

Supporting Information
for
Eosin Y catalyzed visible-light-mediated aerobic
oxidative cyclization of *N,N*-dimethylanilines with
maleimides

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Experimental section and characterization of the
synthesized compounds

Table of contents

General remarks	S2
General procedure for the reactions of <i>N,N</i> -dimethylanilines with maleimides	S2
Photo of experimental setup	S3
Test for the production of hydrogen peroxide	S3
Characterization of products	S4
¹ H and ¹³ C NMR spectra of products	S11

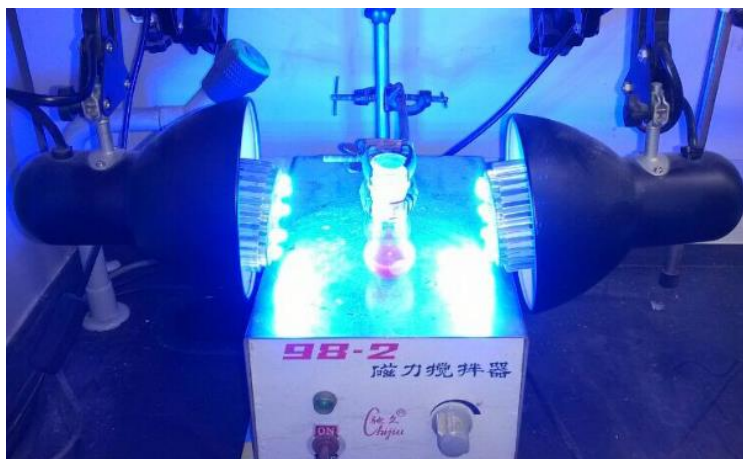
General remarks

N-Arylmaleimides were prepared according to literatures described procedures^{1,2} and other commercially available reagents were used as received without further purification. Flash column chromatography was carried out over silica gel (200–300 mesh). TLC was performed using Silica Gel GF254 plates and was visualized by fluorescence quenching at 254 nm. ¹H NMR and ¹³C NMR spectra were recorded on a Bruker ARX-400 MHz spectrometer in CDCl₃ using TMS as an internal reference with chemical shift values reported in ppm. All coupling constants (*J*) are reported in Hz. The high resolution mass spectra (HRMS) were measured on a Bruker Daltonics micrOTOF II spectrometer using ESI. Melting points were measured on a WRS-2A melting point apparatus. Two 9 W blue LEDs were used as a visible light source.

General procedure for the reactions of *N,N*-dimethylanilines with maleimides

To a 10 mL round bottom flask equipped with magnetic stirring bar was added *N,N*-dimethylaniline (**1**, 0.5 mmol, 2.0 equiv), maleimide (**2**, 0.25 mmol, 1.0 equiv), Eosin Y (water soluble, 0.03 equiv), and MeCN (3 mL). The solution was irradiated with two 9 W blue LEDs (distance app. 5 cm) at room temperature in an air atmosphere. After the completion of the reaction (indicated by TLC), the solvent was removed under reduced pressure. The residue was purified by flash column chromatography on silica gel (petroleum ether/ethyl acetate 15:1–10:1) to give the product **3**.

Photo of experimental setup



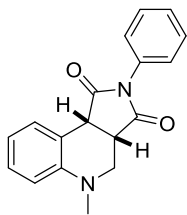
Test for the production of hydrogen peroxide

Hydrogen peroxide (H_2O_2) was detected after the reaction was completed by using KI/starch indicator.

				
(1)	(2)	(3)	(4)	(5)
KI/starch indicator	Addition of reaction mixture ^a at $t = 0$ to (1)	Reaction mixture ^a at $t = 18$ h	Addition of reaction mixture ^a at $t = 18$ h to (1)	Addition of 30% H_2O_2 to (1)
Ivory	Bright red	Red	Puce	Blue

^aReaction mixture of 0.5 mmol of **1a** and 0.25 mmol of **2a** in 3 mL of MeCN under optimized conditions.

Characterization of products



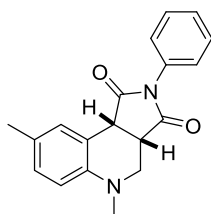
(3aS*,9bR*)-5-Methyl-2-phenyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3a)^{3,4}

White solid; m.p.: 183-185°C.

¹H NMR (400 MHz, CDCl₃): δ 7.52 (d, *J* = 7.4 Hz, 1H), 7.42 (t, *J* = 7.5 Hz, 2H), 7.35 (t, *J* = 7.3 Hz, 1H), 7.29-7.20 (m, 3H), 6.90 (td, *J* = 7.5, 1.1 Hz, 1H), 6.74 (d, *J* = 8.2 Hz, 1H), 4.15 (d, *J* = 9.5 Hz, 1H), 3.60 (dd, *J* = 11.5, 2.7 Hz, 1H), 3.52 (ddd, *J* = 9.5, 4.4, 2.7 Hz, 1H), 3.11 (dd, *J* = 11.5, 4.4 Hz, 1H), 2.83 (s, 3H).

¹³C NMR (101 MHz, CDCl₃): δ 177.7, 175.7, 148.5, 131.9, 130.3, 128.9, 128.6, 128.4, 126.3, 119.6, 118.5, 112.5, 50.6, 43.5, 42.1, 39.4.

HRMS (ESI): calcd for C₁₈H₁₆N₂O₂ + Na = 315.1109, found: 315.1099.



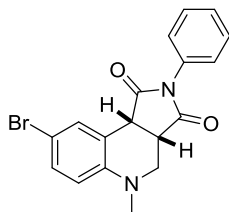
(3aS*,9bR*)-5,8-Dimethyl-2-phenyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3b)^{3,5}

White solid; m.p.: 187-189°C.

¹H NMR (400 MHz, CDCl₃): δ 7.41 (t, *J* = 7.4 Hz, 2H), 7.35 (d, *J* = 9.2 Hz, 2H), 7.26 (d, *J* = 7.2 Hz, 2H), 7.03 (d, *J* = 8.0 Hz, 1H), 6.64 (d, *J* = 8.3 Hz, 1H), 4.09 (d, *J* = 9.5 Hz, 1H), 3.57 (dd, *J* = 11.4, 2.7 Hz, 1H), 3.53-3.46 (m, 1H), 3.03 (dd, *J* = 11.4, 4.4 Hz, 1H), 2.79 (s, 3H), 2.30 (s, 3H).

¹³C NMR (101 MHz, CDCl₃): δ 177.8, 175.8, 146.3, 132.0, 130.8, 129.2, 128.94, 128.4, 126.3, 118.5, 112.5, 50.9, 43.5, 42.1, 39.5, 20.4.

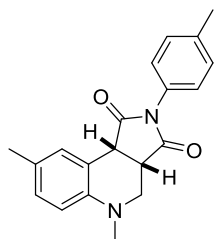
HRMS (ESI): calcd for C₁₉H₁₈N₂O₂ + Na = 329.1266, found: 329.1254.



(3aS*,9bR*)-8-Bromo-5-methyl-2-phenyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3c)⁵

White solid; m.p.: 156-168°C.

^1H NMR (400 MHz, CDCl_3): δ 7.63 (d, J = 1.7 Hz, 1H), 7.47-7.33 (m, 3H), 7.33-7.23 (m, 3H), 6.60 (d, J = 8.8 Hz, 1H), 4.08 (d, J = 9.6 Hz, 1H), 3.59 (dd, J = 11.5, 2.8 Hz, 1H), 3.51 (ddd, J = 9.6, 4.4, 2.8 Hz, 1H), 3.09 (dd, J = 11.6, 4.4 Hz, 1H), 2.80 (s, 3H).
 ^{13}C NMR (101 MHz, CDCl_3): δ 177.2, 175.1, 147.4, 132.6, 131.8, 131.4, 129.0, 128.6, 126.2, 120.3, 114.2, 111.6, 50.3, 43.21, 41.7, 39.4.
 HRMS (ESI): calcd for $\text{C}_{18}\text{H}_{15}\text{BrN}_2\text{O}_2 + \text{Na} = 393.0215$, found: 393.0209.



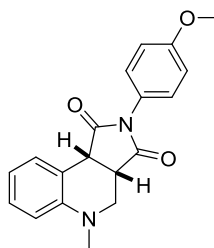
(3aS*,9bR*)-5,8-Dimethyl-2-(p-tolyl)-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3d)⁵

White solid; m.p.: 199-201°C.

^1H NMR (400 MHz, CDCl_3): δ 7.34 (d, J = 2.0 Hz, 1H), 7.27-7.18 (m, 2H), 7.13 (d, J = 8.3 Hz, 2H), 7.03 (dd, J = 8.3, 2.1 Hz, 1H), 6.65 (d, J = 8.3 Hz, 1H), 4.10 (d, J = 9.5 Hz, 1H), 3.57 (dd, J = 11.4, 2.7 Hz, 1H), 3.50 (ddd, J = 9.5, 4.4, 2.7 Hz, 1H), 3.04 (dd, J = 11.4, 4.4 Hz, 1H), 2.80 (s, 3H), 2.35 (s, 3H), 2.30 (s, 3H).

^{13}C NMR (101 MHz, CDCl_3): δ 177.9, 175.9, 146.3, 138.5, 130.8, 129.6, 129.2, 128.9, 126.1, 118.5, 112.5, 50.9, 43.5, 42.1, 39.53, 21.2, 20.42.

HRMS (ESI): calcd for $\text{C}_{20}\text{H}_{20}\text{N}_2\text{O}_2 + \text{Na} = 343.1422$, found: 343.1417.



(new compound)

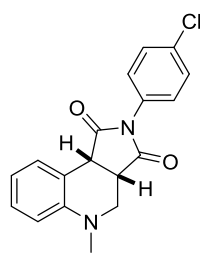
(3aS*,9bR*)-2-(4-Methoxyphenyl)-5-methyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3e)

White solid; m.p.: 192-194°C.

^1H NMR (400 MHz, CDCl_3): δ 7.51 (d, J = 7.2 Hz, 1H), 7.20 (dd, J = 29.7, 9.1 Hz, 3H), 6.95-6.87 (m, 3H), 6.74 (d, J = 8.1 Hz, 1H), 4.13 (d, J = 9.6 Hz, 1H), 3.79 (s, 3H), 3.59 (dd, J = 11.5, 2.7 Hz, 1H), 3.50 (ddd, J = 9.6, 4.4, 2.7 Hz, 1H), 3.10 (dd, J = 11.5, 4.4 Hz, 1H), 2.82 (s, 3H).

^{13}C NMR (101 MHz, CDCl_3): δ 177.9, 176.0, 159.3, 148.4, 130.3, 128.6, 127.5, 124.6, 119.6, 118.6, 114.2, 112.5, 55.4, 50.6, 43.4, 42.0, 39.4.

HRMS (ESI): calcd for $\text{C}_{19}\text{H}_{18}\text{N}_2\text{O}_3 + \text{Na} = 345.1215$, found: 345.1210.



(new compound)

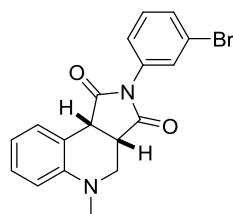
(3aS*,9bR*)-2-(4-Chlorophenyl)-5-methyl-3a,4,5,9b-tetrahydro-1H-pyrrolo-[3,4-c]quinoline-1,3(2H)-dione (3f)

White solid; m.p.: 191-193°C.

^1H NMR (400 MHz, CDCl_3): δ 7.51 (d, J = 7.5 Hz, 1H), 7.40 (d, J = 8.8 Hz, 2H), 7.27-7.21 (m, 3H), 6.91 (td, J = 7.7, 1.1 Hz, 1H), 6.75 (d, J = 8.2 Hz, 1H), 4.17 (d, J = 9.6 Hz, 1H), 3.61 (dd, J = 11.5, 2.6 Hz, 1H), 3.54 (ddd, J = 9.6, 4.3, 2.6 Hz, 1H), 3.12 (dd, J = 11.5, 4.4 Hz, 1H), 2.83 (s, 3H).

^{13}C NMR (101 MHz, CDCl_3): δ 177.4, 175.5, 148.5, 134.2, 130.4, 130.3, 129.1, 128.8, 127.5, 119.7, 118.3, 112.6, 50.6, 43.6, 42.1, 39.4.

HRMS (ESI): calcd for $\text{C}_{18}\text{H}_{15}\text{ClN}_2\text{O}_2 + \text{Na}$ = 349.0720, found: 315.0721.



(new compound)

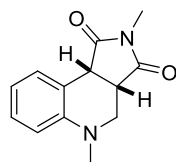
(3aS*,9bR*)-2-(3-Bromophenyl)-5-methyl-3a,4,5,9b-tetrahydro-1H-pyrrolo-[3,4-c]quinoline-1,3(2H)-dione (3g)

White solid; m.p.: 164-166°C.

^1H NMR (400 MHz, CDCl_3): δ 7.54-7.45 (m, 3H), 7.33-7.21 (m, 3H), 6.91 (t, J = 7.4 Hz, 1H), 6.75 (d, J = 8.1 Hz, 1H), 4.16 (d, J = 9.6 Hz, 1H), 3.61 (dd, J = 11.5, 2.6 Hz, 1H), 3.54 (ddd, J = 9.6, 4.3, 2.6 Hz, 1H), 3.11 (dd, J = 11.5, 4.3 Hz, 1H), 2.84 (s, 3H).

^{13}C NMR (101 MHz, CDCl_3): δ 177.3, 175.3, 148.5, 133.1, 131.6, 130.3, 130.1, 129.4, 128.8, 125.0, 122.2, 119.7, 118.3, 112.6, 50.6, 43.6, 42.1, 39.4.

HRMS (ESI): calcd for $\text{C}_{18}\text{H}_{15}\text{BrN}_2\text{O}_2 + \text{Na}$ = 393.0215, found: 393.0215.



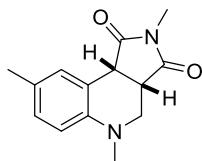
(3aS*,9bR*)-2,5-Dimethyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3h)³

White solid; m.p.: 170-172°C.

^1H NMR (400 MHz, CDCl_3): δ 7.48 (d, J = 7.5 Hz, 1H), 7.24-7.18 (m, 1H), 6.89 (td, J = 7.5, 1.1 Hz, 1H), 6.70 (d, J = 8.2 Hz, 1H), 4.00 (d, J = 9.4 Hz, 1H), 3.54 (dd, J = 11.5, 2.4 Hz, 1H), 3.37 (ddd, J = 9.5, 4.5, 2.4 Hz, 1H), 3.03 (dd, J = 11.5, 4.4 Hz, 1H), 2.99 (s, 3H), 2.79 (s, 3H).

