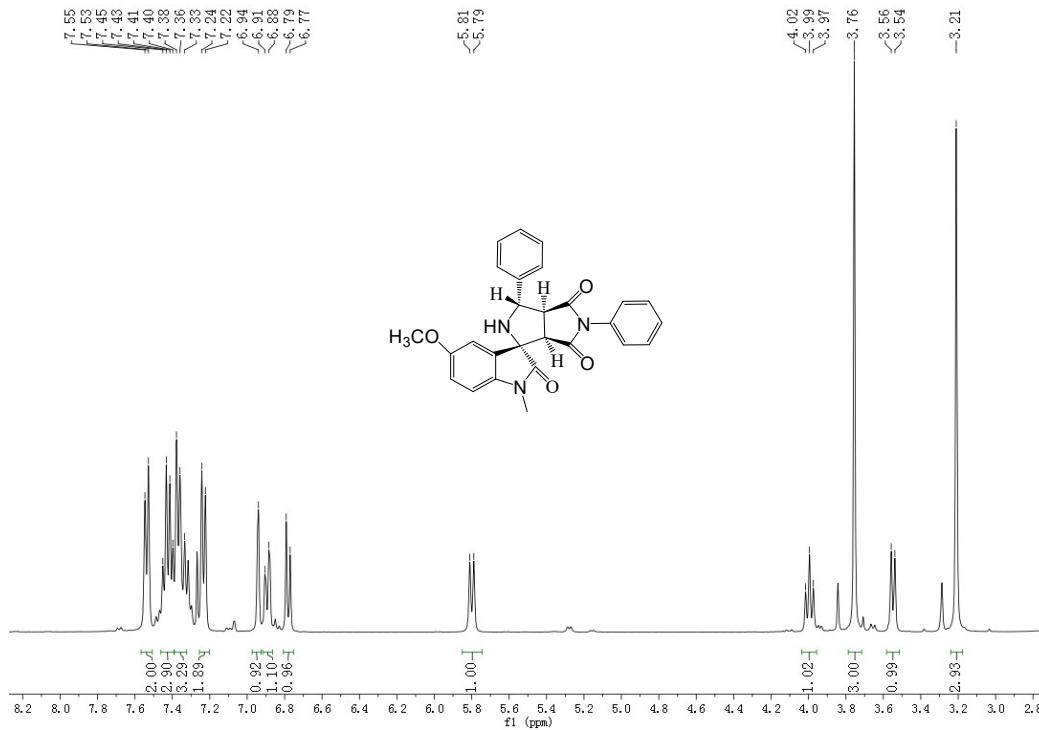
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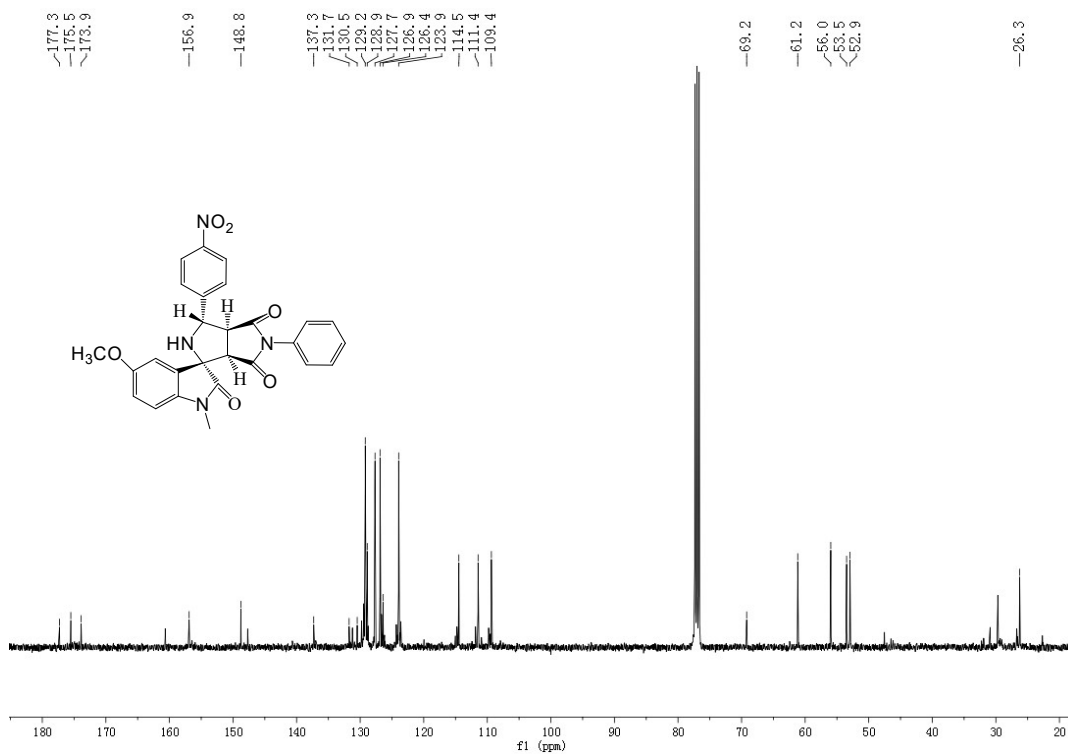
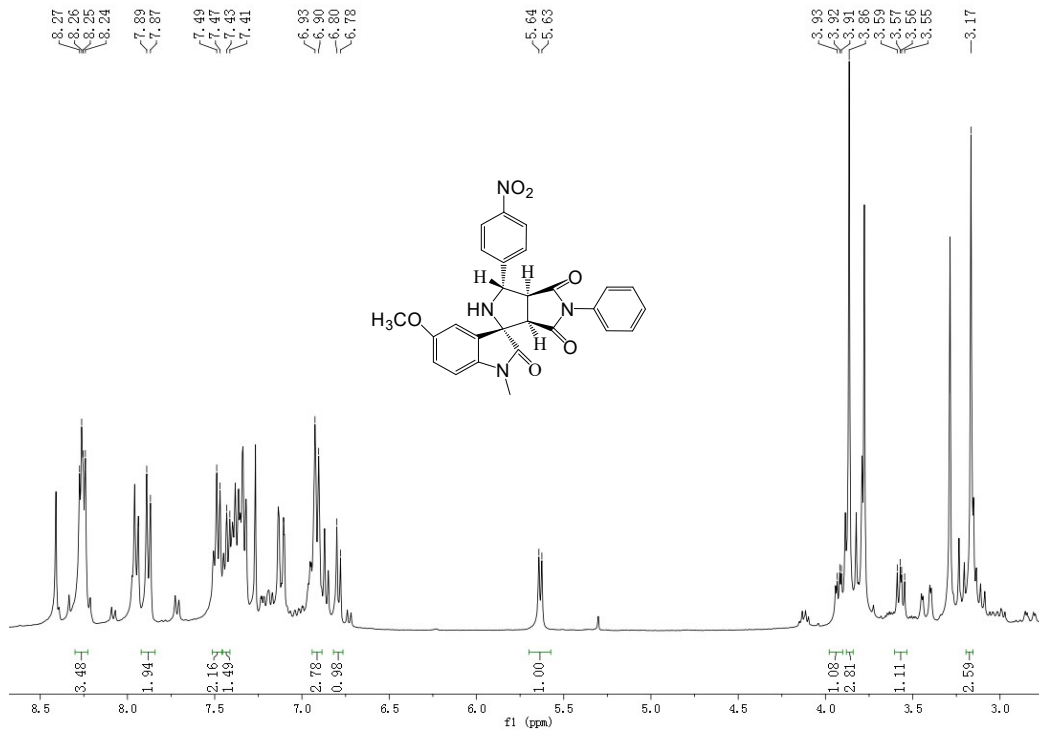
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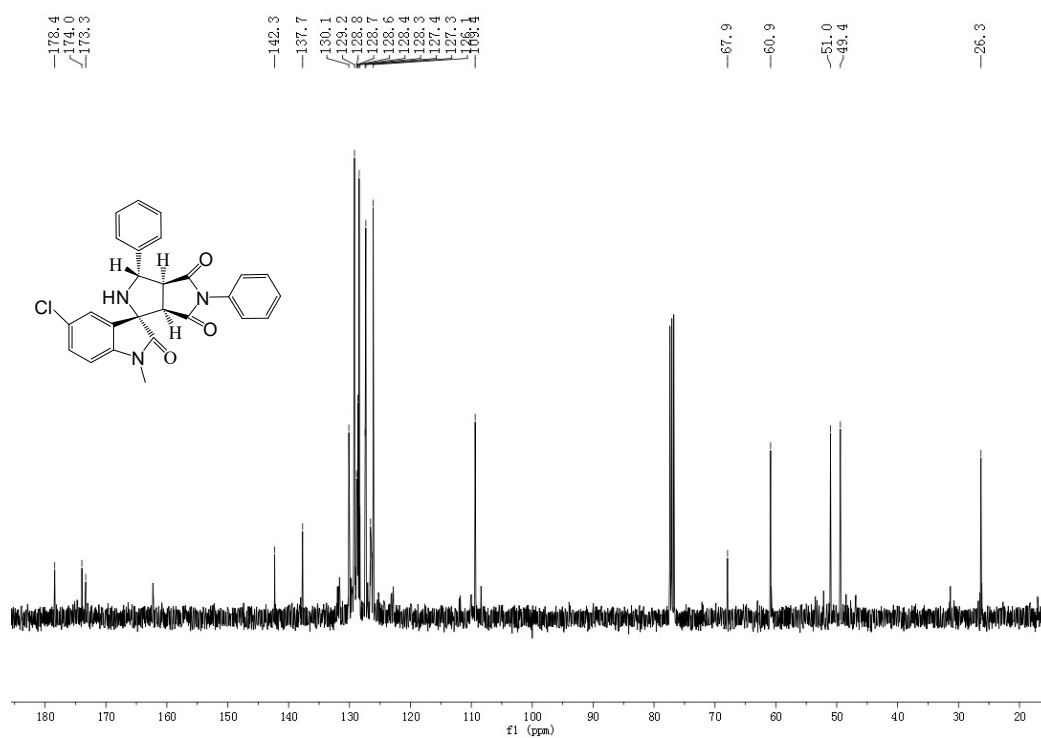
