

Regioselective Bromination of Pyrrolo[1,2-a]quinoxalines

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Supporting Information

1. Materials and instruments	2
2. Experimental Section.....	2
3. The NMR Spectra Data	3
4. Copies of Spectra	13

1. Materials and instruments

Unless otherwise noted, all commercial materials were used directly without further purification, and all reactions were performed in the air. For chromatography, Qingdao Ocean Chemical 200-300 mesh silica gel was employed. Melting points were determined with a fusiometer. ^1H NMR and ^{13}C NMR spectra were recorded on Bruker Avance III HD 400 MHz and Bruker AscendTM 600 MHZ spectrometer in CDCl_3 and the chemical shifts are reported in ppm (δ) relative to the internal standard tetramethylsilane (TMS) (0 ppm). High-resolution mass spectra (HRMS) were acquired in atmospheric pressure chemical ionization (APCI) mode using a TOF mass analyzer.

2. Experimental Section

General procedure for the C3-bromination of pyrrolo[1,2-*a*]quinoxaline with TBATB

A 10 mL Schlenk tube was charged with pyrrolo[1,2-*a*]quinoxaline 1 (0.2 mmol), TBATB (0.24 mmol) and MeCN (2.5 mL). The mixture was stirred at 80 °C for 12 hours. After completion of the reaction, the solution was quenched with a saturated solution of sodium thiosulfate (10 mL) and extracted with ethyl acetate (15 mL × 3). The organic layer was dried with anhydrous Na_2SO_4 and the solvent was removed under reduced pressure, and the crude product was purified by rapid chromatography on silica gel (petroleum ether/ethyl acetate = 8:1) to give the final products **3a-3w**.

General procedure for the dibromination of pyrrolo[1,2-*a*]quinoxaline with TBATB

A 10 mL Schlenk tube was charged with pyrrolo[1,2-*a*]quinoxaline 1 (0.2 mmol), TBATB (0.24 mmol) and DMSO (2.5 mL). The mixture was stirred at 80°C for 12 hours. After completion of the reaction, the solution was quenched with a saturated solution of sodium thiosulfate (10 mL) and extracted with ethyl acetate (15 mL × 3). The organic layer was dried with anhydrous Na_2SO_4 and the solvent was removed under reduced pressure. The residue was purified from the crude product by rapid chromatography on silica gel (petroleum ether/ethyl acetate = 10:1) to give the final

products **5a-5f**.

Gram-scale synthesis procedure for 3a

To a 200 mL round-bottomed flask was added 1a (6 mmol), TBATB (7.2 mmol) and MeCN (50 mL). The solution was stirred at 100 °C for 12 hours. After completion of the reaction, the solution was quenched with a saturated solution of sodium thiosulfate (25 mL) and extracted with dichloromethane (30 mL × 3). The organic layer was dried with anhydrous Na₂SO₄ and the solvent was removed under reduced pressure. The residue was purified by rapid chromatography on silica gel (petroleum ether/ethyl acetate = 8:1) to afford the final product **3a** (1.401 g, 95% yield).

Palladium-catalyzed Suzuki-Miyaura reaction of 3a with phenylboronic acid

To a 10 mL Schlenk tube was added **3a** (0.2 mmol) phenylboronic acid added to acid (0.4 mmol), Pd(PPh₃)₄ (5 mol%), Na₂CO₃ (2 equiv.) and 1,4-dioxane/H₂O (4:1, 2.5 mL). The reaction vessel was stirred at 80 °C for 5 h under air atmosphere. Upon completion of the reaction, saturated aqueous NaCl solution (10 mL) was added to the reaction solution, followed by extraction with dichloromethane (15 mL × 3). The organic phase was dried with Na₂SO₄ and the solvent was removed under reduced pressure. The residue was purified by rapid chromatography on silica gel (petroleum ether/ethyl acetate = 5:1) to give the final product **3aa**.

Palladium-catalyzed C1-arylation of 3a with 4-iodotoluene

A 10 mL Schlenk tube was filled with **3a** (0.25 mmol), 4-iodotoluene (0.5 mmol), Pd(OAc)₂ (10 mol%), X-Phos (15 mol%) and toluene (1 mL). The mixture was stirred in air at 120 °C for 24 hours. The reaction solution was concentrated under reduced pressure and the residue was purified by silica gel column chromatography (petroleum ether/ethyl acetate = 30:1) to give the final product **3ab**.