^{13}C NMR (101 MHz, CDCl_3): δ 178.7, 176.8, 148.4, 130.2, 128.5, 119.6, 118.7, 112.4, 50.4, 43.6, 42.0, 39.4, 25.3.

HRMS (ESI): calcd for $\text{C}_{13}\text{H}_{14}\text{N}_2\text{O}_2 + \text{Na} = 253.0953$, found: 253.0947.



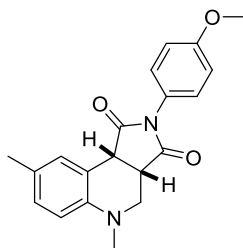
(3aS*,9bR*)-2,5,8-Trimethyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3i)⁴

White solid; m.p.: 178-180°C.

^1H NMR (400 MHz, CDCl_3): δ 7.29 (d, $J = 2.1$ Hz, 1H), 7.01 (dd, $J = 8.3, 2.1$ Hz, 1H), 6.60 (d, $J = 8.3$ Hz, 1H), 3.95 (d, $J = 9.4$ Hz, 1H), 3.51 (dd, $J = 11.4, 2.4$ Hz, 1H), 3.33 (ddd, $J = 9.3, 4.3, 2.2$ Hz, 1H), 3.00-2.93 (m, 4H), 2.75 (s, 3H), 2.29 (s, 3H).

^{13}C NMR (101 MHz, CDCl_3): δ 178.8, 176.8, 146.2, 130.6, 129.1, 128.9, 118.6, 112.4, 50.7, 43.5, 42.0, 39.5, 25.3, 20.4.

HRMS (ESI): calcd for $\text{C}_{14}\text{H}_{16}\text{N}_2\text{O}_2 + \text{Na} = 267.1109$, found: 267.1104.



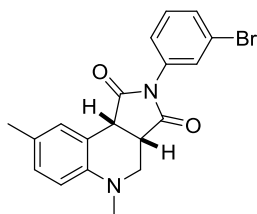
(3aS*,9bR*)-2-(4-Methoxyphenyl)-5,8-dimethyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3j)⁴

White solid; m.p.: 176-178°C.

^1H NMR (400 MHz, CDCl_3): δ 7.33 (d, $J = 2.1$ Hz, 1H), 7.16 (d, $J = 9.0$ Hz, 2H), 7.03 (d, $J = 8.8$ Hz, 1H), 6.92 (d, $J = 9.0$ Hz, 2H), 6.64 (d, $J = 8.3$ Hz, 1H), 4.07 (d, $J = 9.5$ Hz, 1H), 3.78 (s, 3H), 3.56 (dd, $J = 11.4, 2.7$ Hz, 1H), 3.47 (ddd, $J = 9.6, 4.4, 2.6$ Hz, 1H), 3.02 (dd, $J = 11.4, 4.4$ Hz, 1H), 2.78 (s, 3H), 2.29 (s, 3H).

^{13}C NMR (101 MHz, CDCl_3): δ 178.0, 176.0, 159.3, 146.3, 130.7, 129.1, 128.9, 127.5, 124.6, 118.5, 114.2, 112.4, 55.4, 50.9, 43.4, 42.0, 39.5, 20.4.

HRMS (ESI): calcd for $\text{C}_{20}\text{H}_{20}\text{N}_2\text{O}_3 + \text{Na} = 359.1372$, found: 359.1366.



(new compound)

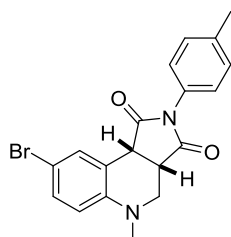
(3aS*,9bR*)-2-(3-Bromophenyl)-5,8-dimethyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3k)

White solid; m.p.: 174-176°C.

¹H NMR (400 MHz, CDCl₃): δ 7.52-7.45 (m, 2H), 7.34-7.23 (m, 3H), 7.04 (d, *J* = 7.5 Hz, 1H), 6.65 (d, *J* = 8.3 Hz, 1H), 4.11 (d, *J* = 9.6 Hz, 1H), 3.58 (dd, *J* = 11.4, 2.6 Hz, 1H), 3.51 (ddd, *J* = 9.6, 4.3, 2.5 Hz, 1H), 3.04 (dd, *J* = 11.4, 4.3 Hz, 1H), 2.80 (s, 3H), 2.30 (s, 3H).

¹³C NMR (101 MHz, CDCl₃): δ 177.3, 175.4, 146.3, 131.5, 130.7, 130.1, 129.4, 129.3, 129.0, 125.0, 122.2, 118.2, 112.6, 50.9, 43.6, 42.2, 39.5, 20.4.

HRMS (ESI): calcd for C₁₉H₁₇BrN₂O₂ + Na = 407.0371, found: 407.0366.



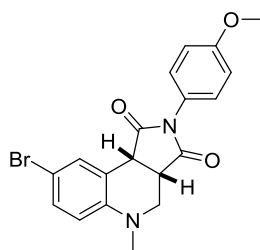
(3a*S,9b*R**)-8-Bromo-5-methyl-2-(p-tolyl)-3a,4,5,9b-tetrahydro-1*H*-pyrrolo[3,4-*c*]quinoline-1,3(2*H*)-dione (3l)⁵**

White solid; m.p.: 193-195°C.

¹H NMR (400 MHz, CDCl₃): δ 7.62 (d, *J* = 1.5 Hz, 1H), 7.30 (dd, *J* = 8.8, 2.3 Hz, 1H), 7.26-7.20 (m, 2H), 7.12 (d, *J* = 8.4 Hz, 2H), 6.59 (d, *J* = 8.8 Hz, 1H), 4.06 (d, *J* = 9.6 Hz, 1H), 3.58 (dd, *J* = 11.5, 2.8 Hz, 1H), 3.50 (ddd, *J* = 9.6, 4.4, 2.8 Hz, 1H), 3.08 (dd, *J* = 11.5, 4.4 Hz, 1H), 2.80 (s, 3H), 2.35 (s, 3H).

¹³C NMR (101 MHz, CDCl₃): δ 177.3, 175.2, 147.4, 138.6, 132.7, 131.3, 129.6, 126.0, 120.4, 114.1, 50.3, 43.2, 41.7, 39.4, 21.1.

HRMS (ESI): calcd for C₁₉H₁₇BrN₂O₂ + Na = 407.0371, found: 407.0366.



(new compound)

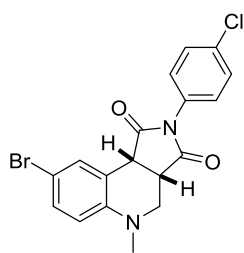
(3a*S,9b*R**)-8-Bromo-2-(4-methoxyphenyl)-5-methyl-3a,4,5,9b-tetrahydro-1*H*-pyrrolo[3,4-*c*]quinoline-1,3(2*H*)-dione (3m)**

White solid; m.p.: 205-207°C.

¹H NMR (400 MHz, CDCl₃): δ 7.62 (d, *J* = 1.7 Hz, 1H), 7.30 (dd, *J* = 8.8, 2.4 Hz, 1H), 7.16 (d, *J* = 9.0 Hz, 2H), 6.93 (d, *J* = 9.0 Hz, 2H), 6.59 (d, *J* = 8.8 Hz, 1H), 4.06 (d, *J* = 9.6 Hz, 1H), 3.79 (s, 3H), 3.58 (dd, *J* = 11.6, 2.8 Hz, 1H), 3.49 (ddd, *J* = 9.6, 4.4, 2.8 Hz, 1H), 3.08 (dd, *J* = 11.5, 4.4 Hz, 1H), 2.80 (s, 3H).

¹³C NMR (101 MHz, CDCl₃): δ 177.4, 175.3, 159.4, 147.4, 132.7, 131.4, 127.5, 120.4, 114.3, 114.2, 111.6, 55.5, 50.4, 43.2, 41.7, 39.4.

HRMS (ESI): calcd for C₁₉H₁₇BrN₂O₃ + Na = 423.0320, found: 423.0315.



(new compound)

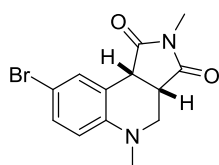
(3aS*,9bR*)-8-Bromo-2-(4-chlorophenyl)-5-methyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3n)

White solid; m.p.: 218-220°C.

^1H NMR (400 MHz, CDCl_3): δ 7.62 (d, J = 1.8 Hz, 1H), 7.40 (d, J = 8.7 Hz, 2H), 7.31 (dd, J = 8.7, 2.3 Hz, 1H), 7.27-7.20 (m, 2H), 6.60 (d, J = 8.8 Hz, 1H), 4.09 (d, J = 9.6 Hz, 1H), 3.60 (dd, J = 11.5, 2.7 Hz, 1H), 3.52 (ddd, J = 9.6, 4.3, 2.7 Hz, 1H), 3.09 (dd, J = 11.5, 4.4 Hz, 1H), 2.81 (s, 3H).

^{13}C NMR (101 MHz, CDCl_3): δ 176.9, 174.8, 147.4, 134.3, 132.6, 131.5, 129.2, 127.5, 120.1, 114.3, 111.7, 50.3, 43.3, 41.7, 39.4.

HRMS (ESI): calcd for $\text{C}_{18}\text{H}_{14}\text{BrClN}_2\text{O}_2 + \text{Na}$ = 426.9825, found: 426.9819.



(new compound)

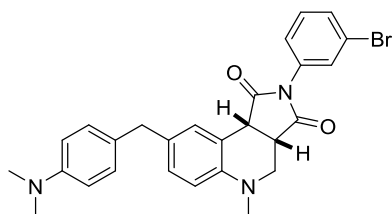
(3aS*,9bR*)-8-Bromo-2,5-dimethyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3o)

White solid; m.p.: 186-188°C.

^1H NMR (400 MHz, CDCl_3): δ 7.58 (d, J = 1.6 Hz, 1H), 7.28 (dd, J = 8.8, 2.3 Hz, 1H), 6.56 (d, J = 8.7 Hz, 1H), 3.94 (d, J = 9.5 Hz, 1H), 3.52 (dd, J = 11.6, 2.5 Hz, 1H), 3.36 (ddd, J = 9.5, 4.5, 2.5 Hz, 1H), 3.05-2.97 (m, 4H), 2.77 (s, 3H).

^{13}C NMR (101 MHz, CDCl_3): δ 178.2, 176.1, 147.4, 132.5, 131.2, 120.5, 114.1, 111.5, 50.2, 43.3, 41.6, 39.4, 25.4.

HRMS (ESI): calcd for $\text{C}_{13}\text{H}_{13}\text{BrN}_2\text{O}_2 + \text{Na}$ = 331.0058, found: 331.0014.



(new compound)

(3aS*,9bR*)-2-(3-Bromophenyl)-8-(4-(dimethylamino)benzyl)-5-methyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3p)

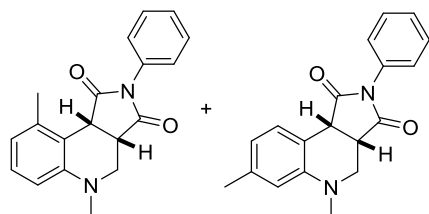
White solid; m.p.: 187-189°C.

^1H NMR (400 MHz, CDCl_3): δ 7.48 (d, J = 8.7 Hz, 2H), 7.36-7.23 (m, 4H), 7.08 (d, J = 8.7 Hz, 2H), 7.03 (d, J = 8.6 Hz, 1H), 6.67 (dd, J = 16.8, 8.5 Hz, 3H), 4.10 (d, J =

9.6 Hz, 1H), 3.84 (s, 2H), 3.56 (dd, $J = 11.4, 2.6$ Hz, 1H), 3.48 (ddd, $J = 9.7, 4.4, 2.5$ Hz, 1H), 3.03 (dd, $J = 11.4, 4.4$ Hz, 1H), 2.90 (s, 6H), 2.78 (s, 3H).

^{13}C NMR (101 MHz, CDCl_3): δ 177.4, 175.4, 146.7, 133.5, 133.2, 131.5, 130.5, 130.1, 129.6, 129.4, 129.0, 125.0, 122.2, 118.3, 113.0, 112.7, 50.9, 43.6, 42.2, 40.8, 40.0, 39.5.

HRMS (ESI): calcd for $\text{C}_{27}\text{H}_{26}\text{BrN}_3\text{O}_2 + \text{Na} = 526.1106$, found: 526.1101.



(3aS*,9bR*)-5,9-Dimethyl-2-phenyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3q1)⁵

(3aS*,9bR*)-5,7-Dimethyl-2-phenyl-3a,4,5,9b-tetrahydro-1H-pyrrolo[3,4-c]quinoline-1,3(2H)-dione (3q2)⁵

^1H NMR (400 MHz, CDCl_3): δ 7.44-7.37 (m, 2.4H), 7.36-7.30 (m, 1.1H), 7.28-7.21 (m, 2.1H), 7.20-7.09 (m, 1.6H), 6.87 (t, $J = 7.3$ Hz, 0.4H), 6.81 (d, $J = 7.5$ Hz, 0.6H), 6.73 (d, $J = 7.5$ Hz, 1.2H), 6.62 (d, $J = 8.1$ Hz, 0.7H), 6.55 (s, 0.3H), 4.49 (d, $J = 9.8$ Hz, 0.7H), 4.08 (d, $J = 9.6$ Hz, 0.4H), 3.60-3.53 (m, 1H), 3.52-3.44 (m, 1H), 3.05 (dd, $J = 11.5, 4.4$ Hz, 0.3H), 2.91 (dd, $J = 11.4, 4.8$ Hz, 0.7H), 2.80 (s, 1H), 2.76 (s, 2H), 2.57 (s, 2H), 2.31 (s, 1H).

^{13}C NMR (101 MHz, CDCl_3): δ 178.7, 178.0, 176.1, 175.7, 155.8, 150.1, 148.3, 138.5, 138.4, 131.9, 131.8, 130.0, 129.4, 128.9, 128.9, 128.5, 128.4, 127.9, 126.4, 126.3, 122.6, 120.4, 120.2, 119.4, 115.5, 115.2, 113.2, 110.6, 52.5, 50.6, 44.7, 43.4, 41.8, 39.8, 39.4, 39.3, 21.6, 20.3.