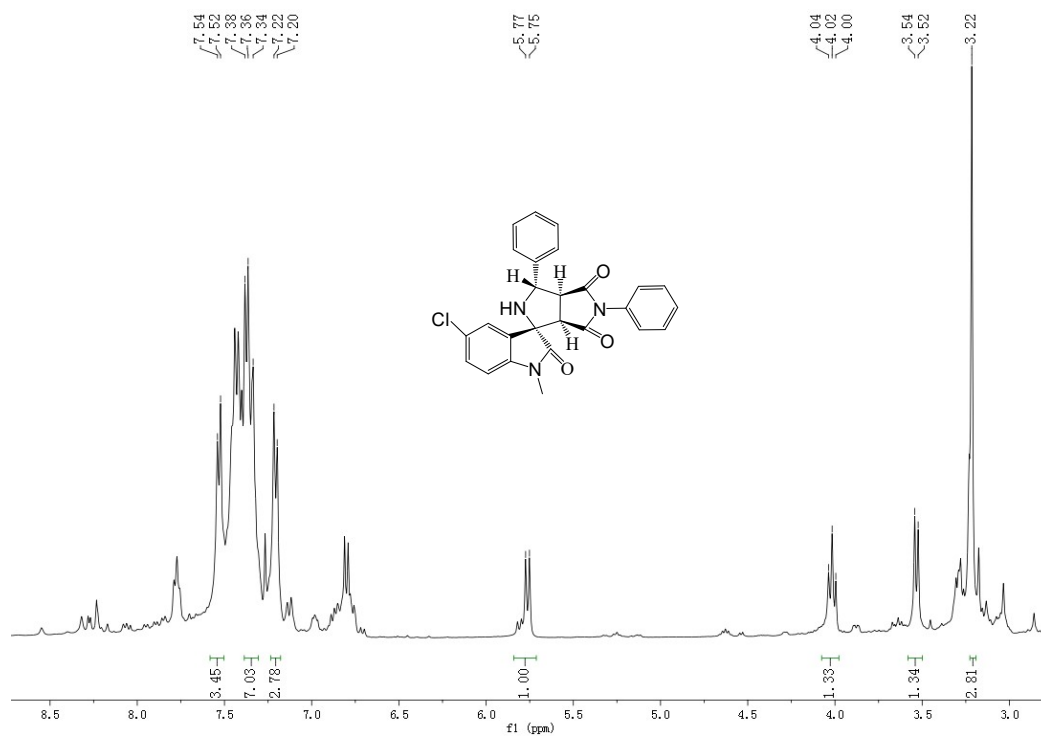
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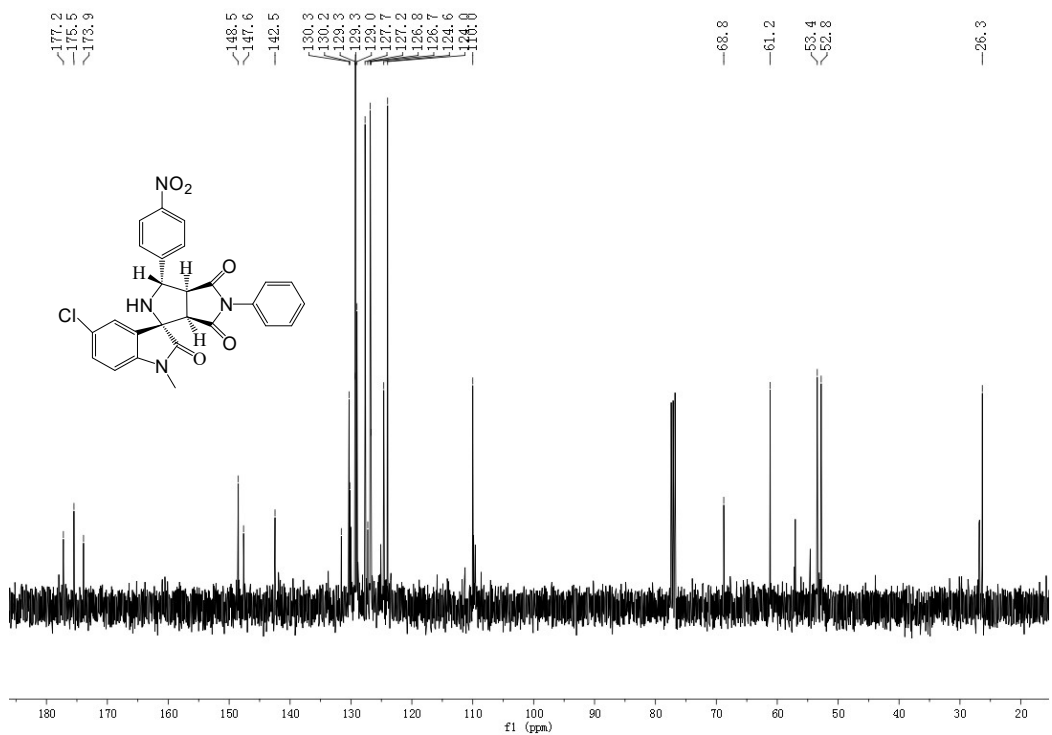
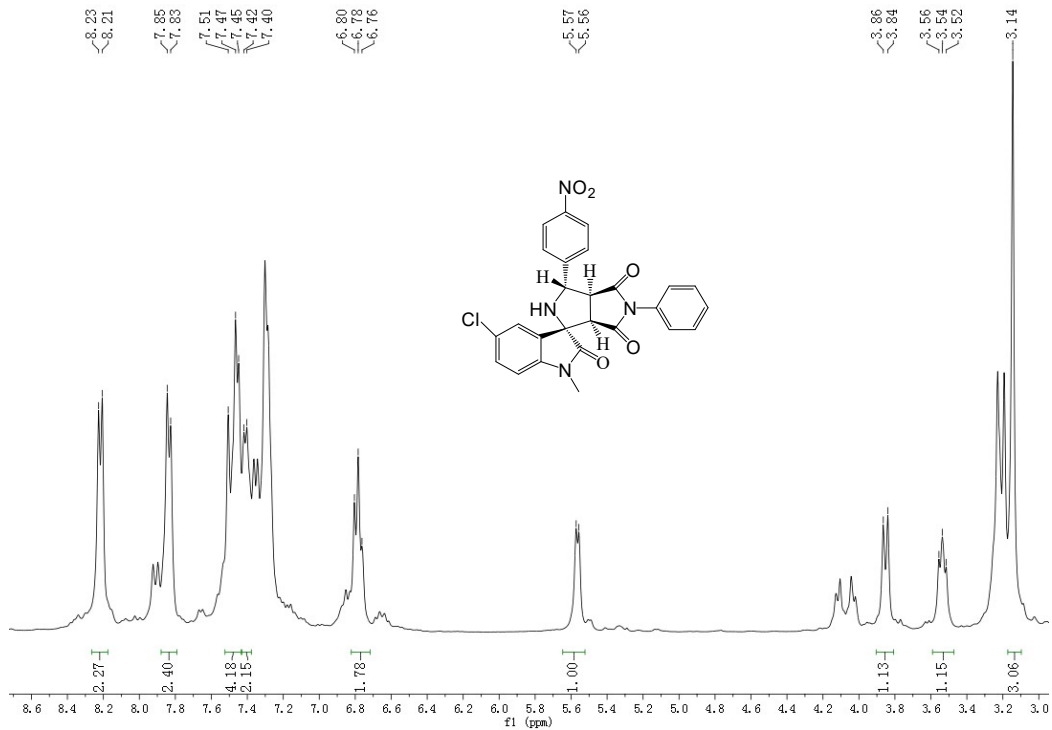