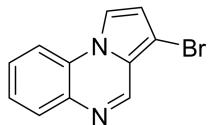
C1-Chlorination of 3a

A 10 mL Schlenk tube was charged with **3a** (0.2 mmol), 1-chloro-1,2-phenyliodonio-3-one (0.24 mmol) and MeCN (2 mL). The mixture was then stirred at room temperature for 12 hours. After completion of the reaction, the solution was quenched with a saturated solution of sodium bicarbonate (10 mL) and extracted with dichloromethane (15 mL × 3). The organic layer was dried with

anhydrous Na₂SO₄ and the solvent was removed under reduced pressure, and the residue was purified by rapid chromatography on silica gel (petroleum ether/ethyl acetate = 10:1) to give the final product **3ac**.

3. The NMR Spectra Data

3-bromopyrrolo[1,2-*a*]quinoxaline [3a]



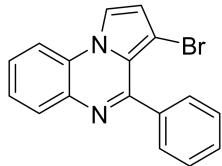
Mp. 166-171 °C. White solid (46.4 mg, 94%).

¹H NMR (400 MHz, CDCl₃) δ 8.81 (s, 1H), 8.01 (dd, *J* = 8.0, 1.6 Hz, 1H), 7.88 (d, *J* = 2.8 Hz, 1H), 7.84 (dd, *J* = 8.0, 1.6 Hz, 1H), 7.56 (td, *J* = 7.2, 1.2 Hz, 1H), 7.49 (td, *J* = 8.0 1.2 Hz, 1H), 6.92 (d, *J* = 2.9 Hz, 1H).

¹³C NMR (101 MHz, CDCl₃) δ 143.33, 137.03, 130.41, 128.85, 127.39, 126.50, 125.95, 120.86, 115.32, 99.35, 96.15.

HRMS (APCI): m/z calcd for C₁₁H₇BrN₂ [M+H]⁺ : 246.9865, found: 246.9863.

3-bromo-4-phenylpyrrolo[1,2-*a*]quinoxaline [3b]



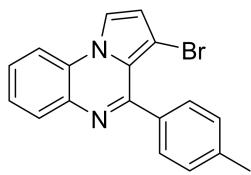
Mp. 161-166 °C. Light yellow solid (48.6 mg, 76%).

¹H NMR (400 MHz, CDCl₃) δ 8.03 (dd, *J* = 8.0, 1.2 Hz, 1H), 7.94 (d, *J* = 3.2 Hz, 1H), 7.83 (dd, *J* = 8.0, 1.2 Hz, 1H), 7.65-7.61 (m, 2H), 7.54 (dd, *J* = 7.2, 1.6 Hz, 1H), 7.51 (dd, *J* = 4.0, 2.4 Hz, 3H), 7.47 (td, *J* = 8.0, 1.2 Hz, 1H), 6.90 (d, *J* = 3.2 Hz, 1H).

¹³C NMR (101 MHz, CDCl₃) δ 154.93, 137.49, 135.58, 130.36, 129.49, 129.37, 128.08, 128.05, 127.02, 125.87, 121.87, 118.21, 114.64, 113.14, 95.74.

HRMS (APCI): m/z calcd for C₁₇H₁₁BrN₂ [M+H]⁺ : 323.0178, found: 323.0176.

3-bromo-4-(p-tolyl)pyrrolo[1,2-*a*]quinoxaline [3c]



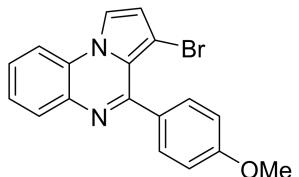
Mp. 174-179 °C. Light yellow solid (24.2 mg, 36%).

¹H NMR (400 MHz, CDCl₃) δ 8.02 (dd, *J* = 8.0, 1.6 Hz, 1H), 7.94 (d, *J* = 2.8 Hz, 1H), 7.83 (dd, *J* = 8.0, 1.2 Hz, 1H), 7.56-7.50 (m, 3H), 7.46 (td, *J* = 7.6, 1.6 Hz, 1H), 7.31 (d, *J* = 8.0 Hz, 2H), 6.91 (d, *J* = 3.2 Hz, 1H), 2.46 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 154.37, 139.44, 136.77, 134.42, 130.32, 129.37, 128.80, 128.35, 126.79, 126.36, 124.46, 123.00, 115.24, 99.78, 96.75, 21.65.

HRMS (APCI): m/z calcd for C₁₈H₁₃BrN₂ [M+H]⁺ : 337.0335, found: 337.0330.

3-bromo-4-(4-methoxyphenyl)pyrrolo[1,2-a]quinoxaline [3d]



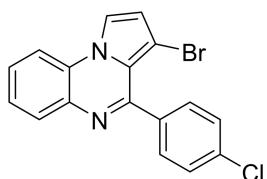
Mp. 167-169 °C. Light yellow solid (53.1 mg, 76%).