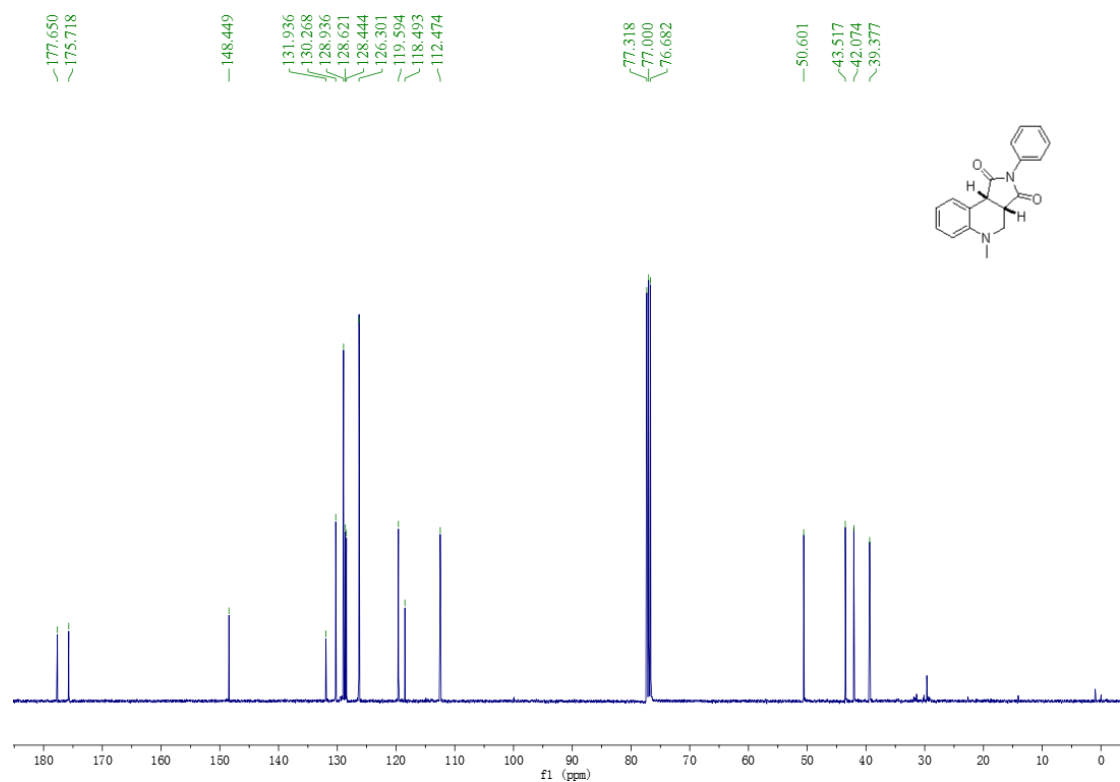
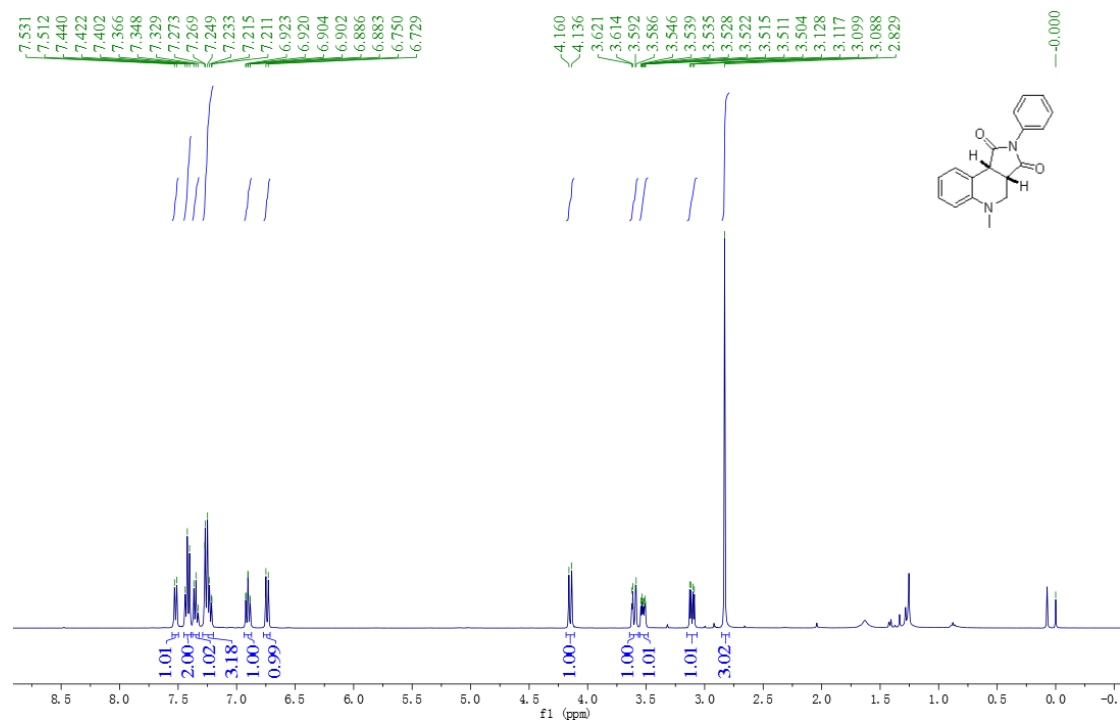
HRMS (ESI): calcd for $\text{C}_{19}\text{H}_{18}\text{N}_2\text{O}_2 + \text{Na} = 329.1266$, found: 329.1257.

References:

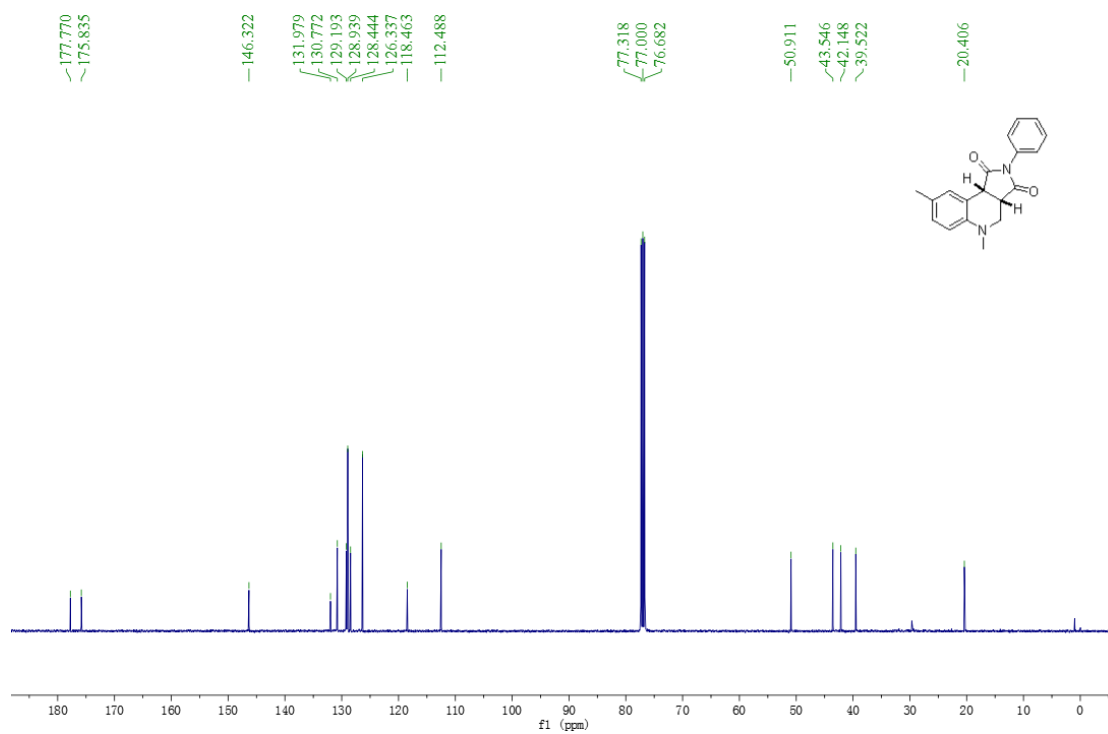
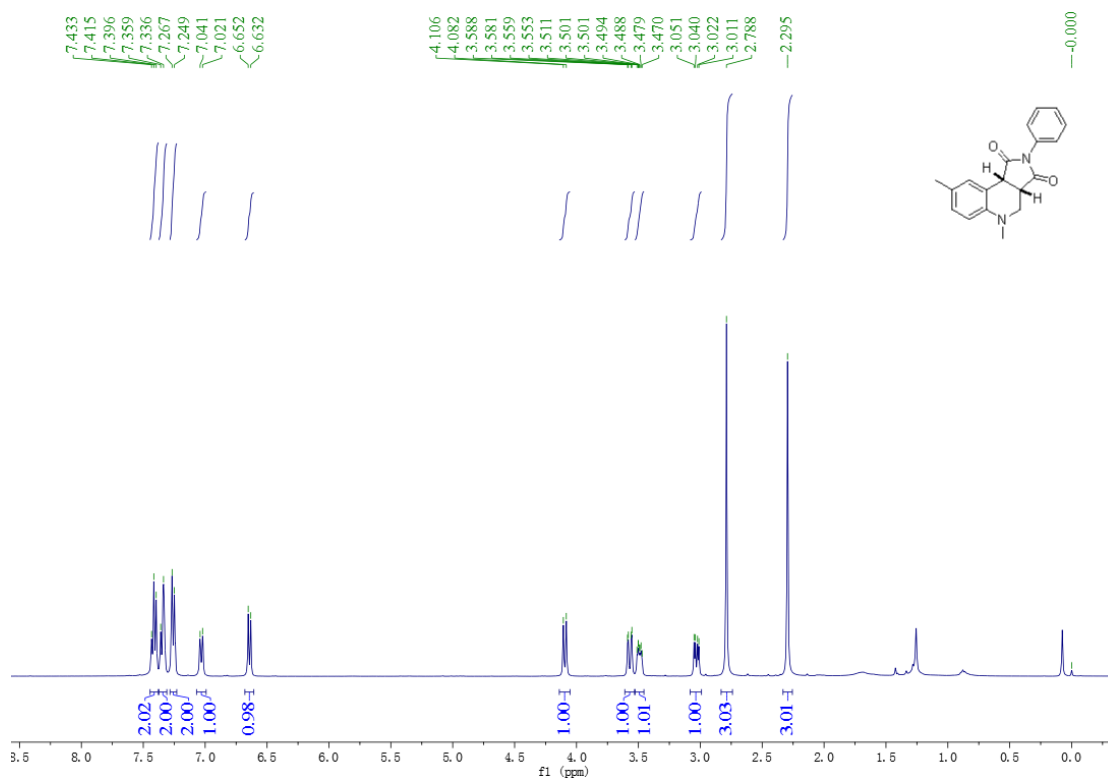
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^1H and ^{13}C NMR spectra of products