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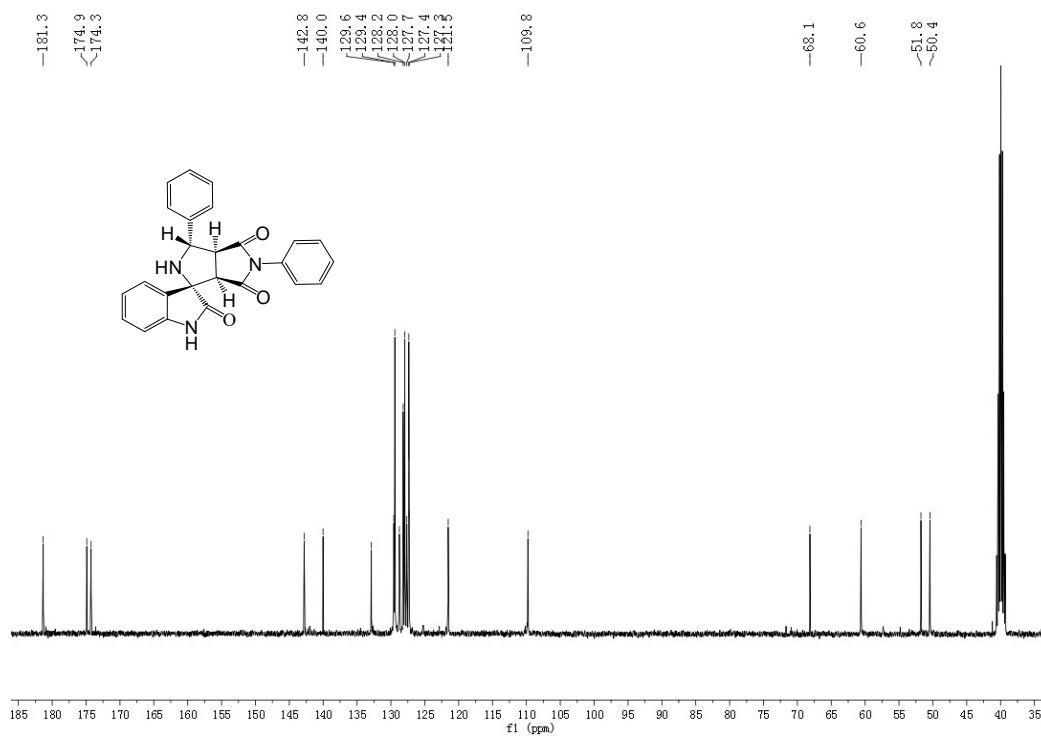
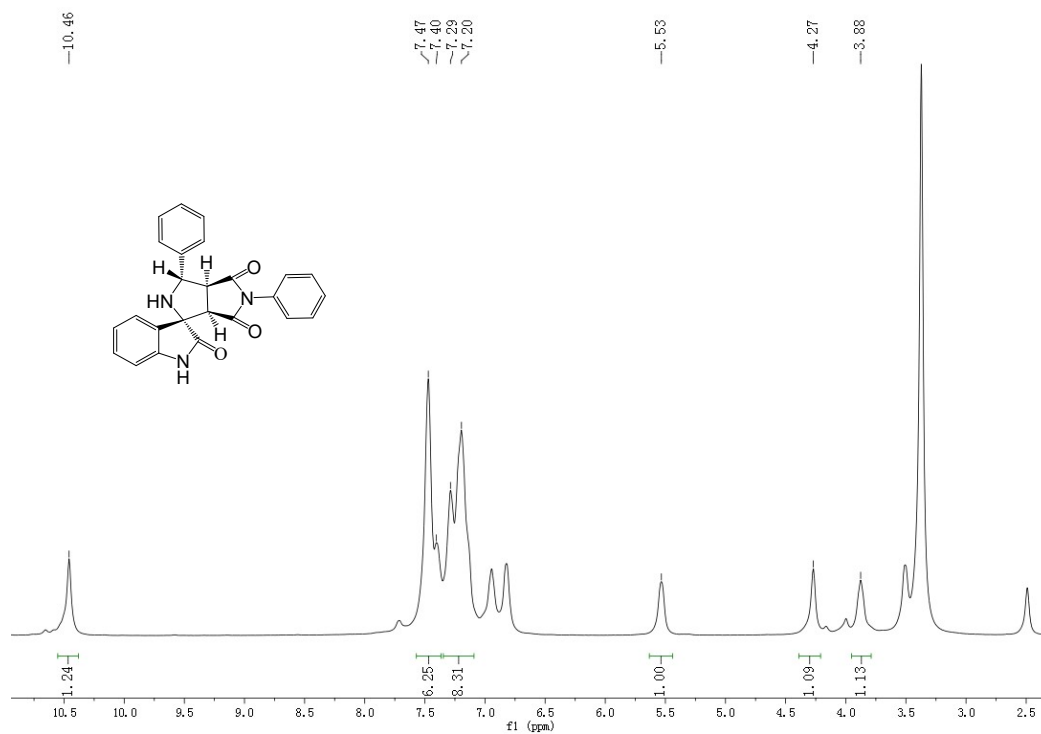
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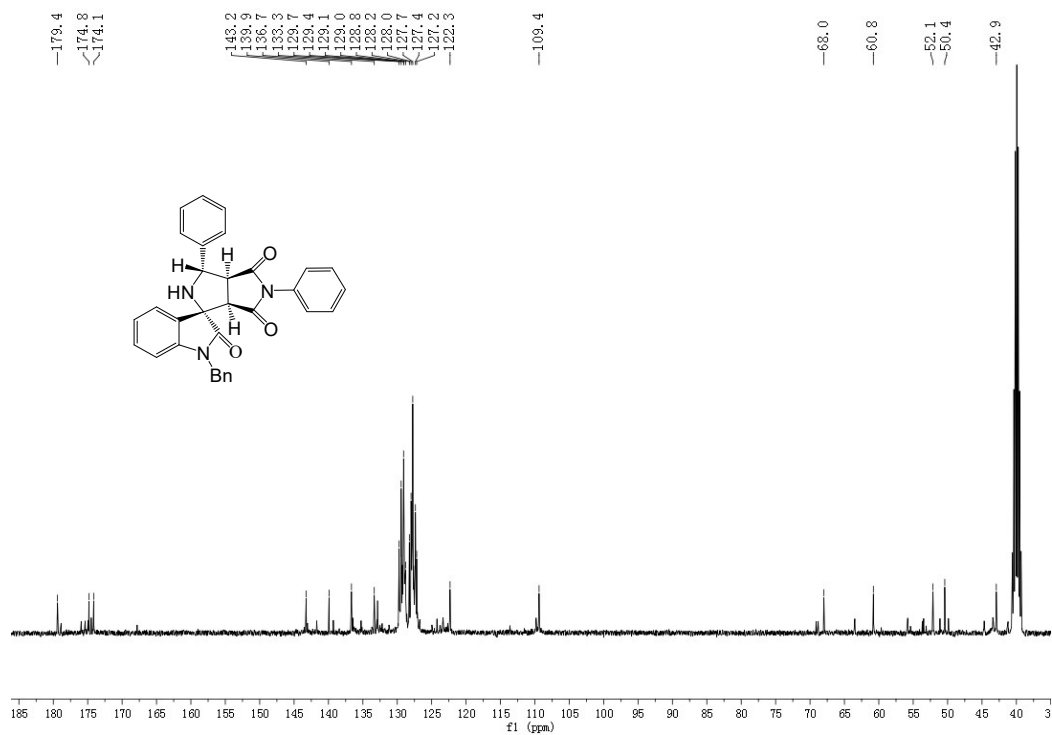
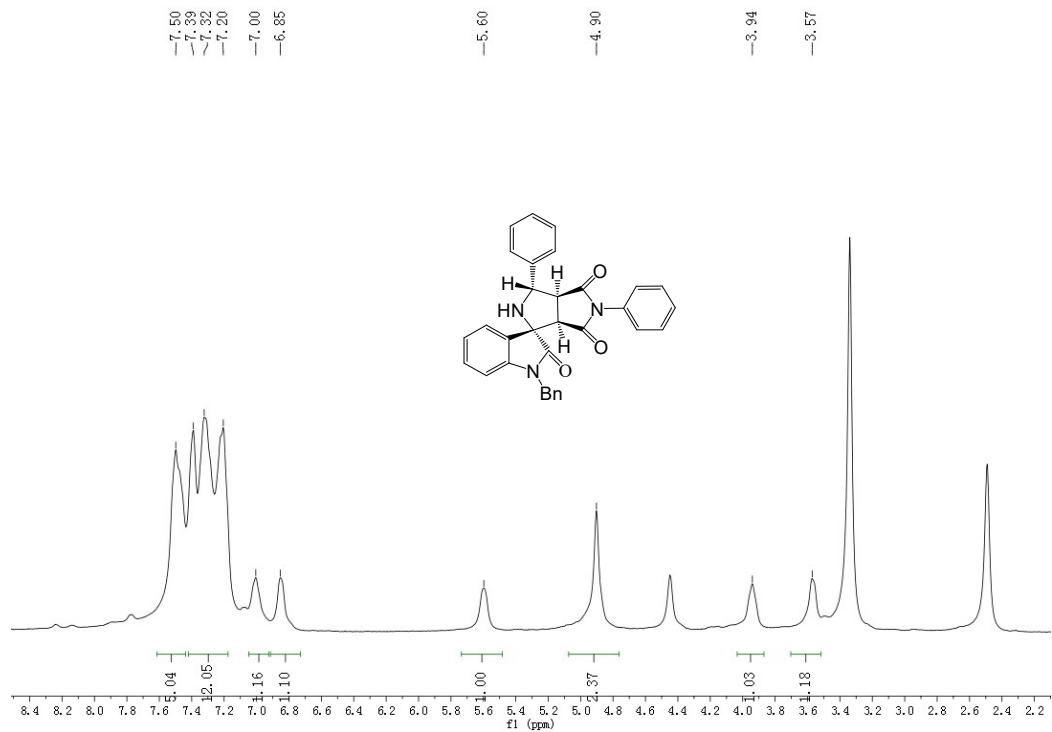