¹H NMR (400 MHz, CDCl₃) δ 8.00 (dd, *J* = 8.0, 1.6 Hz, 1H), 7.91 (d, *J* = 2.8 Hz, 1H), 7.80 (dd, *J* = 8.0, 1.6 Hz, 1H), 7.61 (d, *J* = 2.0 Hz, 1H), 7.59 (d, *J* = 2.0 Hz, 1H), 7.49 (td, *J* = 8.0, 1.6 Hz, 1H), 7.44 (td, *J* = 7.6, 1.2 Hz, 1H), 7.04 (d, *J* = 2.4 Hz, 1H), 7.02 (d, *J* = 2.0 Hz, 1H), 6.89 (d, *J* = 2.8 Hz, 1H), 3.89 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 160.73, 154.54, 135.55, 131.08, 130.12, 129.87, 127.82, 126.85, 125.78, 121.94, 118.17, 114.64, 113.41, 113.07, 95.70, 55.45.

HRMS (APCI): m/z calcd for C₁₈H₁₃BrN₂O [M+H]⁺ : 353.0284, found: 353.0280.

3-bromo-4-(4-chlorophenyl)pyrrolo[1,2-a]quinoxaline [3e]



Mp. 199-201°C. Light yellow solid (49.5 mg, 70%).

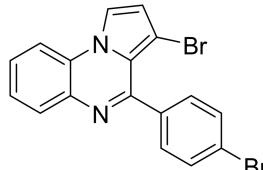
¹H NMR (400 MHz, CDCl₃) δ 8.01 (dd, *J* = 8.0, 1.6 Hz, 1H), 7.96 (d, *J* = 2.8 Hz, 1H), 7.85 (dd, *J* = 8.0, 1.2 Hz, 1H), 7.60-7.57 (m, 2H), 7.54 (dd, *J* = 8.4,

1.6 Hz, 1H), 7.50-7.45 (m, 3H), 6.92 (d, J = 3.2 Hz, 1H).

^{13}C NMR (101 MHz, CDCl_3) δ 153.72, 135.92, 135.59, 135.51, 131.04, 130.40, 128.36, 128.33, 127.04, 126.03, 121.69, 118.33, 114.87, 113.22, 95.71 .

HRMS (APCI): m/z calcd for $\text{C}_{17}\text{H}_{10}\text{BrClN}_2$ [M+H] $^+$: 356.9789, found: 356.9762.

3-bromo-4-(4-bromophenyl)pyrrolo[1,2-a]quinoxaline [3f]



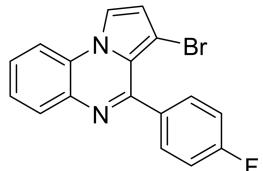
Mp. 202-205 °C. Light yellow solid (45.9 mg, 58%).

^1H NMR (400 MHz, CDCl_3) δ 8.01 (dd, J = 8.0, 1.6 Hz, 1H), 7.97 (d, J = 2.8 Hz, 1H), 7.85 (dd, J = 8.0, 1.2 Hz, 1H), 7.67-7.62 (m, 2H), 7.58-7.53 (m, 2H), 7.52-7.50 (m, 1H), 7.49-7.46 (m, 1H), 6.92 (d, J = 2.8 Hz, 1H).

^{13}C NMR (101 MHz, CDCl_3) δ 153.69, 136.26, 135.39, 131.30, 131.26, 130.32, 128.39, 127.00, 126.05, 123.91, 121.59, 118.36, 114.95, 113.23, 95.82.

HRMS (APCI): m/z calcd for $\text{C}_{17}\text{H}_{10}\text{Br}_2\text{N}_2$ [M+H] $^+$: 400.9284, found: 400.9261.

3-bromo-4-(4-fluorophenyl)pyrrolo[1,2-a]quinoxaline [3g]



Mp. 174-176 °C. Light yellow solid (59.3 mg, 87%).

^1H NMR (400 MHz, CDCl_3) δ 7.91 (dd, J = 8.0, 1.6 Hz, 1H), 7.85 (d, J = 2.8 Hz, 1H), 7.74 (dd, J = 8.0, 1.6 Hz, 1H), 7.57-7.49 (m, 2H), 7.44 (td, J = 8.0, 7.6, 1.6 Hz, 1H), 7.38 (td, J = 8.0, 7.6, 1.6 Hz, 1H), 7.14-7.06 (m, 2H), 6.82 (d, J = 2.8 Hz, 1H).