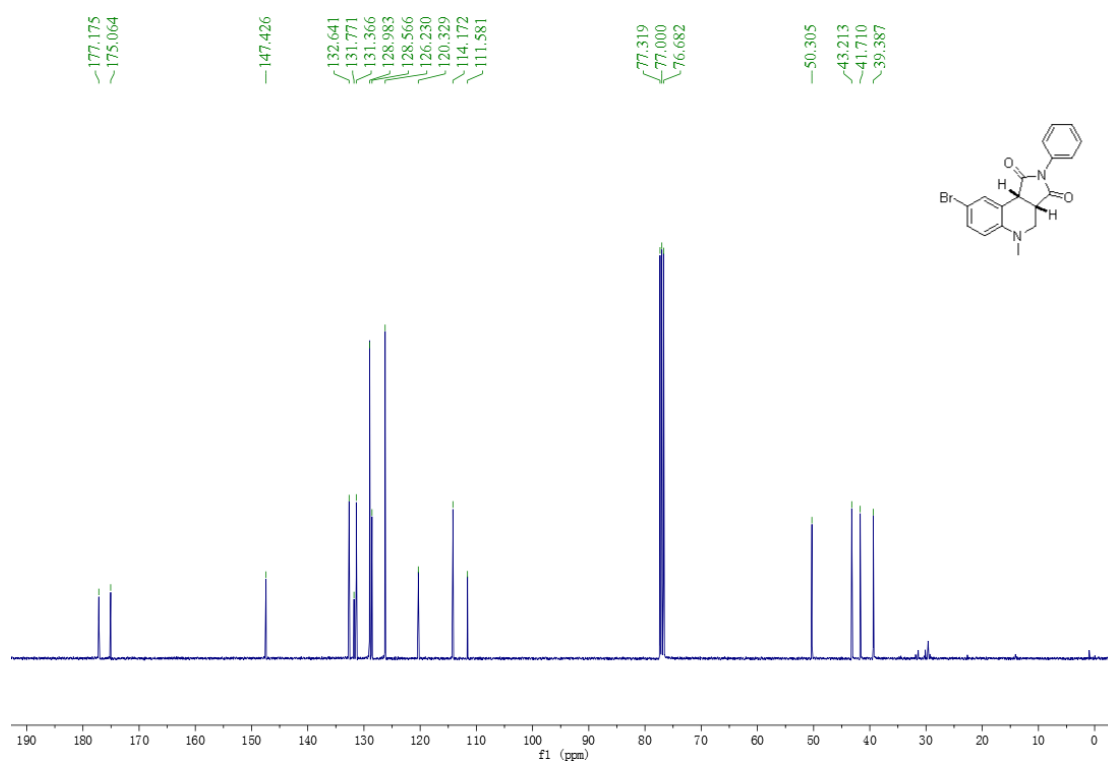
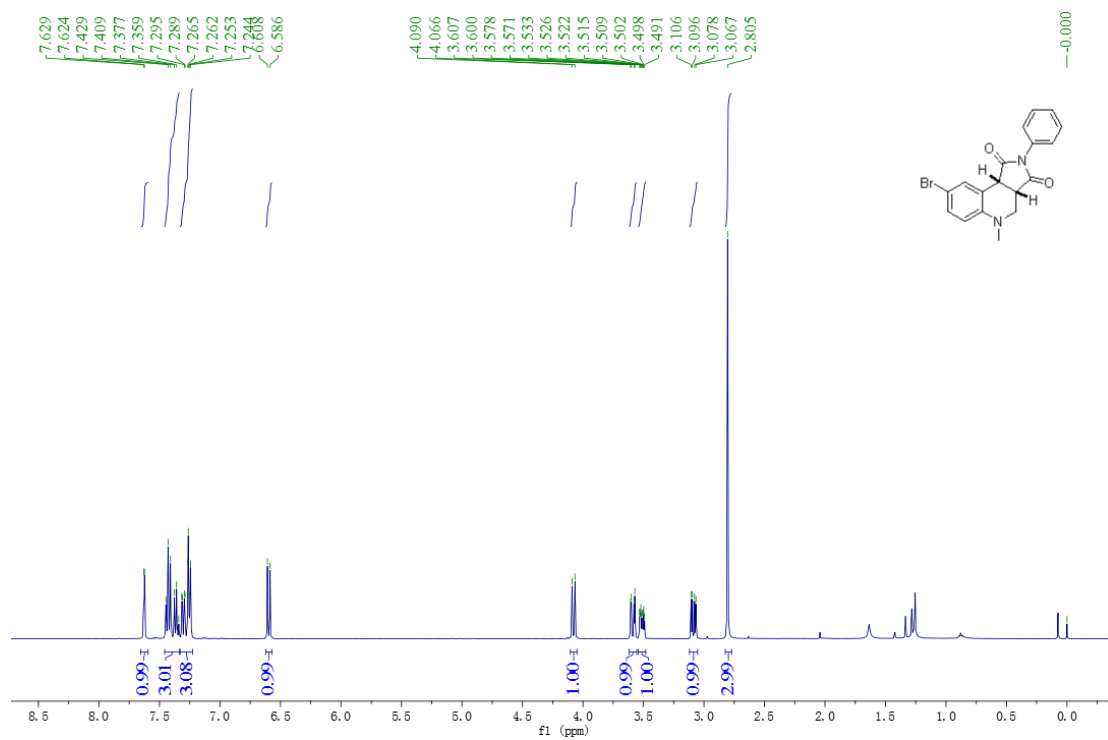
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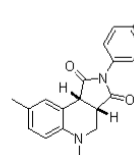
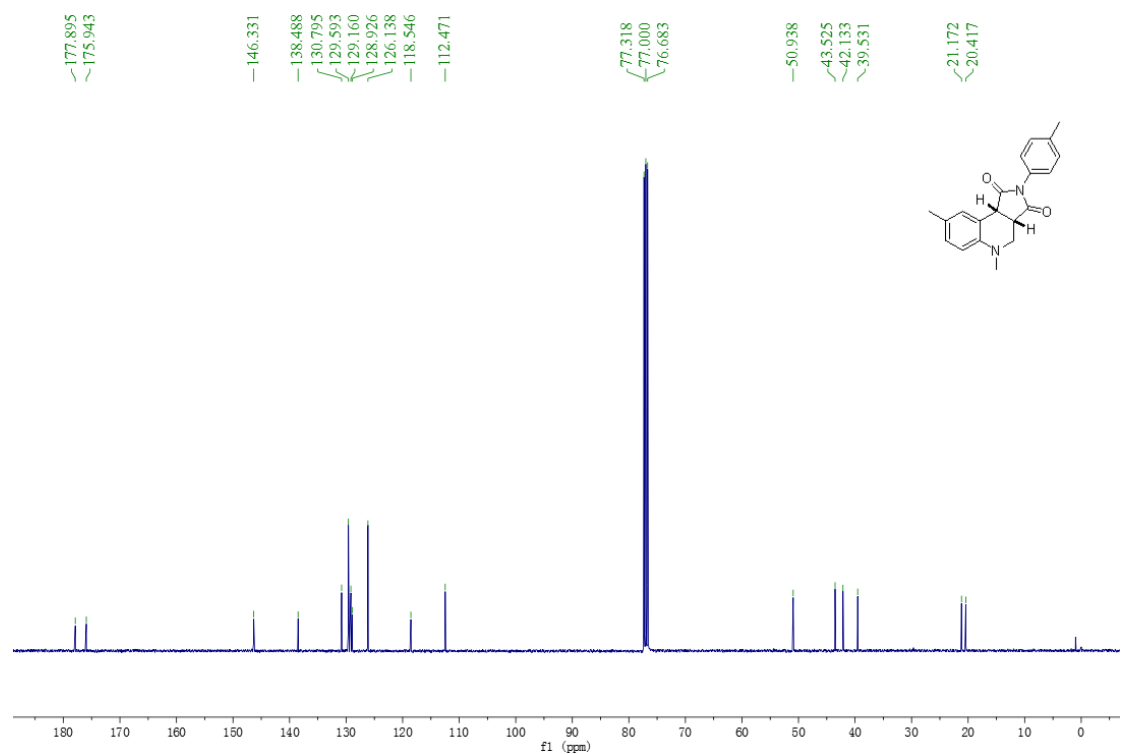
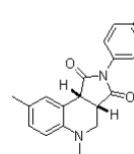
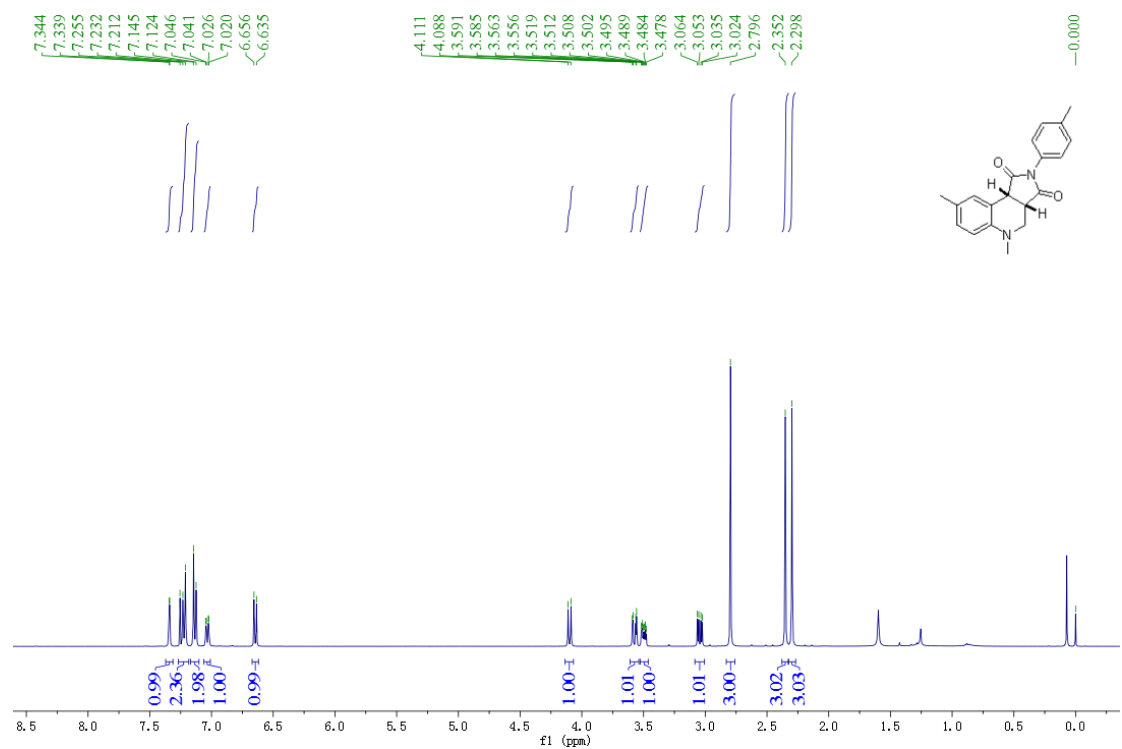
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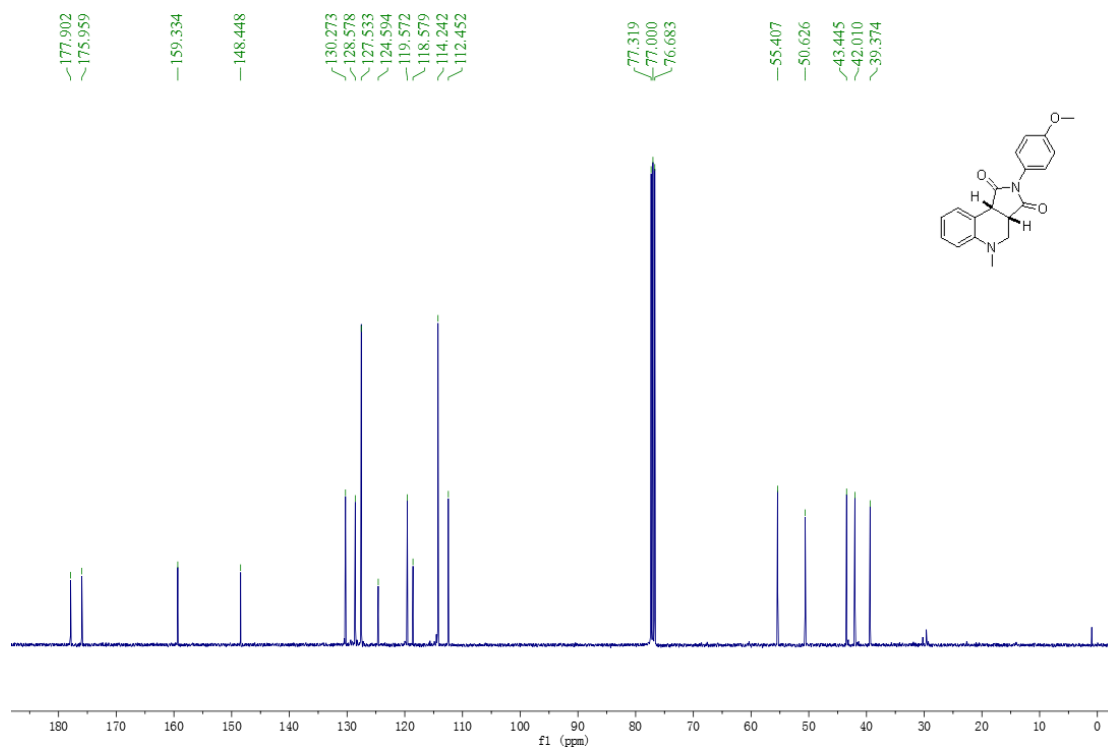
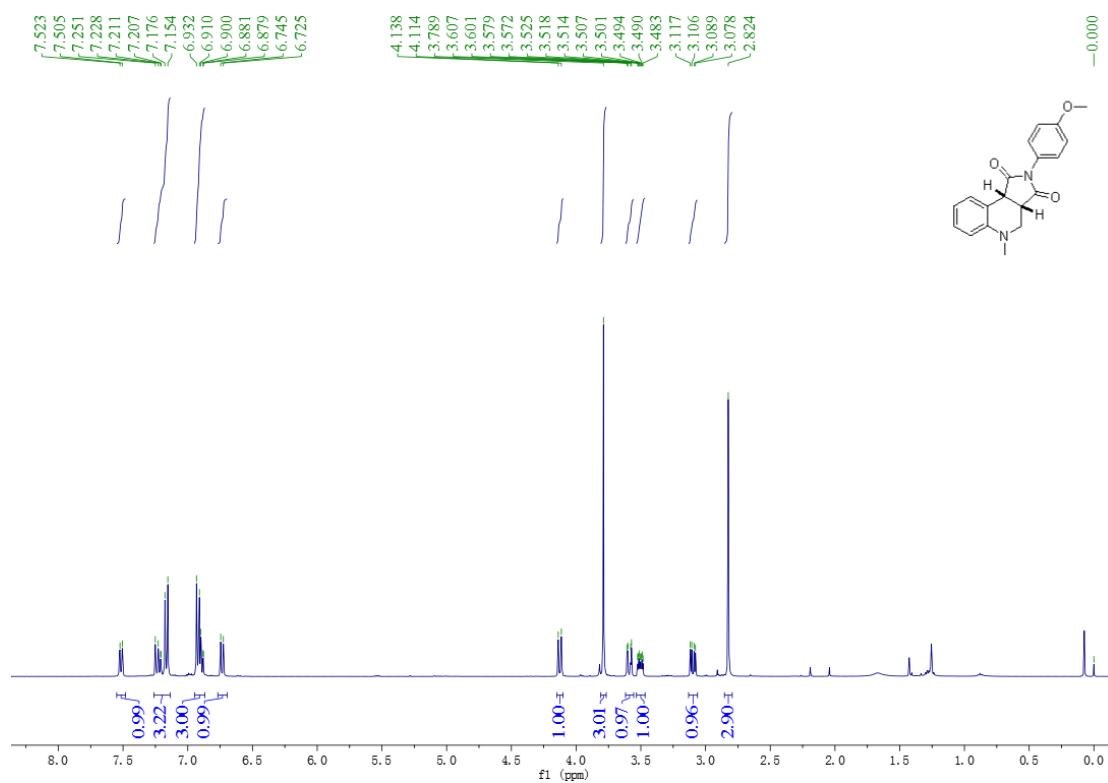
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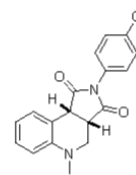
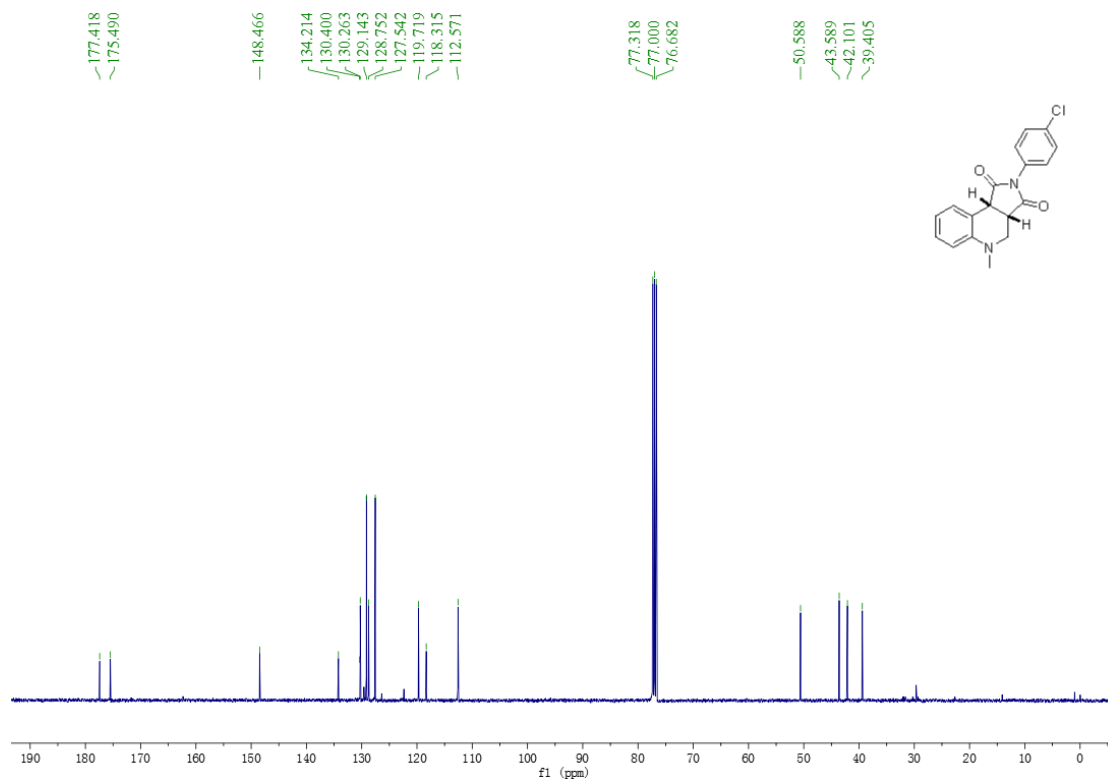
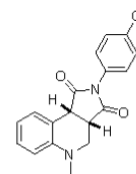