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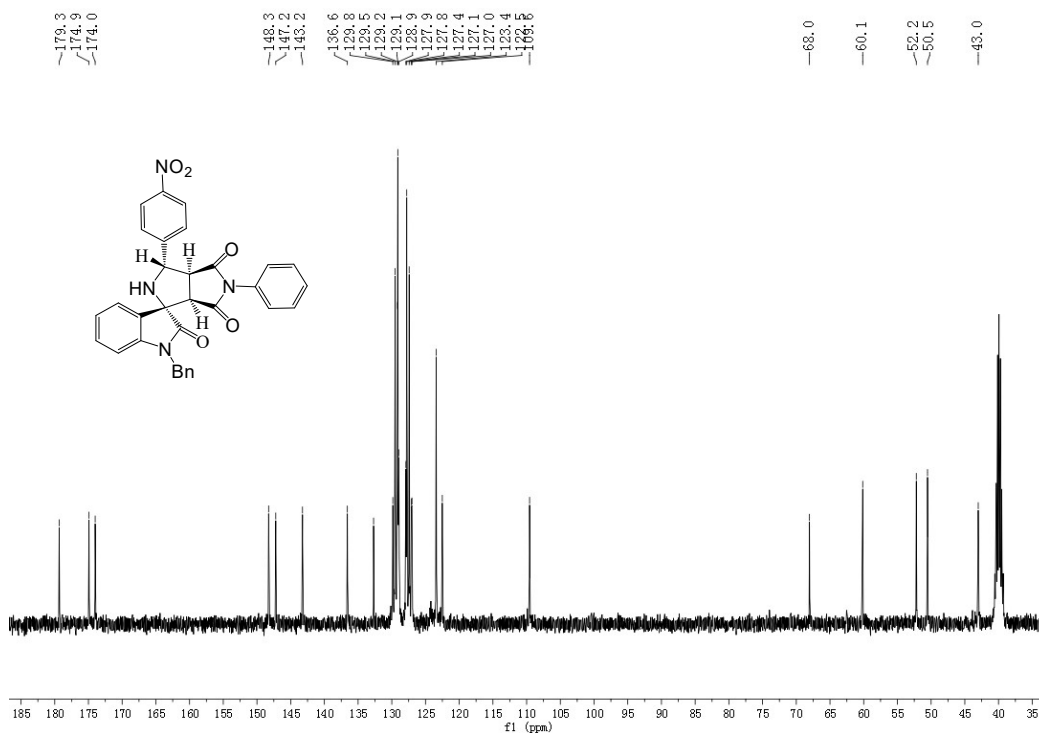
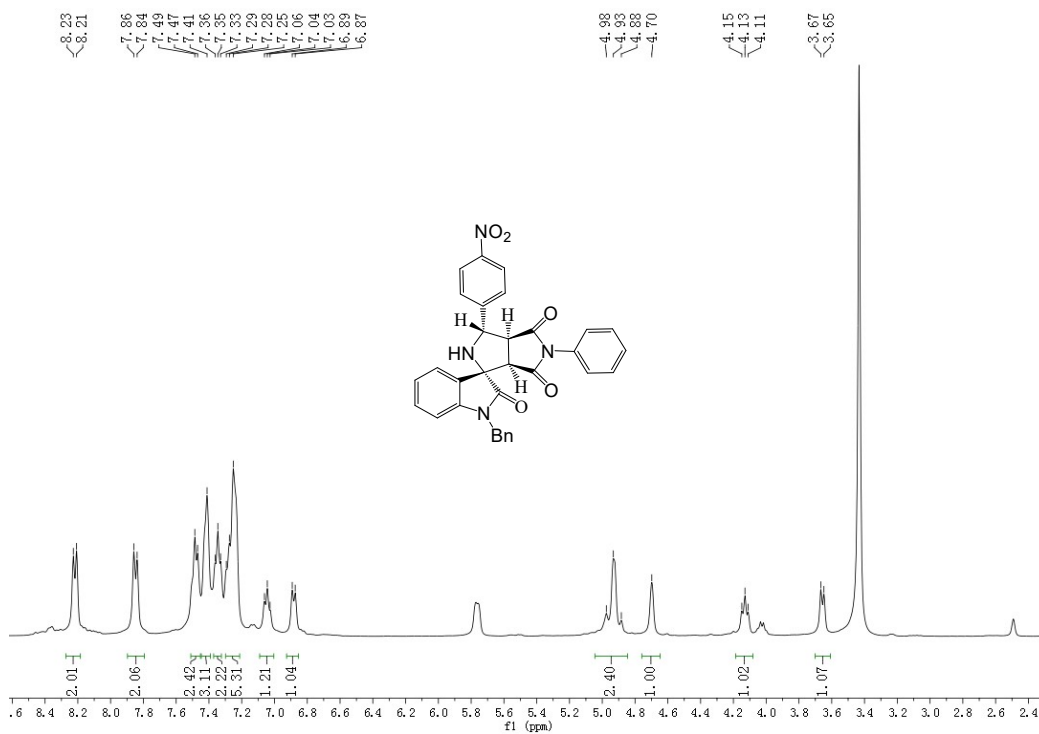
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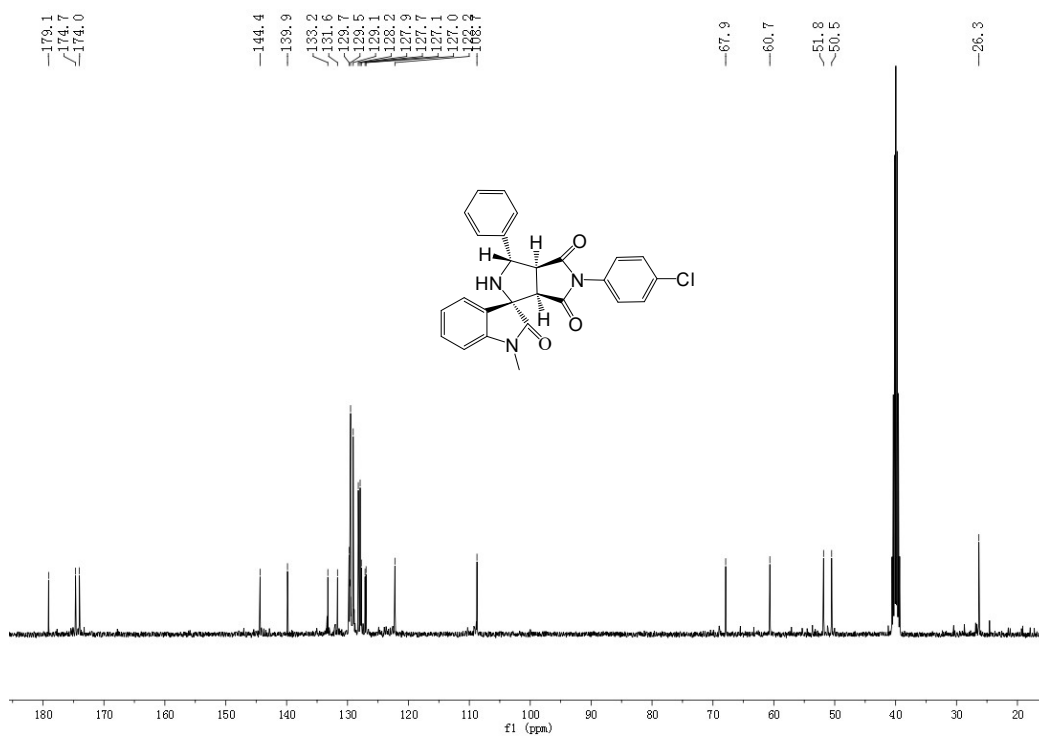