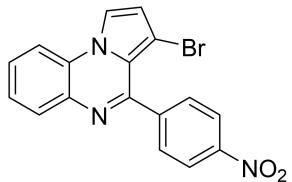
^{13}C NMR (101 MHz, CDCl_3) δ 163.70 (d, J = 249.3 Hz), 153.81, 135.43, 133.47 (d, J = 3.3 Hz), 131.54 (d, J = 8.4 Hz), 130.26, 128.20, 126.95, 125.94, 121.75, 118.26, 115.08 (d, J = 21.9 Hz), 114.80, 113.15, 95.70.

^{19}F NMR (376 MHz, CDCl_3) δ -111.96.

HRMS (APCI): m/z calcd for $\text{C}_{17}\text{H}_{10}\text{BrFN}_2$ [M+H] $^+$: 341.0084, found:

341.0083.

3-bromo-4-(4-nitrophenyl)pyrrolo[1,2-*a*]quinoxaline [3h]



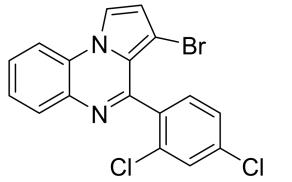
Mp. 244-249 °C. Yellow solid (69.1 mg, 94%).

¹H NMR (400 MHz, CDCl₃) δ 8.38 (d, *J* = 8.8 Hz, 2H), 8.07-8.00 (m, 2H), 7.90 (dd, *J* = 8.4, 0.8 Hz, 1H), 7.83 (d, *J* = 8.4 Hz, 2H), 7.61 (td, *J* = 8.4, 7.2, 1.2 Hz, 1H), 7.53 (td, *J* = 8.4, 7.2, 1.2 Hz, 1H), 6.97 (d, *J* = 2.8 Hz, 1H).

¹³C NMR (151 MHz, CDCl₃) δ 152.57, 148.62, 143.70, 135.34, 130.86, 130.58, 128.98, 127.13, 126.32, 123.37, 121.35, 118.53, 115.27, 113.38, 95.67.

HRMS (APCI): m/z calcd for C₁₇H₁₀BrN₃O₂ [M+H]⁺ : 368.0029, found: 368.0027.

3-bromo-4-(2,4-dichlorophenyl)pyrrolo[1,2-*a*]quinoxaline [3i]



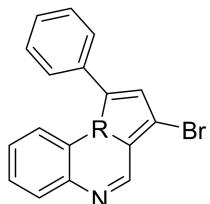
Mp. 137-141 °C. Light yellow solid (63.3 mg, 81%).

¹H NMR (400 MHz, CDCl₃) δ 8.03 (dd, *J* = 8.0, 1.6 Hz, 1H), 7.97 (d, *J* = 2.8 Hz, 1H), 7.87 (dd, *J* = 8.0, 1.2 Hz, 1H), 7.59 (td, *J* = 8.0, 1.2 Hz, 1H), 7.54 (t, *J* = 1.2 Hz, 1H), 7.50 (td, *J* = 8.0, 1.2 Hz, 2H), 7.41 (d, *J* = 1.6 Hz, 2H), 6.91 (d, *J* = 3.2 Hz, 1H);

¹³C NMR (101 MHz, CDCl₃) δ 151.39, 136.02, 134.94, 131.64, 130.35, 129.47, 128.88, 127.37, 126.23, 121.88, 118.17, 115.17, 113.41.

HRMS (APCI): m/z calcd for C₁₇H₉BrCl₂N₂ [M+H]⁺ : 390.9398, found: 390.9395.

3-bromo-1-phenylpyrrolo[1,2-*a*]quinoxaline [3j]



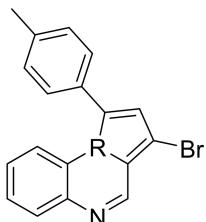
Mp. 153-158 °C. Light yellow solid (49.8 mg, 39%).

¹H NMR (400 MHz, CDCl₃) δ 8.83 (s, 1H), 7.96 (dd, *J* = 8.0, 1.6 Hz, 1H), 7.51 (t, *J* = 3.2 Hz, 5H), 7.40-7.31 (m, 2H), 7.14 (td, *J* = 7.2, 1.6 Hz, 1H), 6.81 (s, 1H).

¹³C NMR (101 MHz, CDCl₃) δ 144.35, 137.21, 132.99, 132.49, 130.33, 129.84, 129.36, 129.02, 128.52, 127.16, 125.67, 124.79, 118.76, 116.50, 95.96.

HRMS (APCI): m/z calcd for C₁₇H₁₁BrN₂ [M+H]⁺ : 323.0178, found: 323.0177.

3-bromo-1-(p-tolyl)pyrrolo[1,2-*a*]quinoxaline [3k]



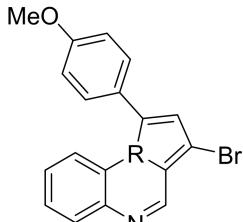
Mp. 118-123 °C. Light yellow solid (55.0 mg, 41%).