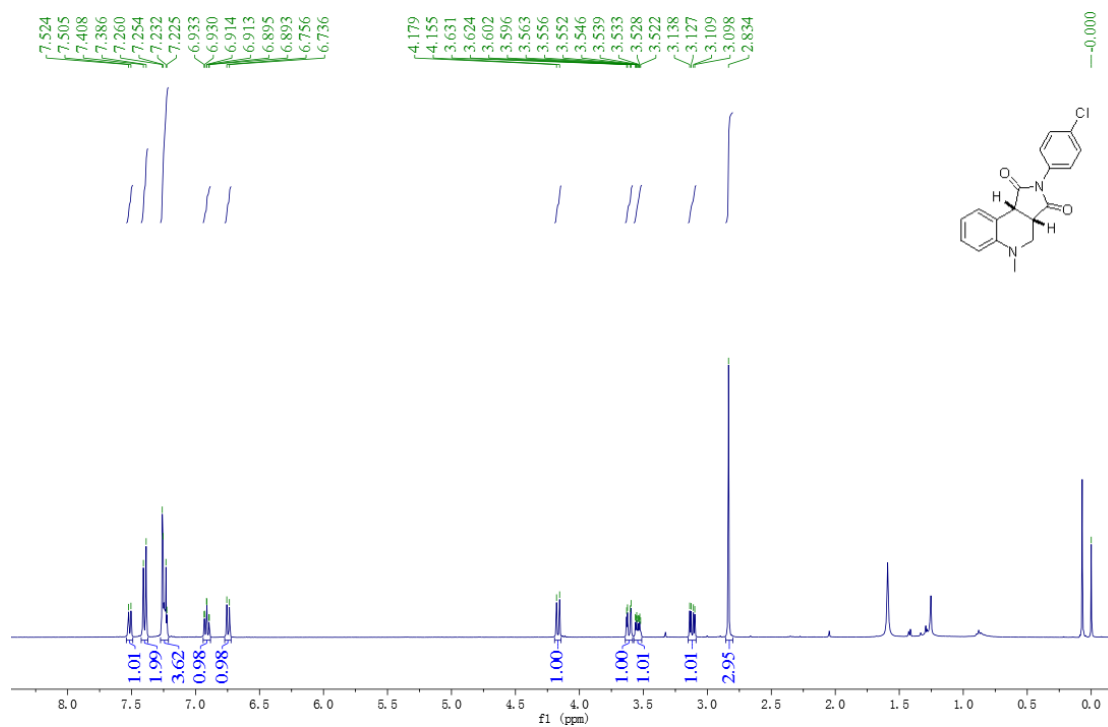
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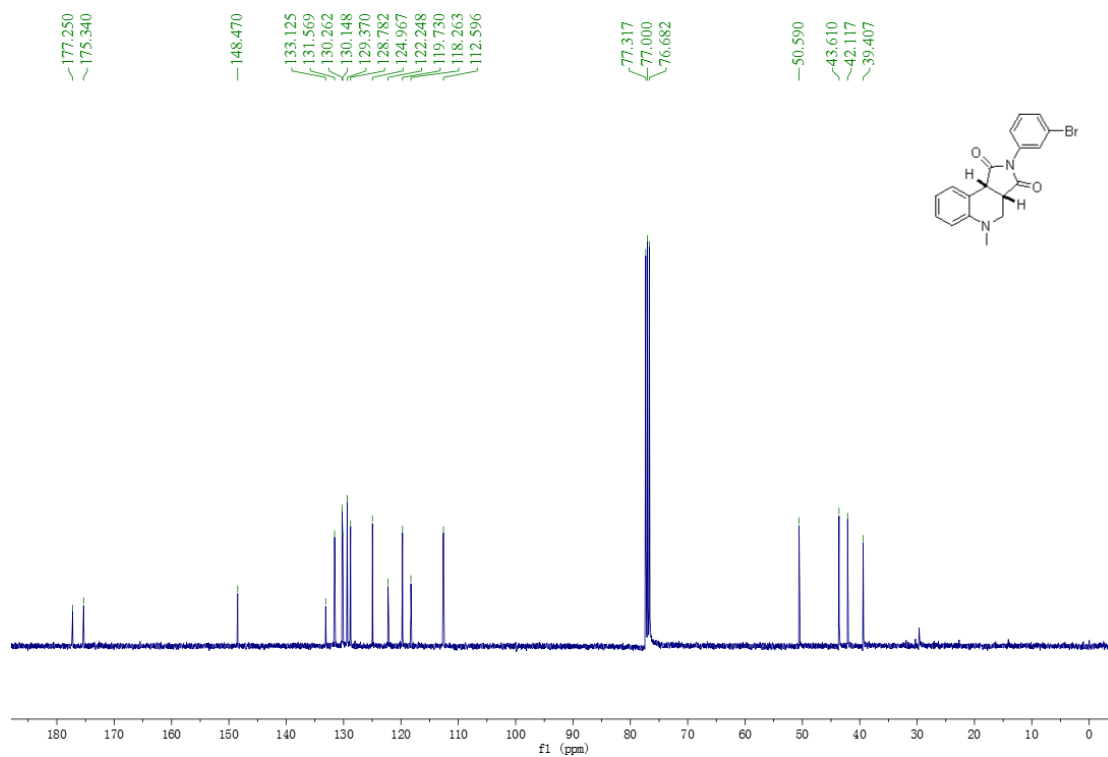
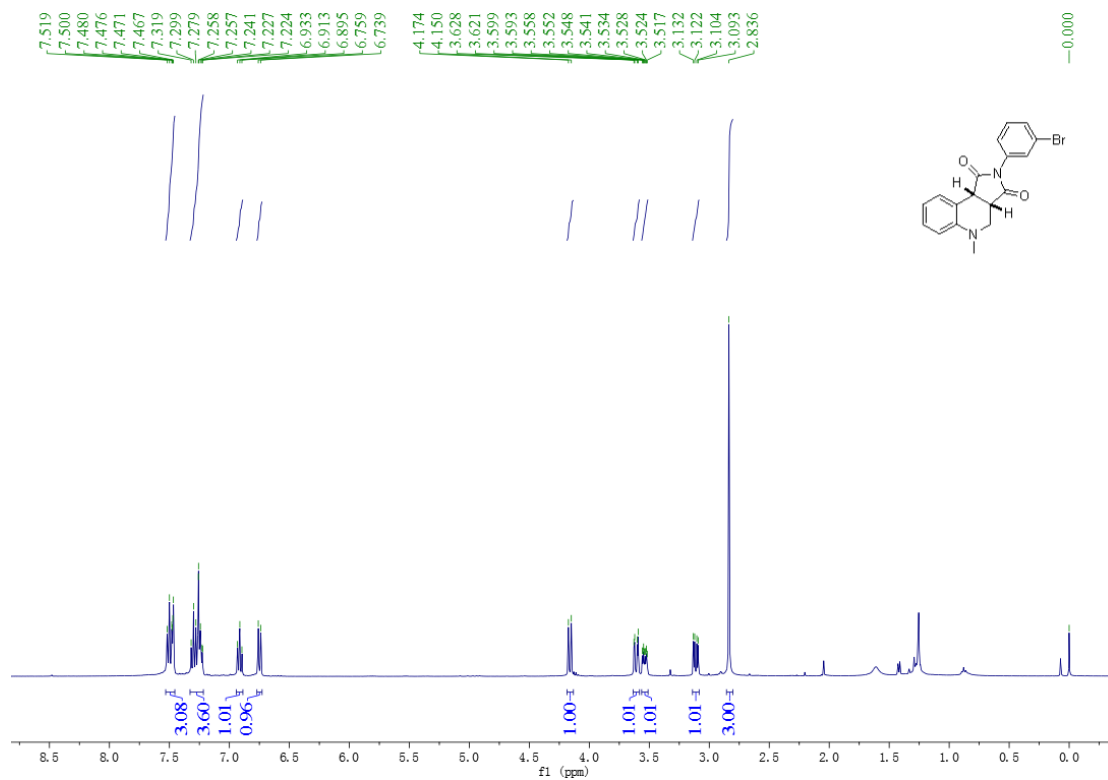
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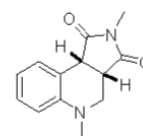
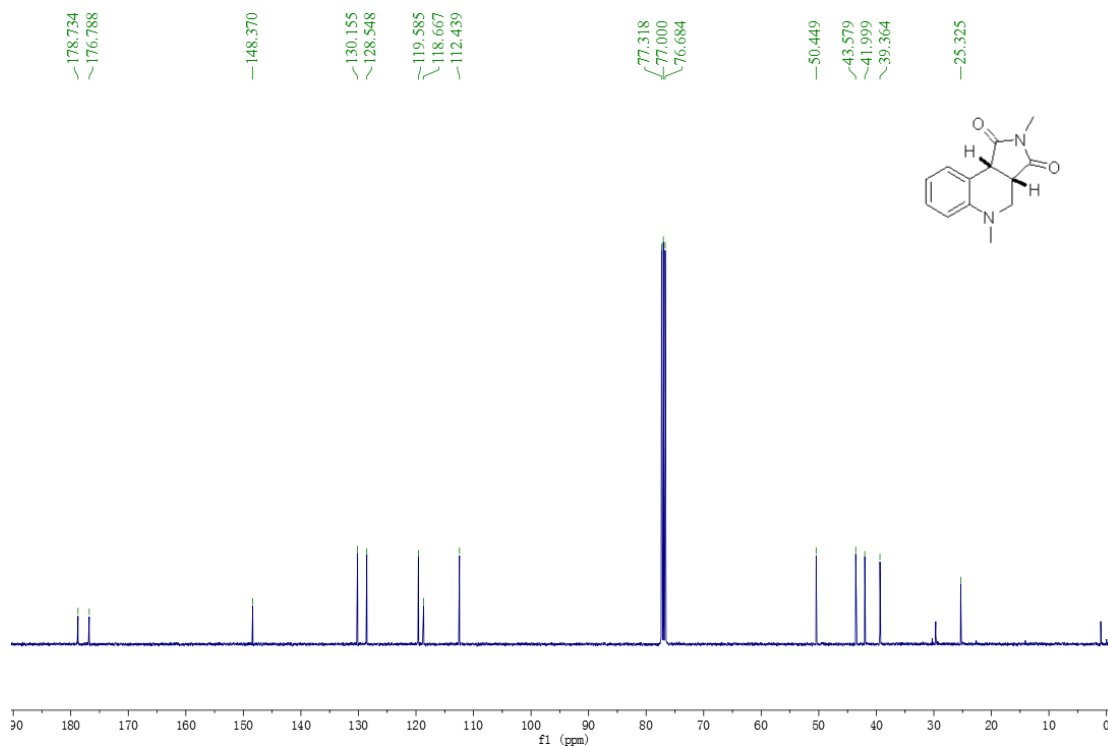
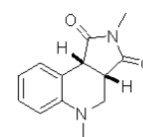
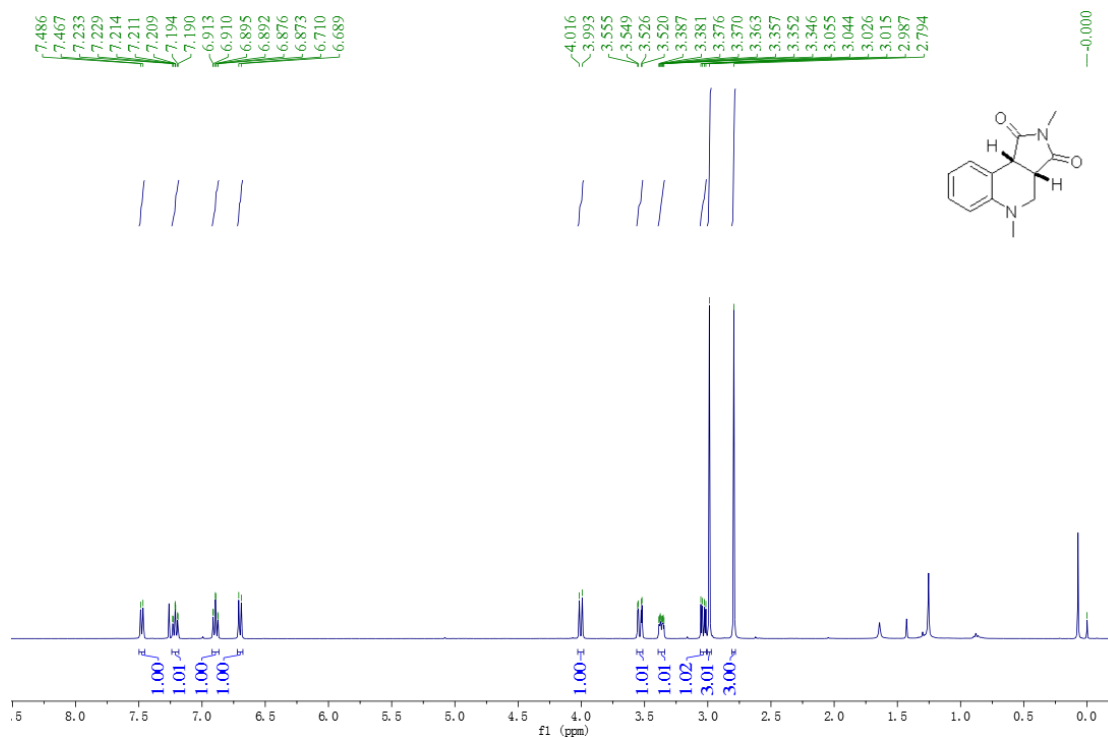
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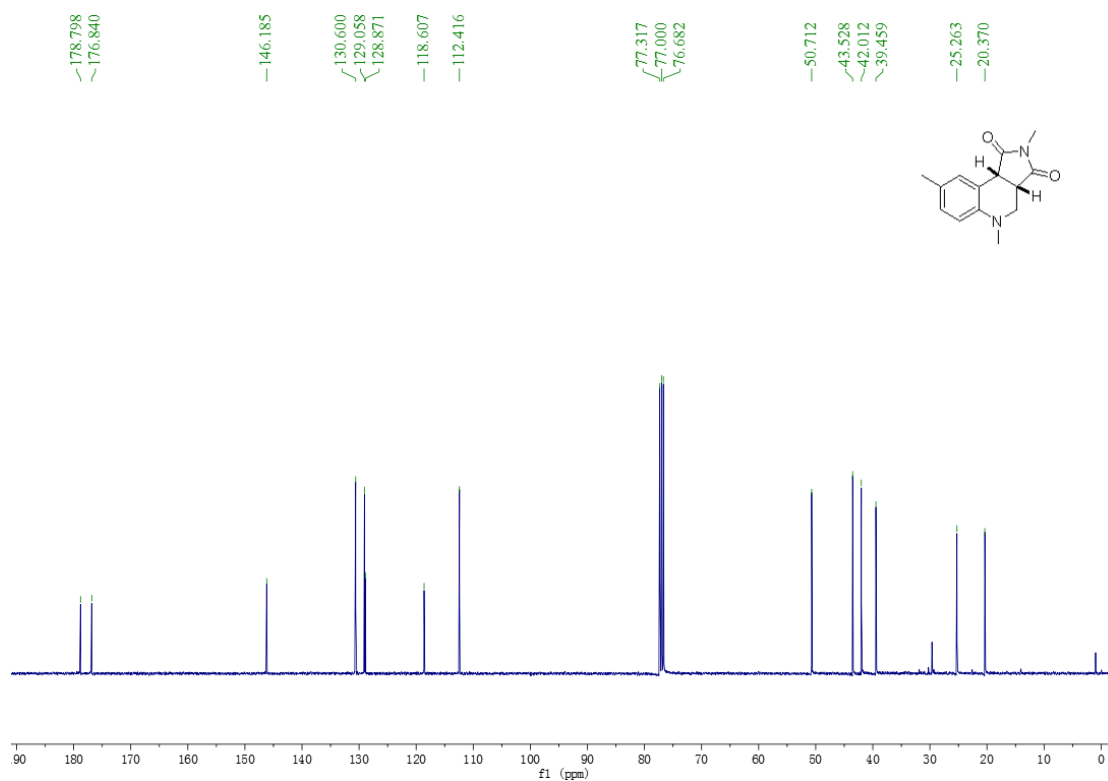
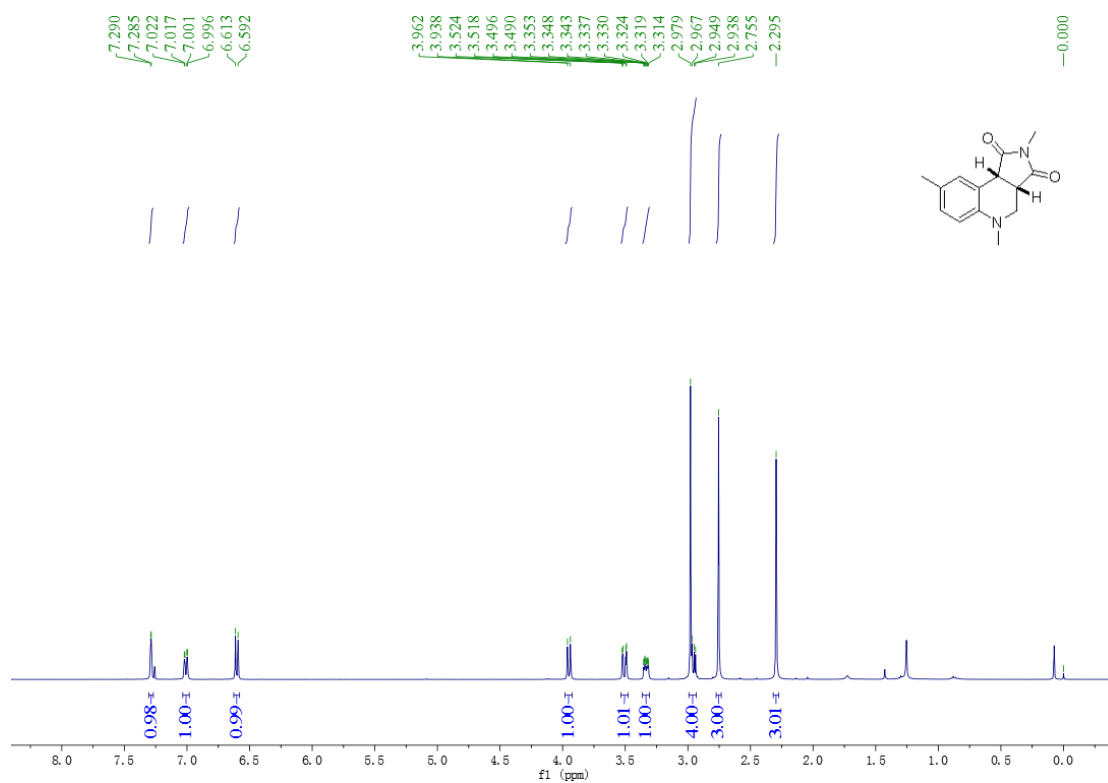
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3h