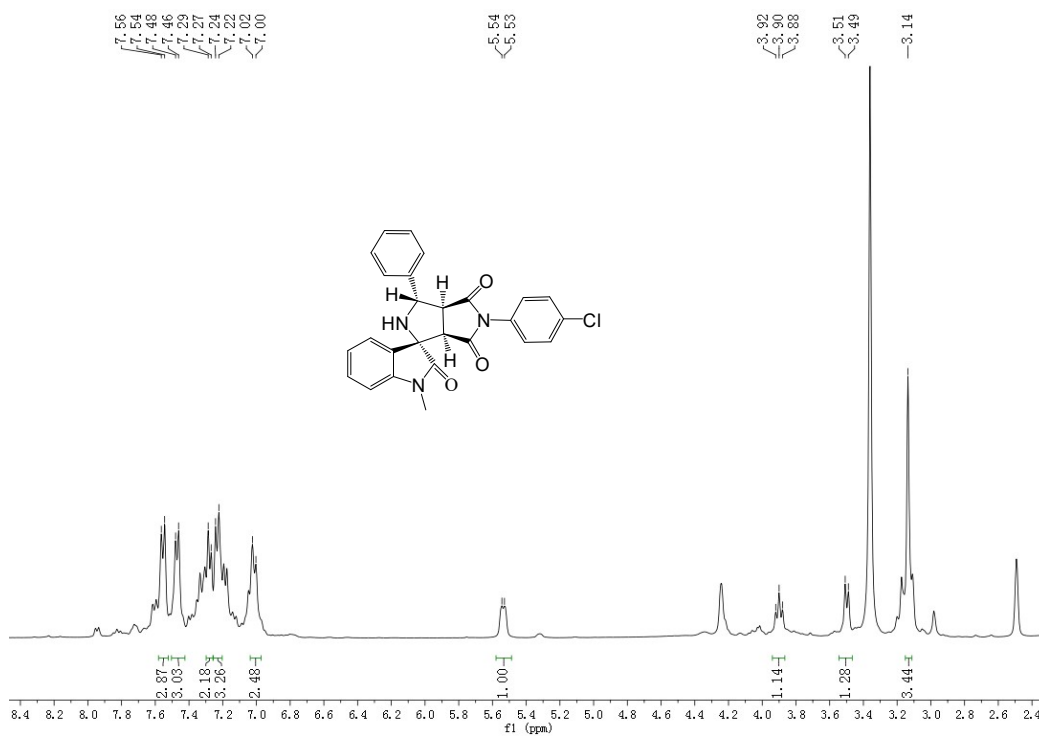
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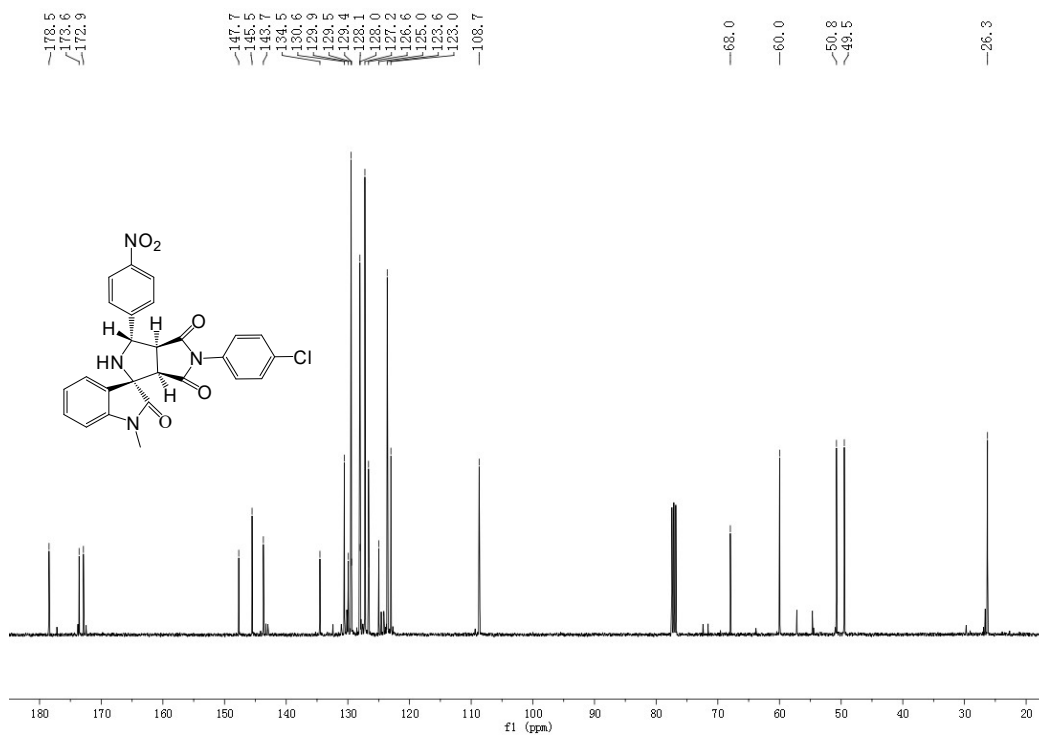
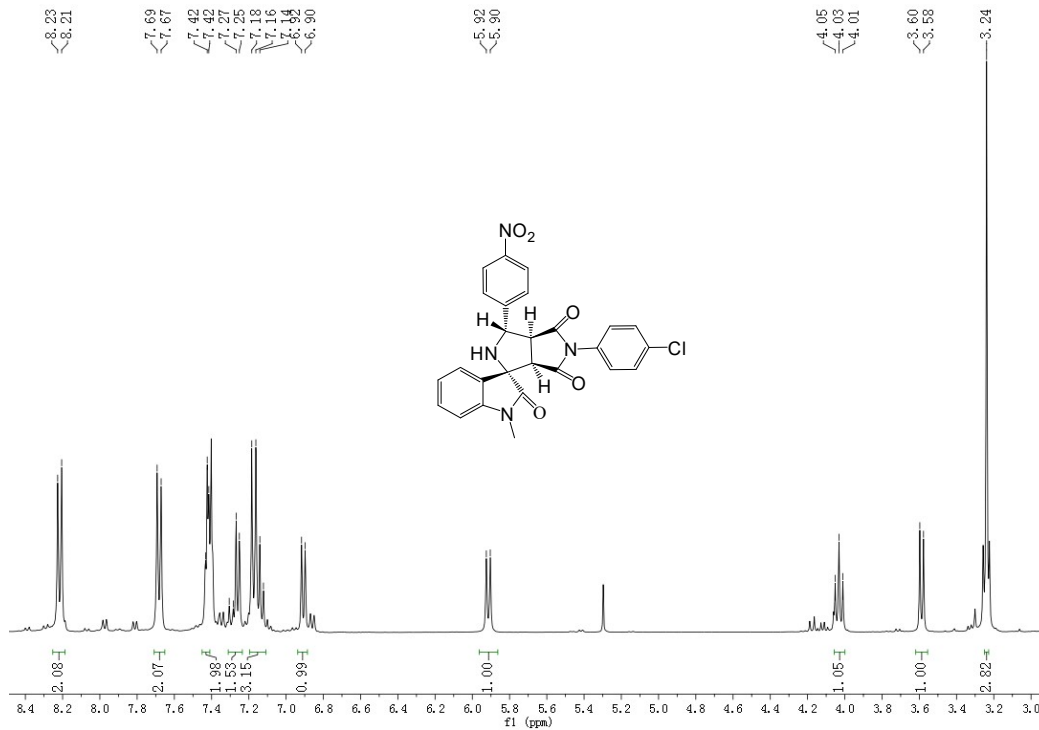
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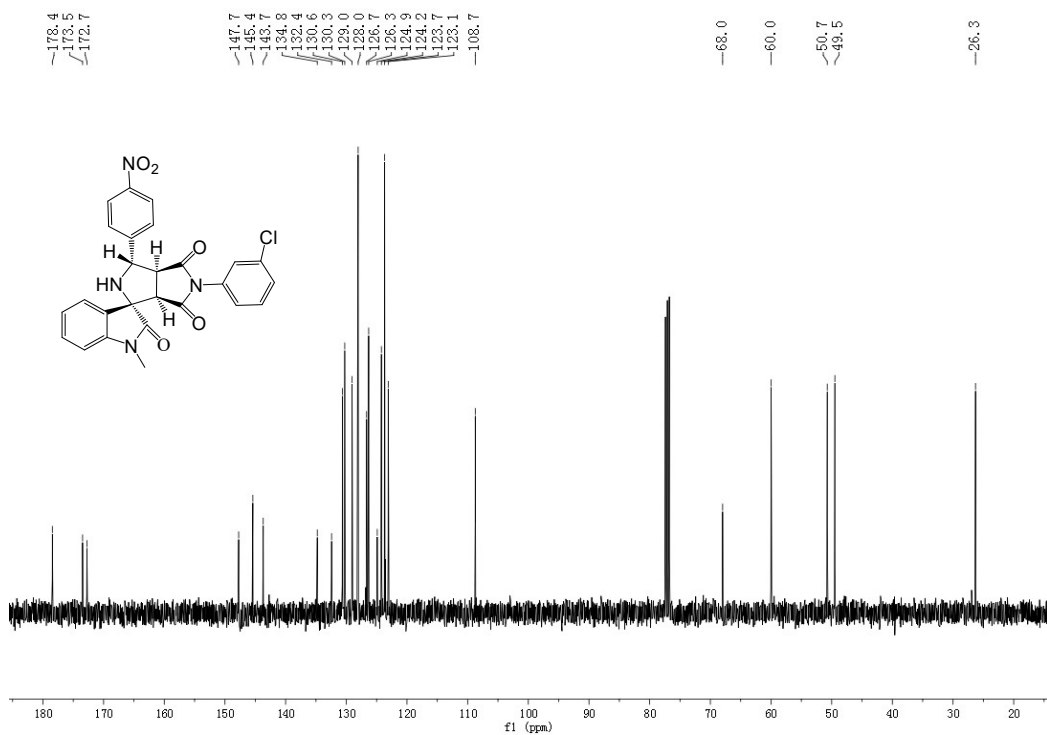