¹H NMR (400 MHz, CDCl₃) δ 8.81 (s, 1H), 7.95 (dd, *J* = 8.0, 1.6 Hz, 1H), 7.41-7.34 (m, 4H), 7.31 (d, *J* = 8.0 Hz, 2H), 7.14 (td, *J* = 8.0, 1.6 Hz, 1H), 6.77 (s, 1H), 2.48 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 144.38, 139.40, 137.27, 132.62, 130.30, 130.03, 129.71, 128.63, 127.05, 125.58, 124.70, 118.63, 116.51, 95.82, 21.60.

HRMS (APCI): m/z calcd for C₁₈H₁₃BrN₂ [M+H]⁺ : 337.0335, found: 337.0333.

3-bromo-1-(4-methoxyphenyl)pyrrolo[1,2-*a*]quinoxaline [3l]



Mp. 148-150 °C. Light yellow solid (65.2 mg, 93%).

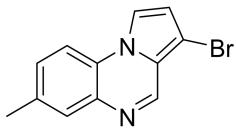
¹H NMR (400 MHz, CDCl₃) δ 8.80 (s, 1H), 7.94 (dd, *J* = 8.0, 1.6 Hz, 1H), 7.42-7.33 (m, 4H), 7.17-7.11 (m, 1H), 7.03 (d, *J* = 8.4 Hz, 2H), 6.75 (s, 1H), 3.91 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 160.42, 144.34, 137.26, 132.37, 131.17, 130.28, 128.67, 127.06, 125.54, 125.12, 124.58, 118.57, 116.38, 114.42, 95.74, 55.53.

HRMS (APCI): m/z calcd for C₁₈H₁₃BrN₂O [M+H]⁺ : 353.0284, found:

353.0283.

3-bromo-7-methylpyrrolo[1,2-*a*]quinoxaline [3m]



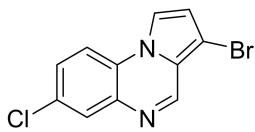
Mp. 174-179 °C. Light yellow solid (51.0 mg, 98%).

¹H NMR (400 MHz, CDCl₃) δ 8.72 (s, 1H), 7.74 (dd, *J* = 7.2, 2.4 Hz, 2H), 7.64 (d, *J* = 8.4 Hz, 1H), 7.30 (dd, *J* = 8.4, 2.0 Hz, 1H), 6.83 (d, *J* = 2.8 Hz, 1H), 2.47 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 143.86, 135.79, 135.71, 130.08, 129.42, 125.16, 123.73, 115.98, 113.89, 113.02, 94.65, 21.20.

HRMS (APCI): m/z calcd for C₁₂H₉BrN₂ [M+H]⁺ : 261.0022, found: 221.0019.

3-bromo-7-chloropyrrolo[1,2-*a*]quinoxaline [3n]



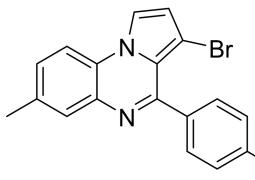
Mp. 167-172 °C. Light yellow solid (33.9 mg, 61%).

¹H NMR (400 MHz, CDCl₃) δ 8.78 (s, 1H), 7.96 (d, *J* = 2.4 Hz, 1H), 7.82 (dd, *J* = 2.8, 0.8 Hz, 1H), 7.74 (d, *J* = 8.8 Hz, 1H), 7.49 (dd, *J* = 8.8, 2.4 Hz, 1H), 6.91 (d, *J* = 2.8 Hz, 1H).

¹³C NMR (101 MHz, CDCl₃) δ 145.02, 136.68, 131.24, 129.73, 128.51, 126.05, 123.82, 116.86, 114.69, 114.61, 96.10.

HRMS (APCI): m/z calcd for C₁₁H₆BrClN₂ [M+H]⁺ : 280.9476, found: 280.9473.

3-bromo-4-(4-methoxyphenyl)-7-methylpyrrolo[1,2-*a*]quinoxaline [3o]



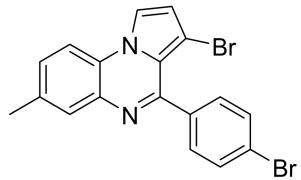
Mp. 162-167 °C. Light yellow solid (47.4 mg, 65%)

¹H NMR (400 MHz, CDCl₃) δ 7.89 (d, *J* = 2.8 Hz, 1H), 7.82 (s, 1H), 7.71 (d, *J* = 8.4 Hz, 1H), 7.60 (d, *J* = 2.0 Hz, 1H), 7.58 (d, *J* = 2.0 Hz, 1H), 7.32 (dd, *J* = 8.4, 2.0 Hz, 1H), 7.03 (d, *J* = 2.0 Hz, 1H), 7.02 (d, *J* = 2.0 Hz, 1H), 6.88 (d, *J* = 3.2 Hz, 1H), 3.89 (s, 3H), 2.48 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 160.79, 154.48, 135.80, 135.33, 131.16, 129.82, 129.08, 124.79, 121.86, 118.06, 114.60, 113.43, 112.86, 95.66, 55.49, 21.23.