3i

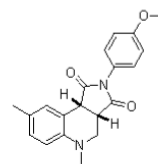
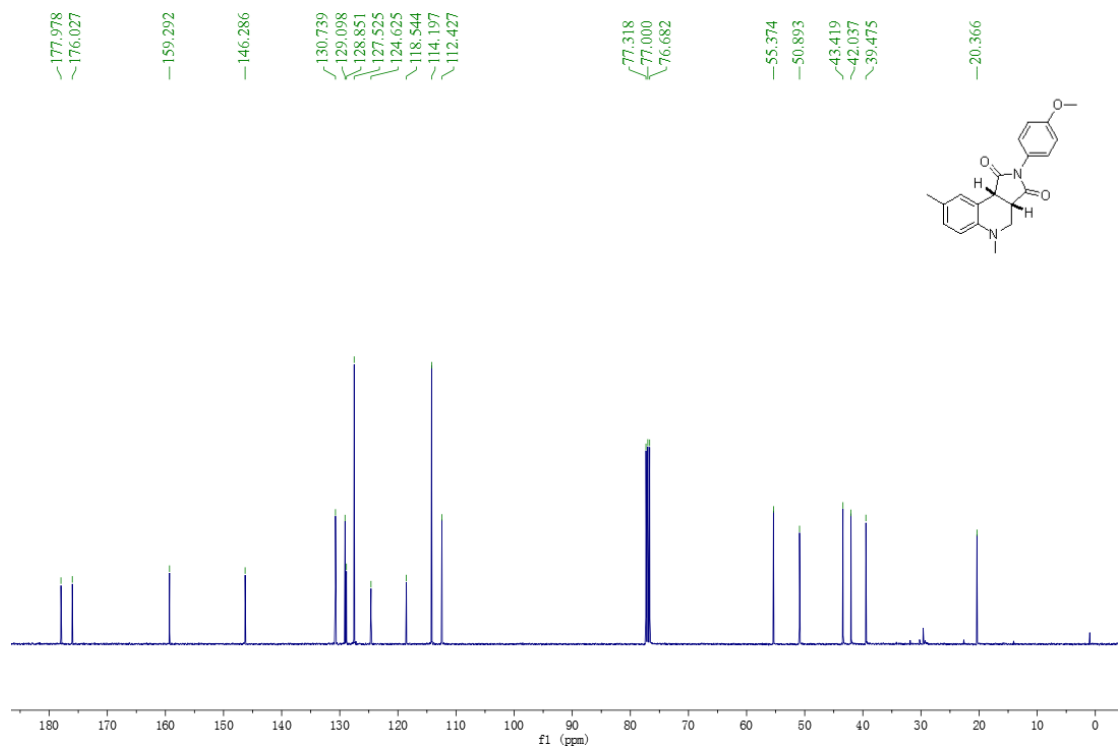
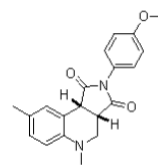


Chemical Structure of Compound 10:

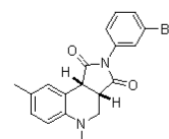
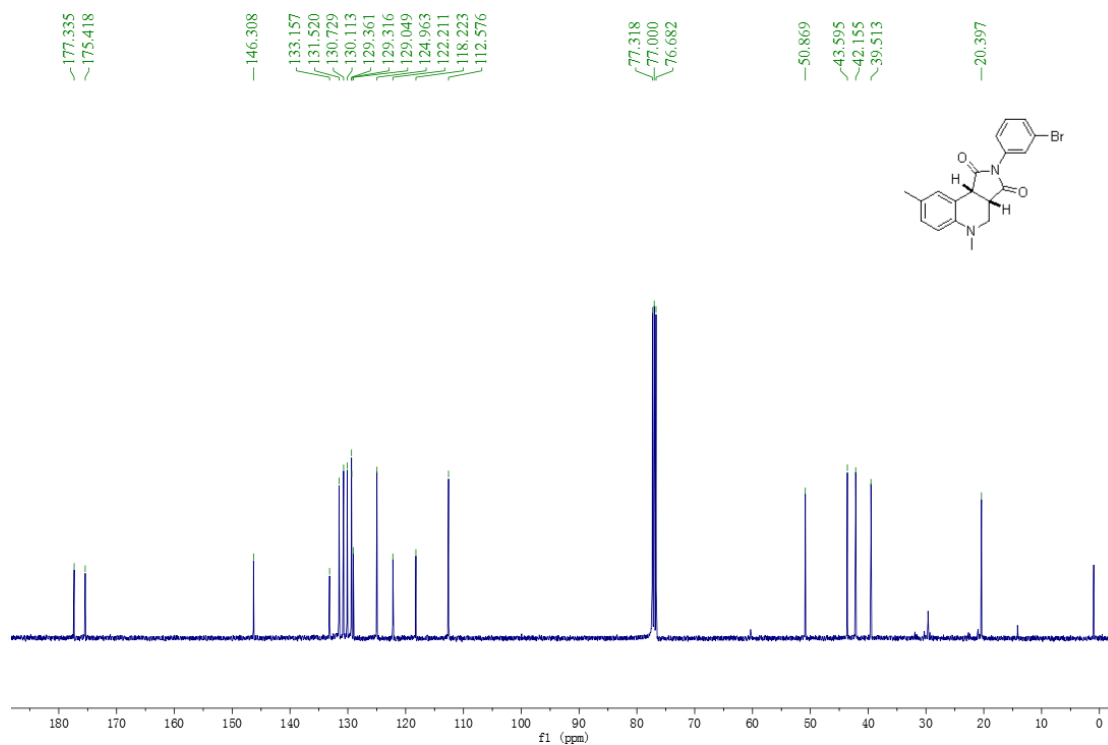
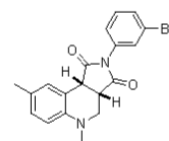
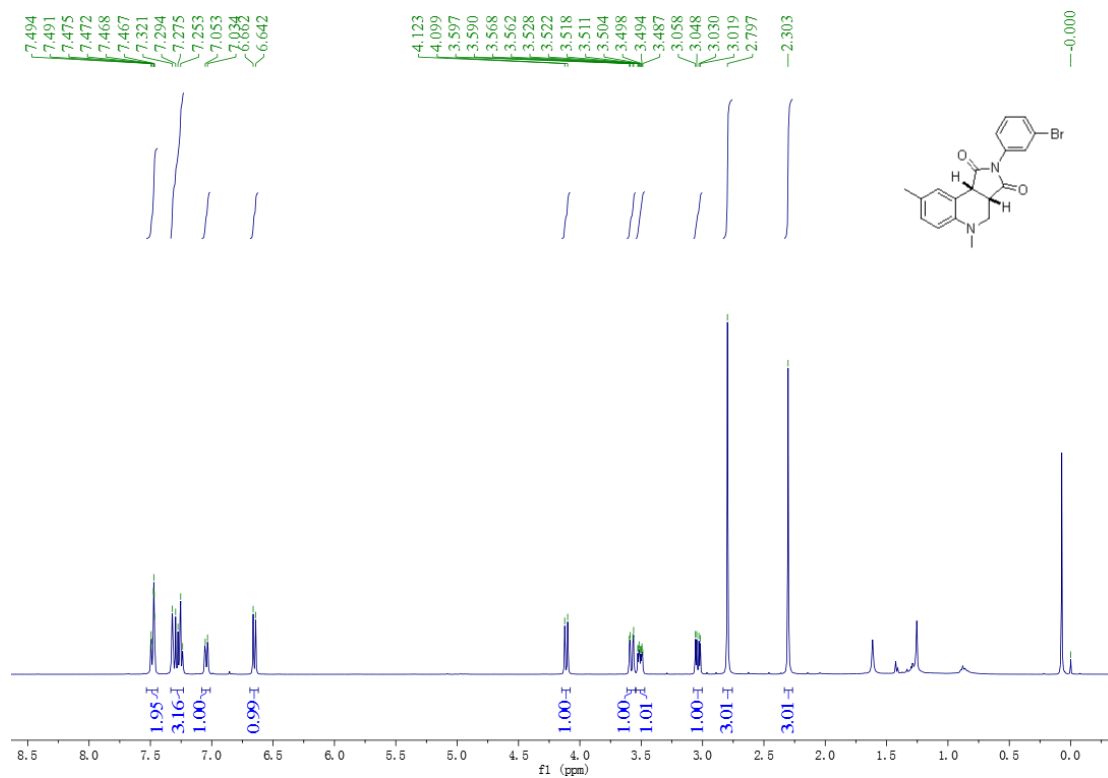
Cc1ccc2c(c1)nc3c2c(=O)n(c3=O)C(=O)c4ccc(OC)cc4