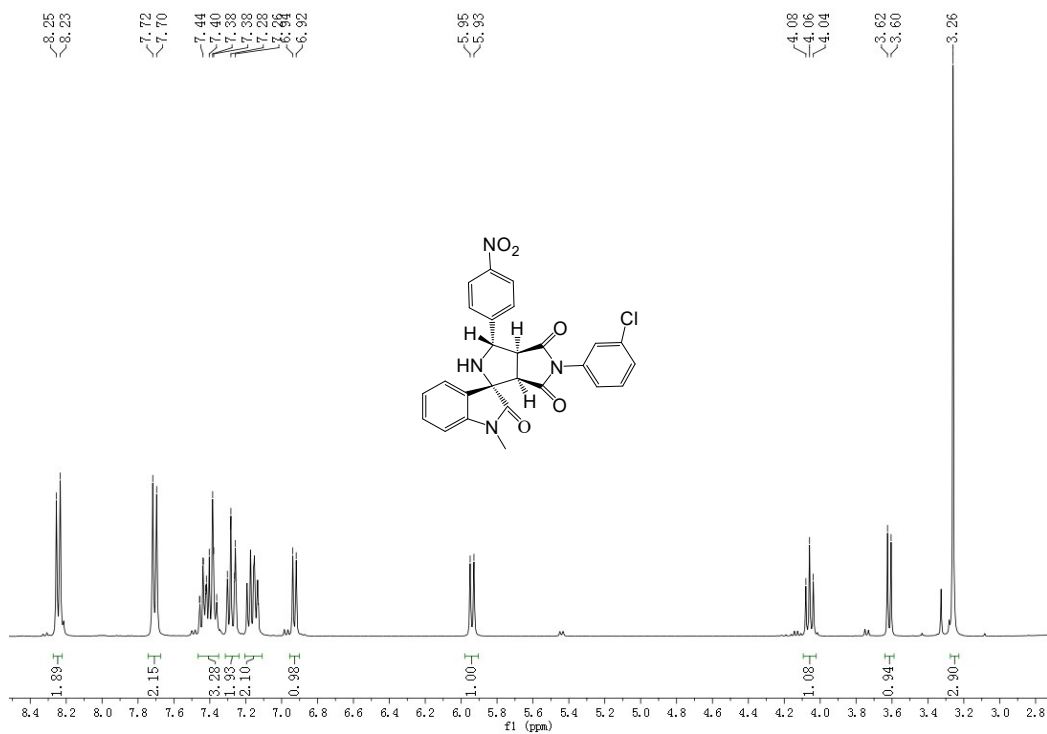
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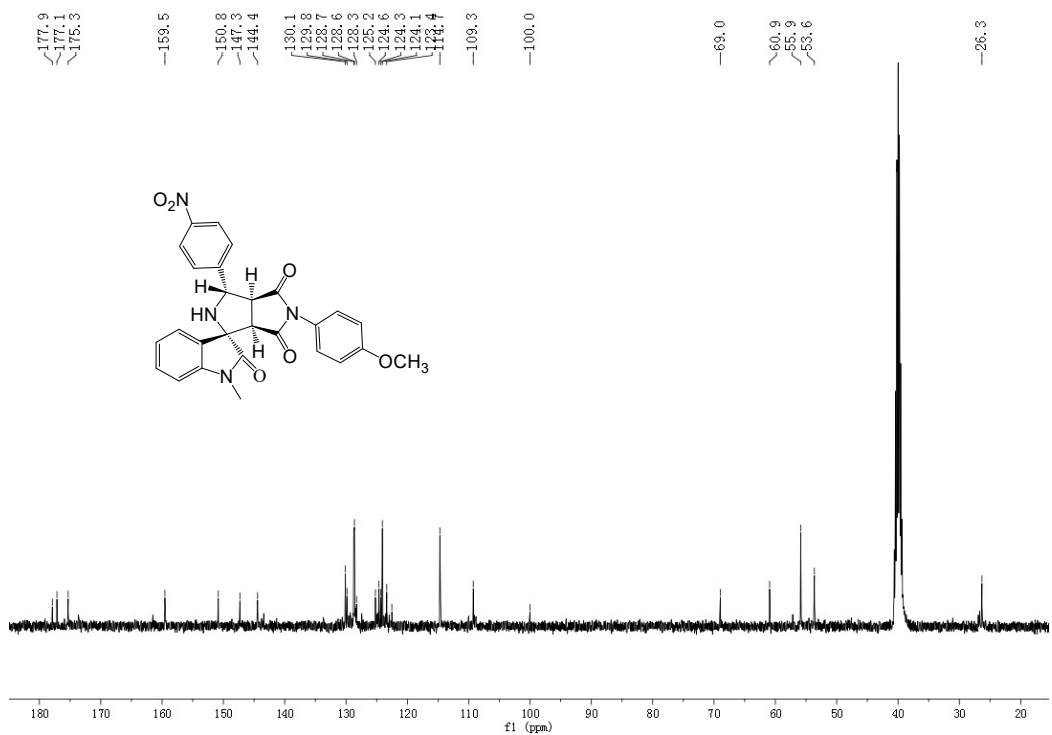
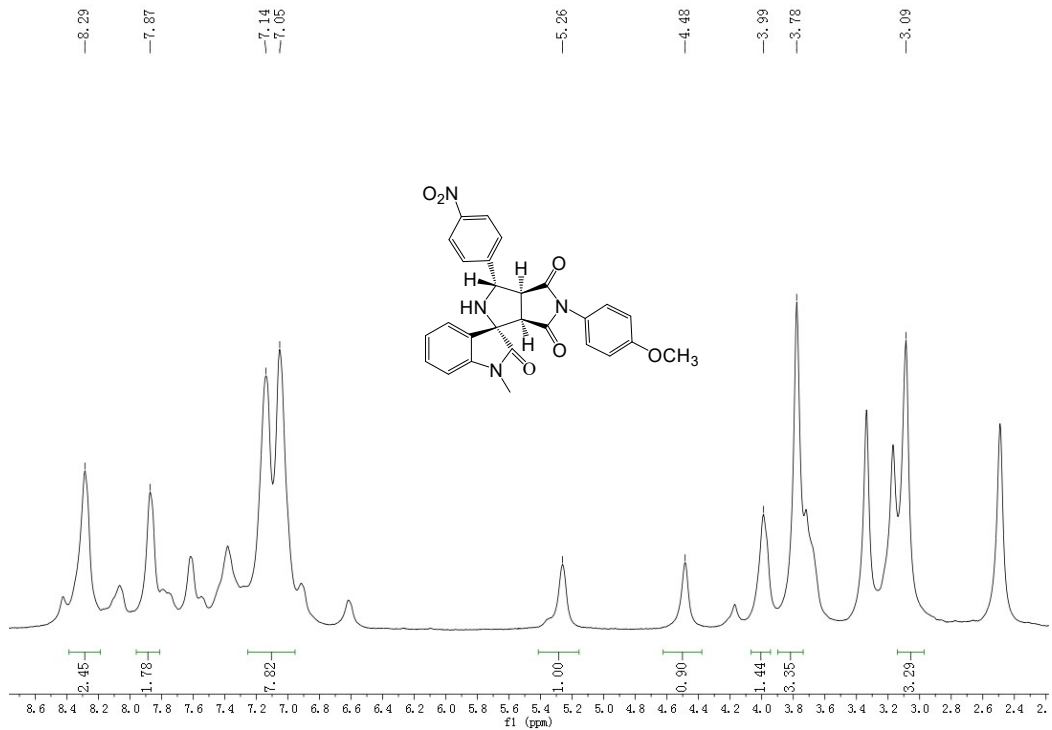
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