HRMS (APCI): m/z calcd for C₁₉H₁₅BrN₂O [M+H]⁺ : 367.0441, found: 367.0437.

3-bromo-4-(4-bromophenyl)-7-methylpyrrolo[1,2-*a*]quinoxaline [3p]



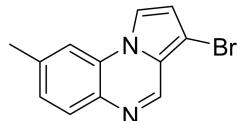
Mp. 216-219 °C. Light yellow solid (39.5 mg, 51%).

¹H NMR (400 MHz, CDCl₃) δ 7.68 (d, *J* = 3.2 Hz, 1H), 7.57 (s, 1H), 7.49 (d, *J* = 8.4 Hz, 1H), 7.39 (dd, *J* = 6.4, 2.0 Hz, 2H), 7.27 (dd, *J* = 6.4, 2.0 Hz, 2H), 7.12 (dd, *J* = 8.4, 2.0 Hz, 1H), 6.65 (d, *J* = 2.8 Hz, 1H), 2.25 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 153.60, 136.40, 135.99, 135.36, 131.23, 130.09, 129.53, 124.90, 123.82, 121.52, 118.09, 114.71, 112.95, 95.41, 21.25.

HRMS (APCI): m/z calcd for C₁₈H₁₂Br₂N₂ [M+H]⁺ : 414.9440, found: 414.9436.

3-bromo-8-methylpyrrolo[1,2-*a*]quinoxaline [3q]



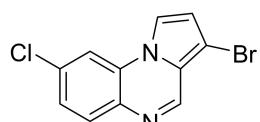
Mp. 169-174 °C. Light yellow solid (51.5 mg, 98%).

¹H NMR (400 MHz, CDCl₃) δ 8.72 (s, 1H), 7.84 (d, *J* = 8.4 Hz, 1H), 7.79 (d, *J* = 3.2 Hz, 1H), 7.58 (d, *J* = 2.0 Hz, 1H), 7.26 (dd, *J* = 8.0, 2.0 Hz, 1H), 6.86 (d, *J* = 2.8 Hz, 1H), 2.53 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 143.03, 139.08, 133.78, 130.02, 127.22, 127.13, 123.92, 116.23, 113.86, 113.42, 94.69, 21.96.

HRMS (APCI): m/z calcd for C₁₂H₉BrN₂ [M+H]⁺ : 261.0022, found: 261.0020.

3-bromo-8-chloropyrrolo[1,2-*a*]quinoxaline [3r]



Mp. 169-174 °C. Light yellow solid (28.7 mg, 51%).

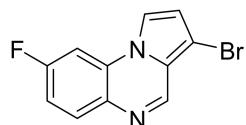
¹H NMR (400 MHz, CDCl₃) δ 8.82 (s, 1H), 8.04 (d, *J* = 8.8 Hz, 1H), 7.88 (d, *J*

= 2.8 Hz, 1H), 7.85 (d, J = 2.0 Hz, 1H), 7.47 (dd, J = 8.4, 2.0 Hz, 1H), 7.00 (d, J = 2.8 Hz, 1H).

^{13}C NMR (101 MHz, CDCl_3) δ 143.51, 134.45, 130.95, 127.97, 127.01, 126.74, 123.73, 117.50, 115.46, 115.36, 113.81.

HRMS (APCI): m/z calcd for $\text{C}_{11}\text{H}_6\text{BrClN}_2$ $[\text{M}+\text{H}]^+$: 280.9476, found: 280.9474.

3-bromo-8-fluoropyrrolo[1,2-a]quinoxaline [3s]



Mp. 164-166 °C. Light yellow solid (43.3 mg, 84%).

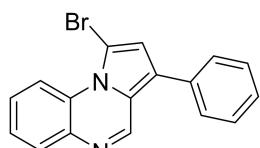
^1H NMR (400 MHz, CDCl_3) δ 8.78 (s, 1H), 8.02 (dd, J = 8.8, 5.6 Hz, 1H), 7.78 (d, J = 3.2 Hz, 1H), 7.51 (dd, J = 8.8, 2.4 Hz, 1H), 7.22 (td, J = 9.2, 2.8 Hz, 1H), 6.95 (d, J = 2.8 Hz, 1H).