¹H NMR Spectrum (CDCl₃):

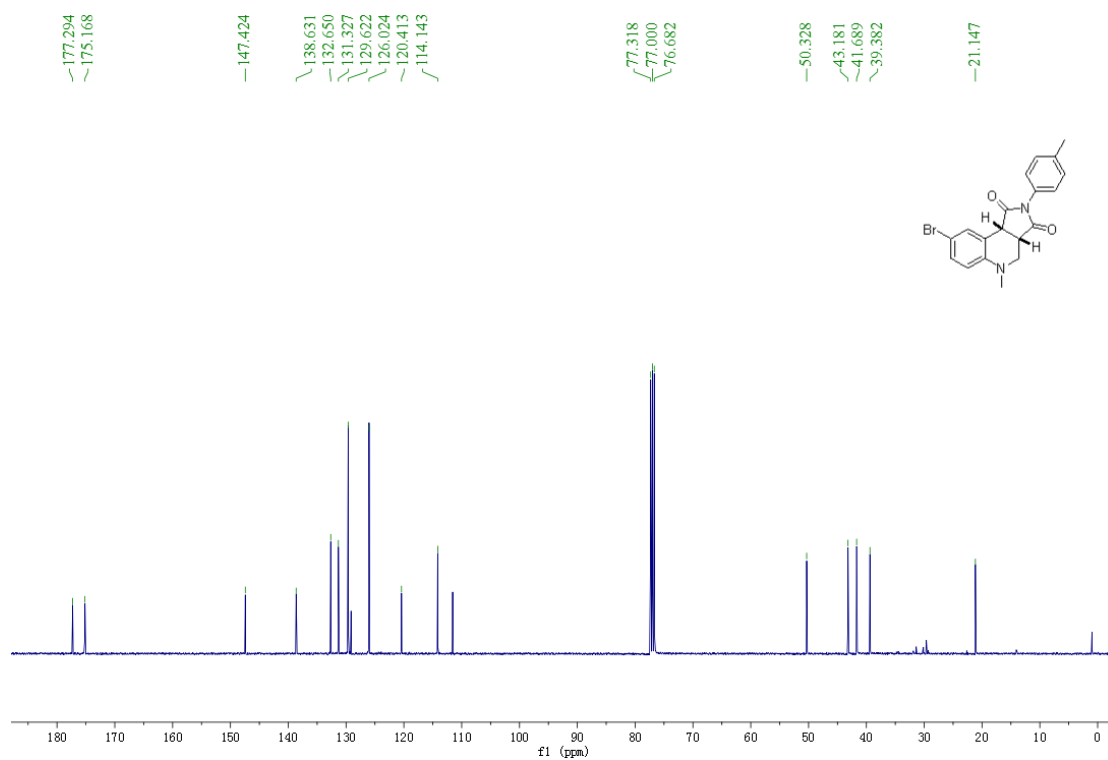
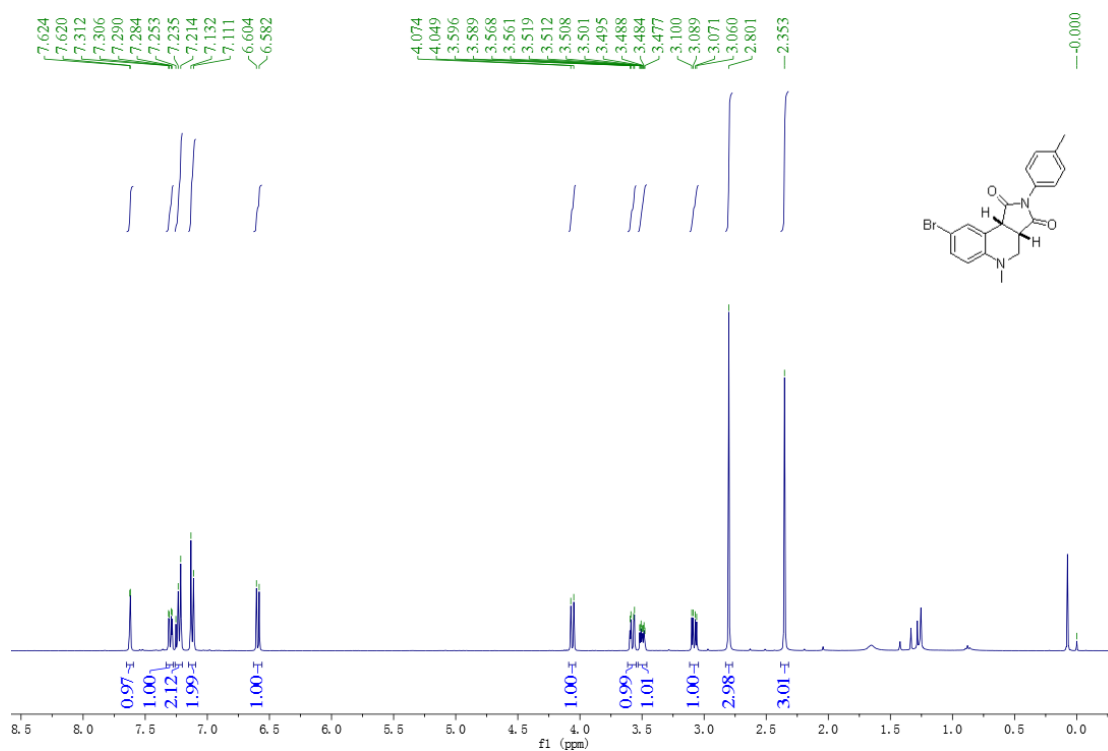
Chemical Shift (ppm)	Integration
7.331, 7.326, 7.175, 7.153, 7.038, 7.017, 6.928, 6.905, 6.648, 6.628	0.99, 2.00, 1.00, 2.00, 1.00
4.084, 4.060, 3.784, 3.575, 3.569, 3.547, 3.540, 3.492, 3.485, 3.481, 3.474, 3.468, 3.461, 3.457, 3.450	1.00, 3.00, 1.00, 1.00
3.042, 3.031, 3.013, 3.002, 2.785	0.99, 2.99
2.293	3.00
0.000	-