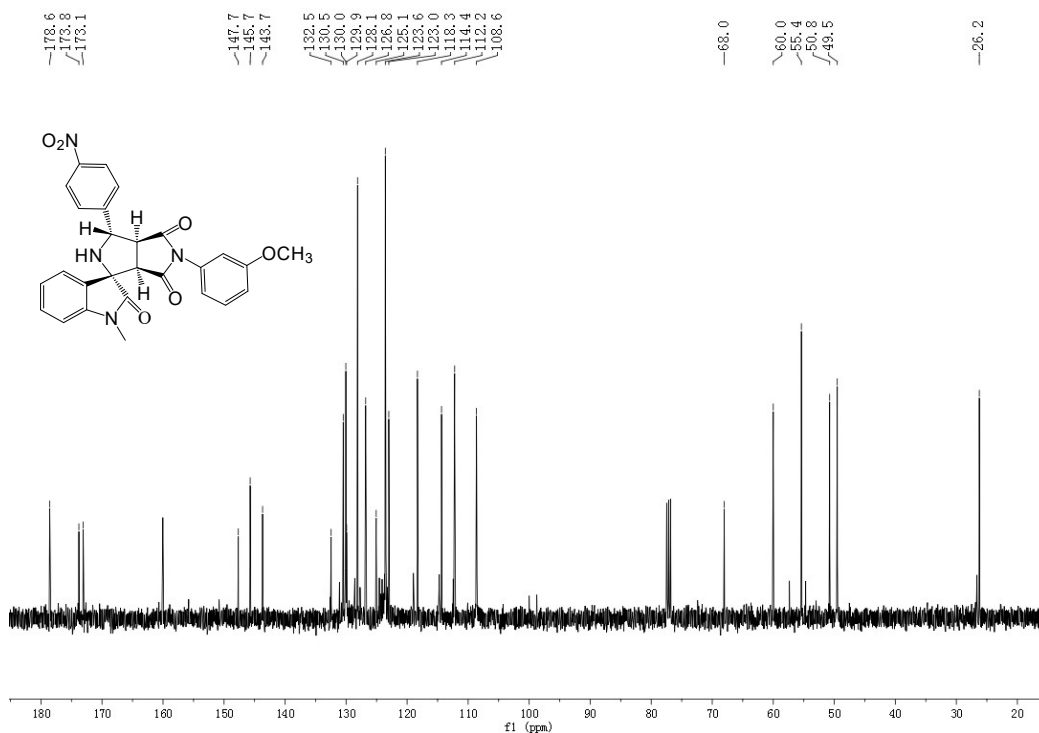
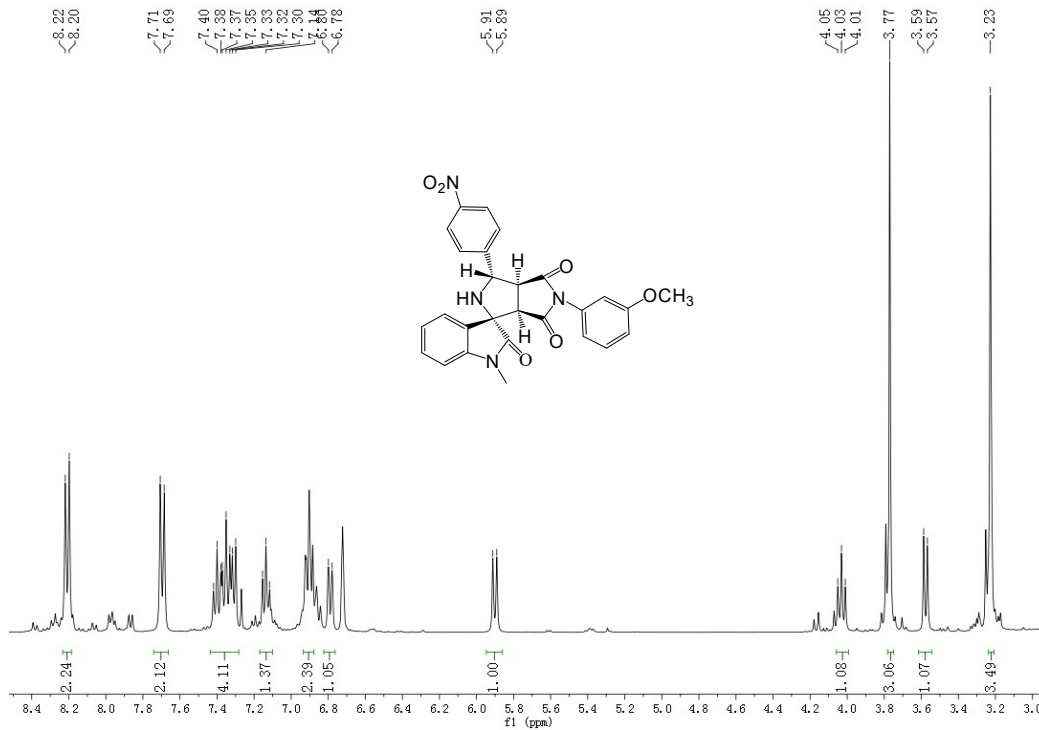
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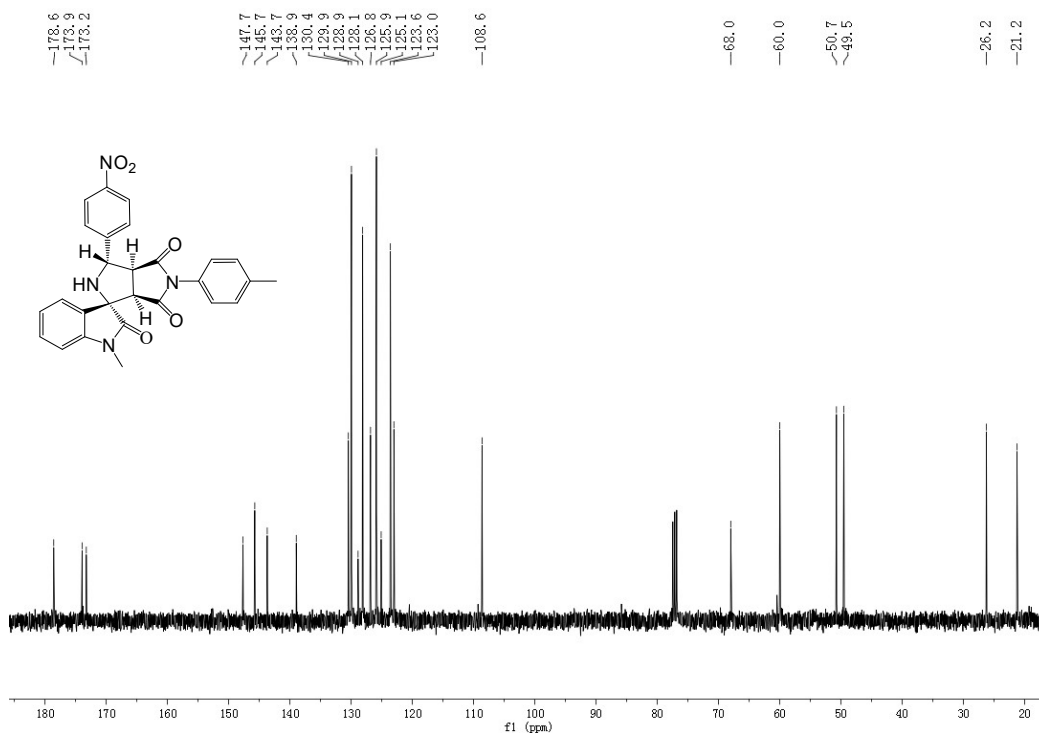
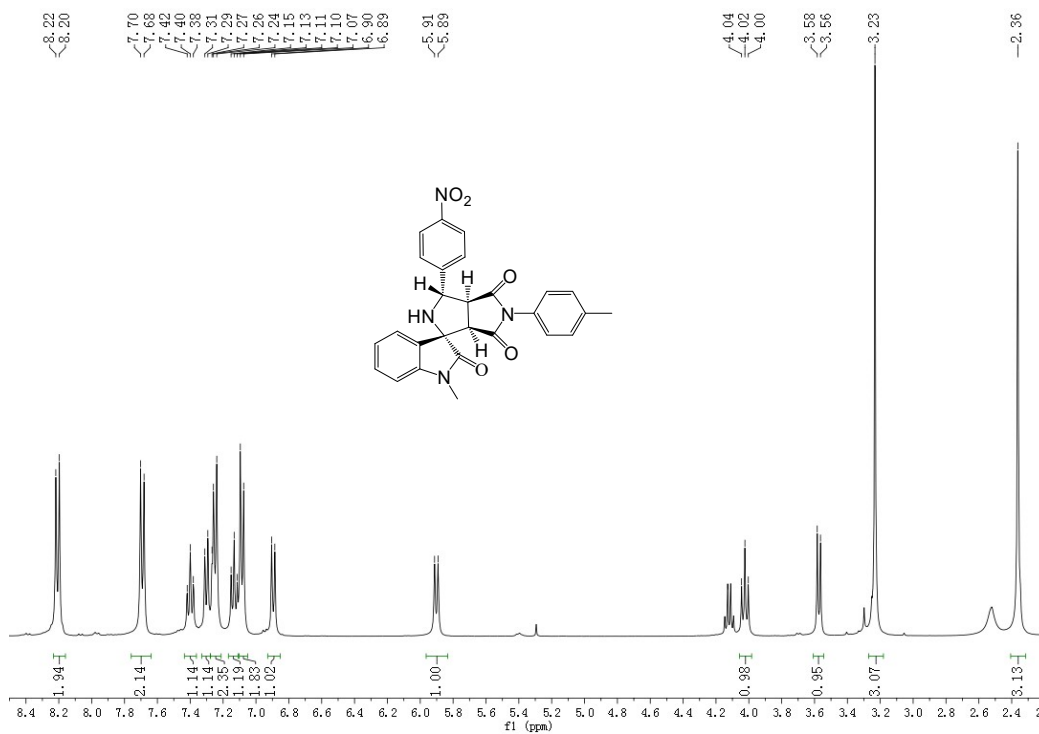
5k