^{13}C NMR (151 MHz, CDCl_3) δ 161.94 (d, J = 250.1 Hz), 143.25, 132.33 (d, J = 9.6 Hz), 116.98, 114.50, 114.10, 100.55, 100.37.

^{19}F NMR (565 MHz, CDCl_3) δ -109.16.

HRMS (APCI): m/z calcd for $\text{C}_{11}\text{H}_6\text{BrFN}_2$ $[\text{M}+\text{H}]^+$: 264.9771, found: 264.9769.

1-bromo-3-phenylpyrrolo[1,2-a]quinoxaline [3t]



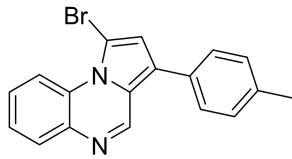
Mp. 137-141°C. Light yellow solid (58.2 mg, 90%).

^1H NMR (400 MHz, CDCl_3) δ 9.27 (dd, J = 8.0, 2.0 Hz, 1H), 8.90 (s, 1H), 7.98 (dd, J = 7.2, 2.4 Hz, 1H), 7.57 (d, J = 7.2 Hz, 2H), 7.53-7.46 (m, 4H), 7.40 (t, J = 7.2 Hz, 1H), 7.03 (s, 1H).

^{13}C NMR (101 MHz, CDCl_3) δ 144.08, 137.05, 133.00, 129.85, 129.23, 128.72, 127.86, 127.09, 126.11, 124.81, 124.70, 119.02, 115.61, 100.11.

HRMS (APCI): m/z calcd for $\text{C}_{17}\text{H}_{11}\text{BrN}_2$ $[\text{M}+\text{H}]^+$: 323.0178, found: 323.0177.

1-bromo-3-(p-tolyl)pyrrolo[1,2-a]quinoxaline [3u]



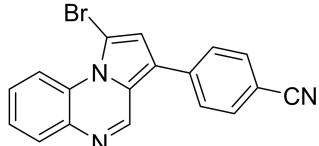
Mp. 150-155 °C. Light yellow solid (59.4 mg, 88%).

¹H NMR (400 MHz, CDCl₃) δ 9.26 (dd, *J* = 8.0, 2.0 Hz, 1H), 8.88 (s, 1H), 7.98 (dd, *J* = 7.2, 2.0 Hz, 1H), 7.54-7.45 (m, 4H), 7.30 (d, *J* = 7.6 Hz, 2H), 7.01 (s, 1H), 2.43 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 144.28, 137.76, 137.20, 130.10, 129.94, 129.90, 129.35, 128.58, 126.99, 126.03, 124.76, 124.67, 118.89, 115.59, 99.91, 21.37.

HRMS (APCI): m/z calcd for C₁₈H₁₃BrN₂ [M+H]⁺ : 337.0335, found: 337.0333.

4-(1-bromopyrrolo[1,2-a]quinoxalin-3-yl)benzonitrile [3v]



Mp. 240-242 °C. Light yellow solid (43.9mg, 63%).

¹H NMR (400 MHz, CDCl₃) δ 9.32 (dd, *J* = 8.0, 2.0 Hz, 1H), 8.90 (s, 1H), 8.04 (dd, *J* = 7.2, 2.0 Hz, 1H), 7.79 (d, *J* = 8.4 Hz, 2H), 7.68 (d, *J* = 8.4 Hz, 2H), 7.57 (pd, *J* = 7.2, 2.0 Hz, 2H), 7.08 (s, 1H).

¹³C NMR (101 MHz, CDCl₃) δ 143.08, 137.69, 133.06, 129.98, 129.10, 127.72, 126.66, 119.12, 118.81, 115.77, 111.40, 101.17.

HRMS (APCI): m/z calcd for C₁₈H₁₀BrN₃ [M+H]⁺ : 348.0131, found: 348.0129.

1,3-dibromopyrrolo[1,2-a]quinoxaline [4a]



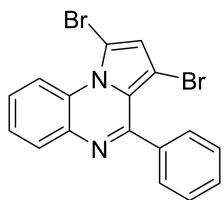
Mp. 157-162 °C. White solid (52.1 mg, 80%).

¹H NMR (400 MHz, CDCl₃) *I* 9.18-9.23 (m, 1H), 8.71 (s, 1H), 8.03-7.96 (m, 1H), 7.51-7.53 (m, 2H), 6.91 (s, 1H).