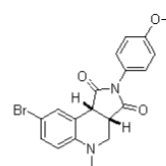
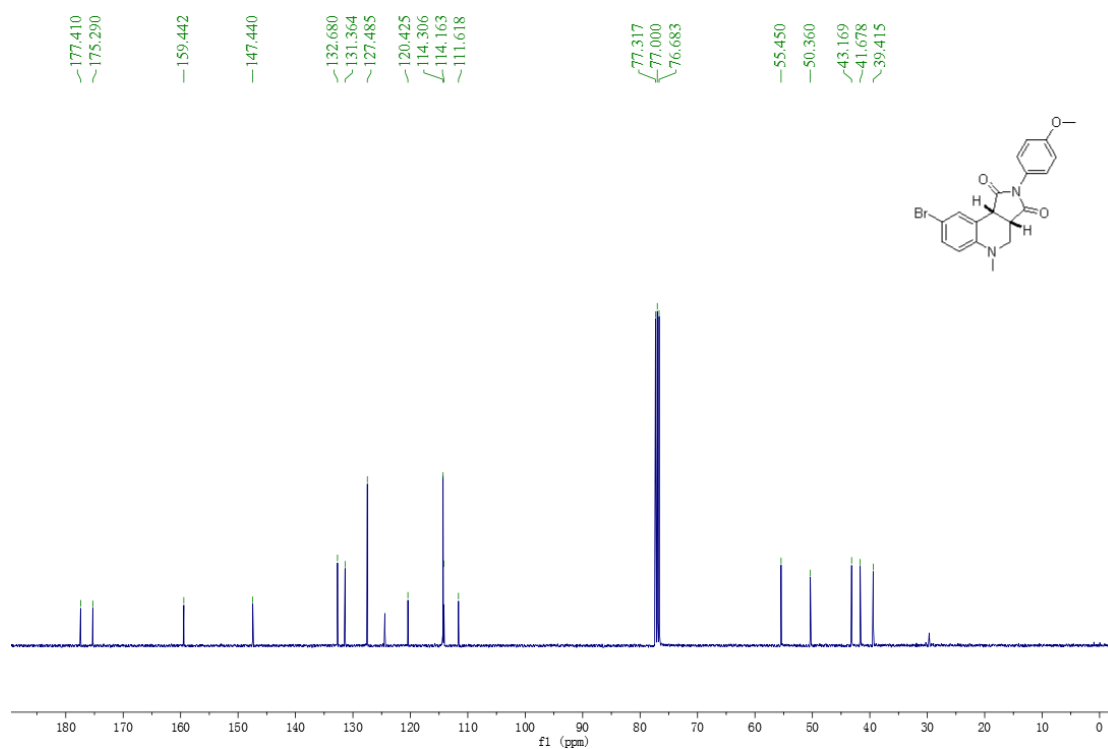
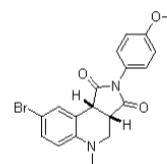
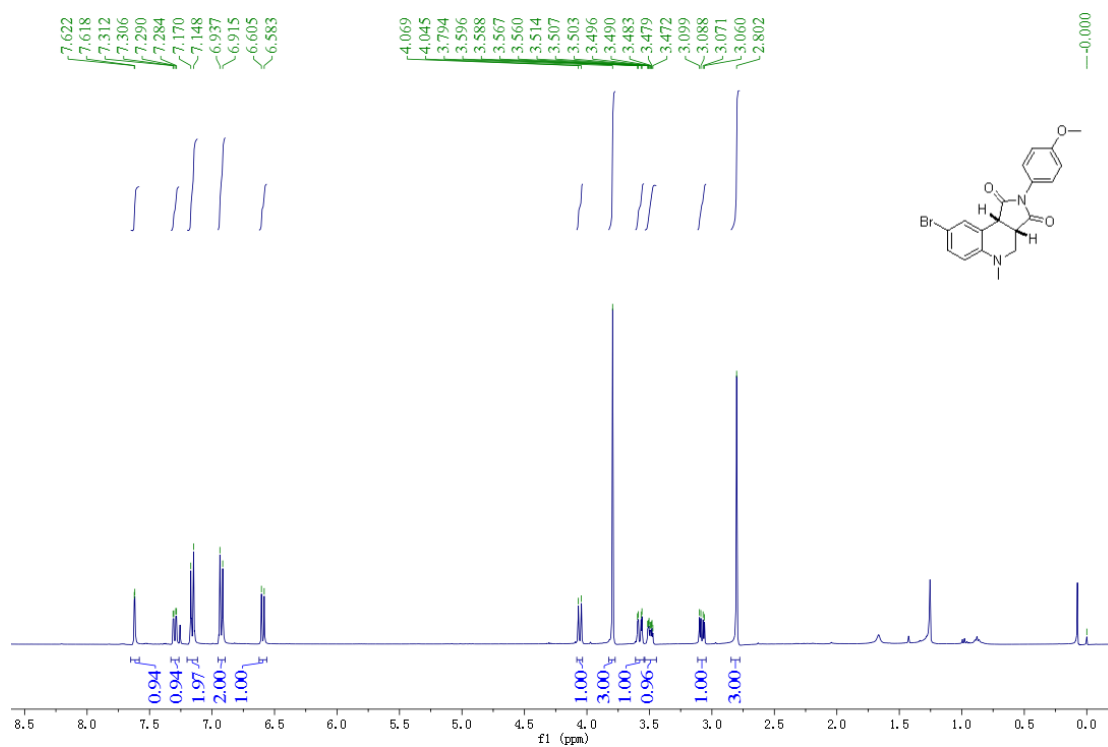
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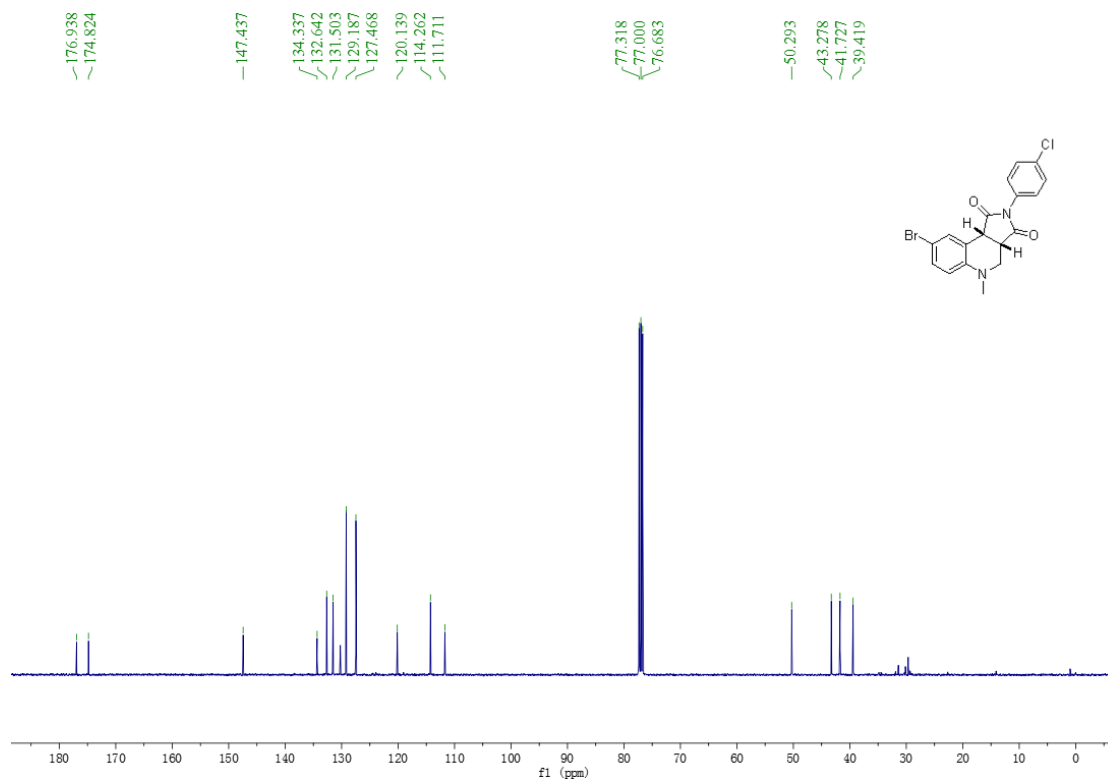
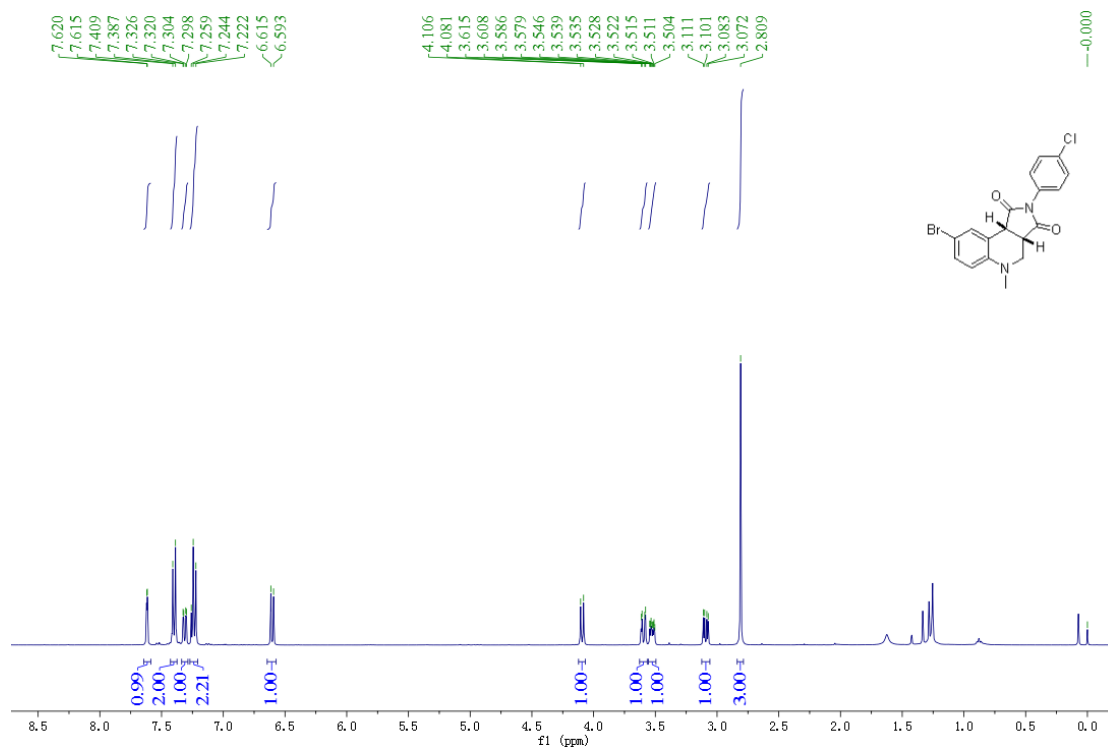
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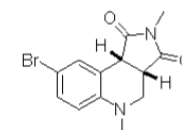
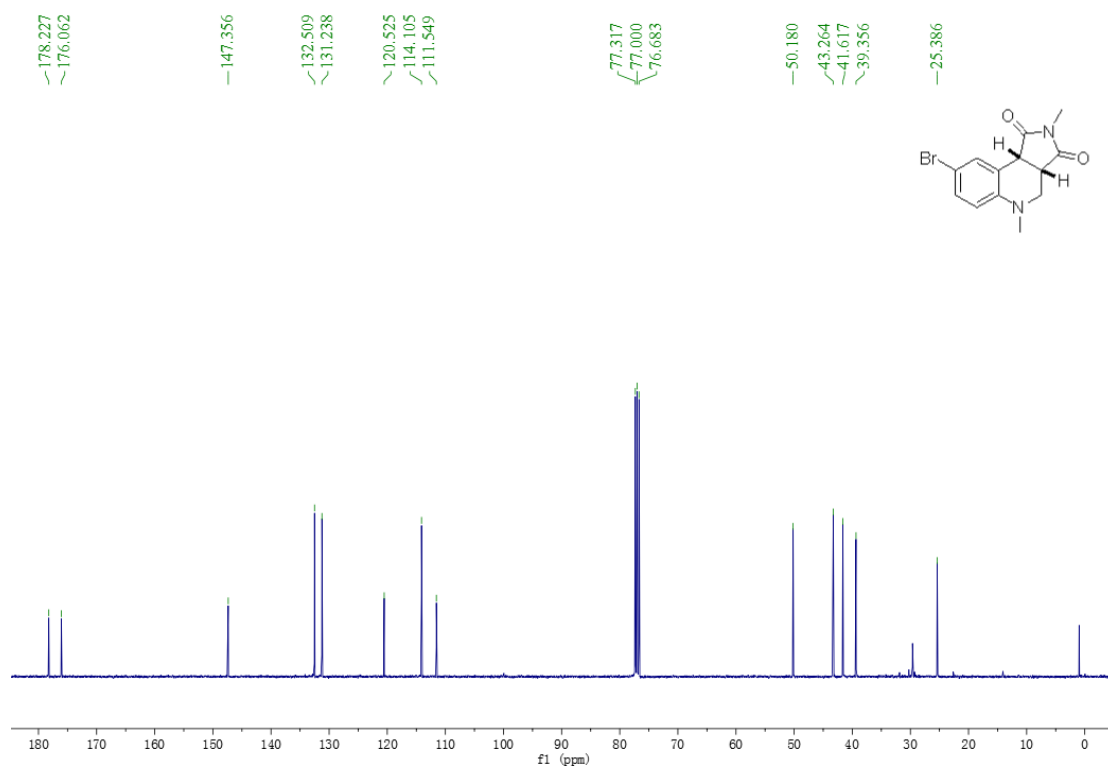
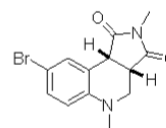
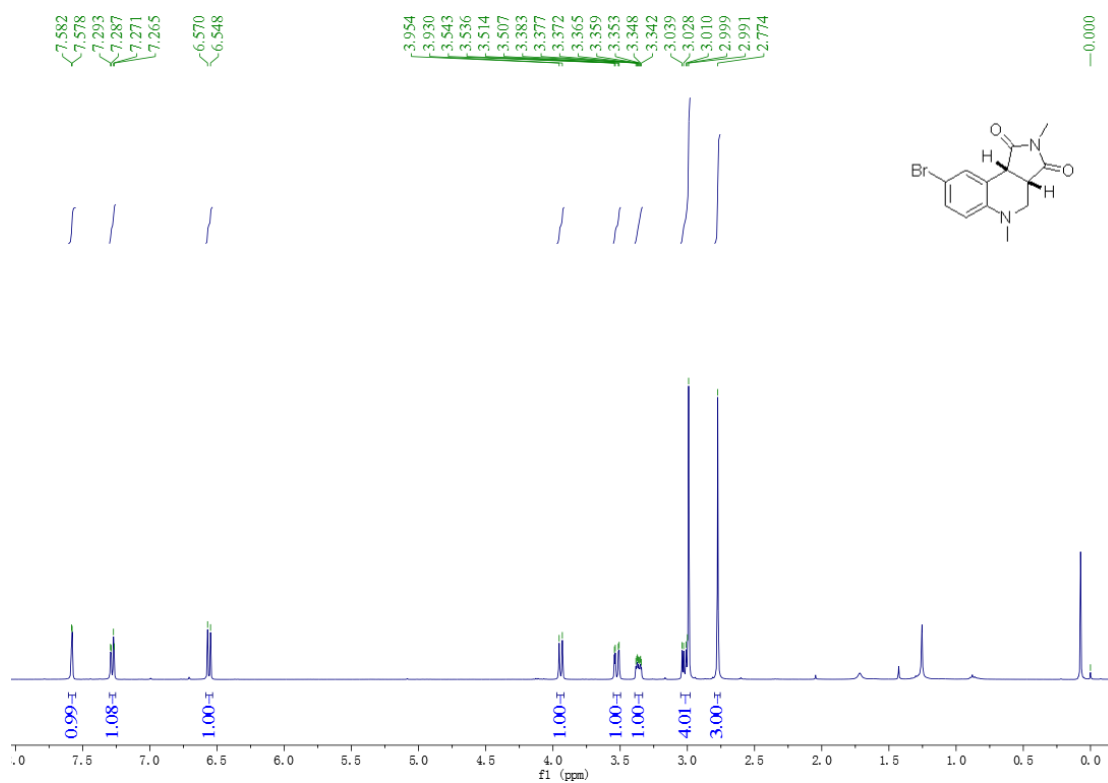
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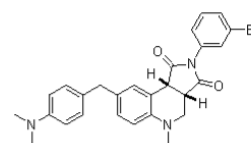
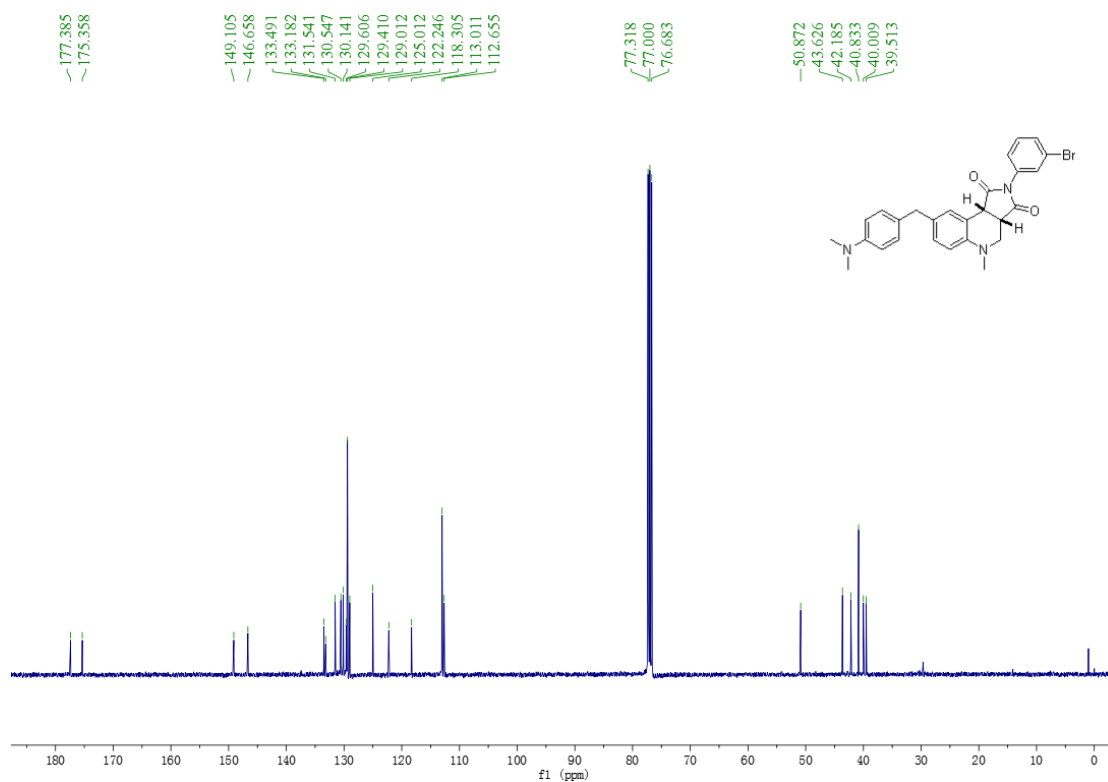
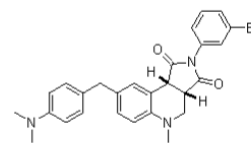
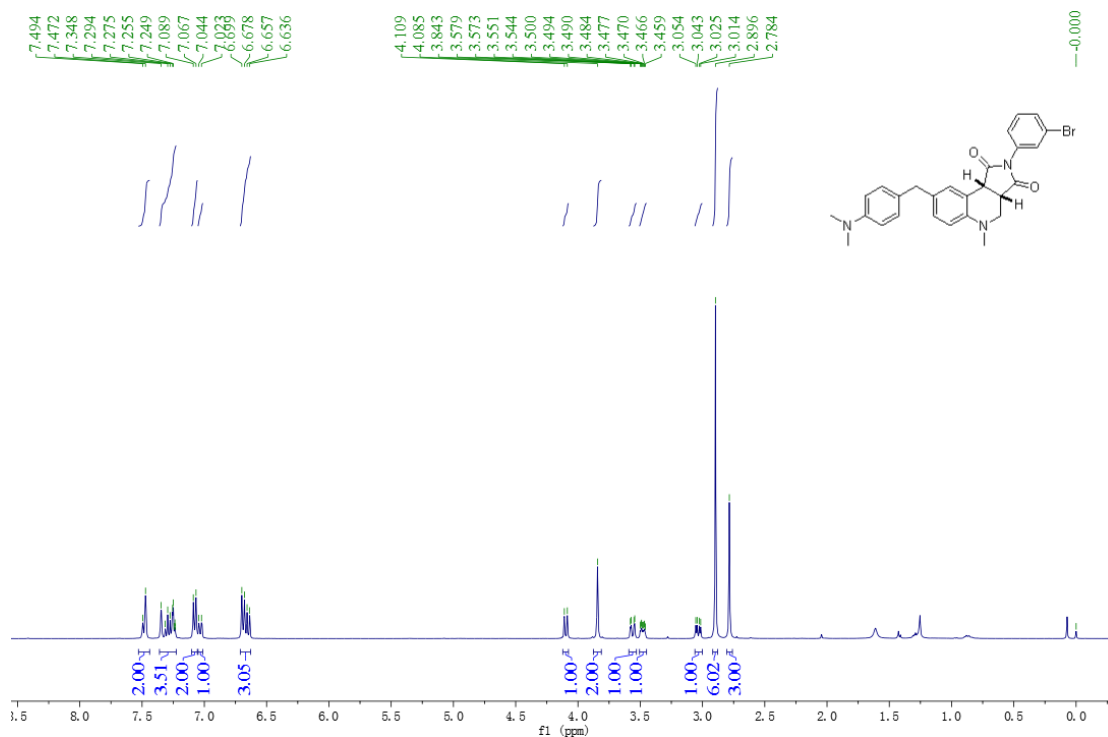
3n



3o



3p



3q1+3q2