51



5m



The crystallographic data of **4k**

Chemical Formula: C₂₆H₂₀BrN₃O₃

Molecular Weight: 502.3680

Temperature (K) 173(2)

Crystal system Monoclinic

Bond precision: C-C = 0.0042 Å Wavelength=0.71073

Cell: a=10.0641(5) b=10.7045(6) c=11.0214(7)

alpha=74.935(2) beta=73.244(2) gamma=81.118(2)

Temperature: 298 K

Calculated Reported

Volume 1093.80(11) 1093.80(11)

Space group P -1 P -1

Hall group -P 1 -P 1

Moiety formula C₂₆ H₂₀ Br N₃ O₃ C₂₆ H₂₀ Br N₃ O₃

Sum formula C₂₆ H₂₀ Br N₃ O₃ C₂₆ H₂₀ Br N₃ O₃

Mr 502.35 502.36

Dx, g cm⁻³ 1.525 1.525

Z 2 2

Mu (mm⁻¹) 1.914 1.914

F000 512.0 512.0

F000' 511.62

h, k, lmax 12,13,13 12,13,13

Nref 4500 4487

Tmin, Tmax 0.638,0.682 0.638,0.682

Tmin' 0.558

Correction method= # Reported T Limits: Tmin=0.638 Tmax=0.682

AbsCorr = MULTI-SCAN

Data completeness= 0.997 Theta(max)= 26.387

R(reflections)= 0.0405(3736) wR2(reflections)= 0.1183(4487)

S = 1.071 Npar= 30