¹³C NMR (101 MHz, CDCl₃) δ 143.32, 137.00, 130.40, 128.87, 127.42, 126.53, 125.97, 120.90, 115.34, 99.42, 96.23.

HRMS (APCI): m/z calcd for C₁₁H₆Br₂N₂ [M+H]⁺ : 324.8971, found: 324.8968.

1,3-dibromo-4-phenylpyrrolo[1,2-*a*]quinoxaline [4b]



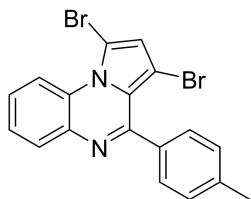
Mp. 151-153 °C. White solid (49.5 mg, 62%).

¹H NMR (400 MHz, CDCl₃) δ 9.21-9.13 (m, 1H), 7.96-7.90 (m, 1H), 7.51-7.45 (m, 2H), 7.44-7.38 (m, 5H), 6.84 (s, 1H).

¹³C NMR (101 MHz, CDCl₃) δ 154.29, 137.32, 136.77, 130.41, 129.43, 129.41, 128.43, 128.18, 126.94, 126.42, 124.38, 123.03, 115.29, 99.87, 96.74.

HRMS (APCI): m/z calcd for C₁₇H₁₀Br₂N₂ [M+H]⁺ : 400.9284, found: 400.9280.

1,3-dibromo-4-(p-tolyl)pyrrolo[1,2-*a*]quinoxaline [4c]



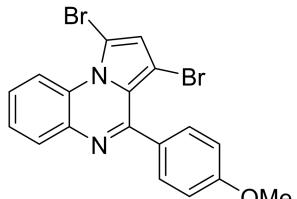
Mp. 159-162 °C. Light yellow solid (46.5 mg, 56%).

¹H NMR (400 MHz, CDCl₃) δ 9.29-9.21 (m, 1H), 8.05-7.99 (m, 1H), 7.53-7.47 (m, 3H), 7.46 (s, 1H), 7.31 (s, 1H), 7.29 (s, 1H), 6.93 (s, 1H), 2.45 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 155.01, 139.42, 135.52, 134.50, 130.24, 129.48, 128.72, 127.98, 126.97, 125.86, 121.94, 118.25, 114.69, 113.13, 95.88, 21.66.

HRMS (APCI): m/z calcd for C₁₈H₁₂Br₂N₂ [M+H]⁺ : 414.9440, found: 414.9435.

1,3-dibromo-4-(4-methoxyphenyl)pyrrolo[1,2-*a*]quinoxaline [4d]



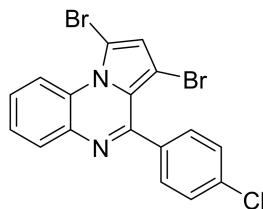
Mp. 159-162 °C. White solid (43.2 mg, 50%)

¹H NMR (400 MHz, CDCl₃) δ 9.22-9.15 (m, 1H), 7.95-7.97 (m, 1H), 7.46 (d, *J* = 2.0 Hz, 1H), 7.45-7.40 (m, 3H), 6.95 (d, *J* = 2.0 Hz, 1H), 6.93 (d, *J* = 2.0 Hz, 1H), 6.87 (s, 1H), 3.81 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 160.89, 154.01, 136.68, 131.05, 130.17, 128.33, 126.82, 126.45, 124.54, 123.14, 115.31, 113.60, 100.01, 97.18, 55.52.

HRMS (APCI): m/z calcd for C₁₈H₁₂Br₂N₂O [M+H]⁺ : 430.9389, found: 430.9385.

1,3-dibromo-4-(4-chlorophenyl)pyrrolo[1,2-*a*]quinoxaline [4e]



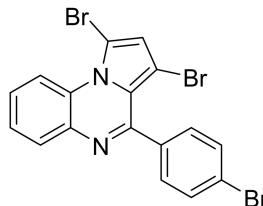
Mp. 189-193 °C. White solid (45.2 mg, 53%).

¹H NMR (400 MHz, CDCl₃) δ 9.31-9.22 (m, 1H), 8.04-7.97 (m, 1H), 7.51-7.55 (m, 4H), 7.46-7.49 (m, 2H), 6.94 (s, 1H).

¹³C NMR (101 MHz, CDCl₃) δ 153.07, 136.59, 135.71, 130.96, 130.36, 128.45, 128.43, 127.22, 126.58, 124.16, 123.15, 115.36, 100.26, 96.80.

HRMS (APCI): m/z calcd for C₁₇H₉Br₂ClN₂ [M+H]⁺ : 434.8894, found: 434.8890.

1,3-dibromo-4-(4-bromophenyl)pyrrolo[1,2-*a*]quinoxaline [4f]



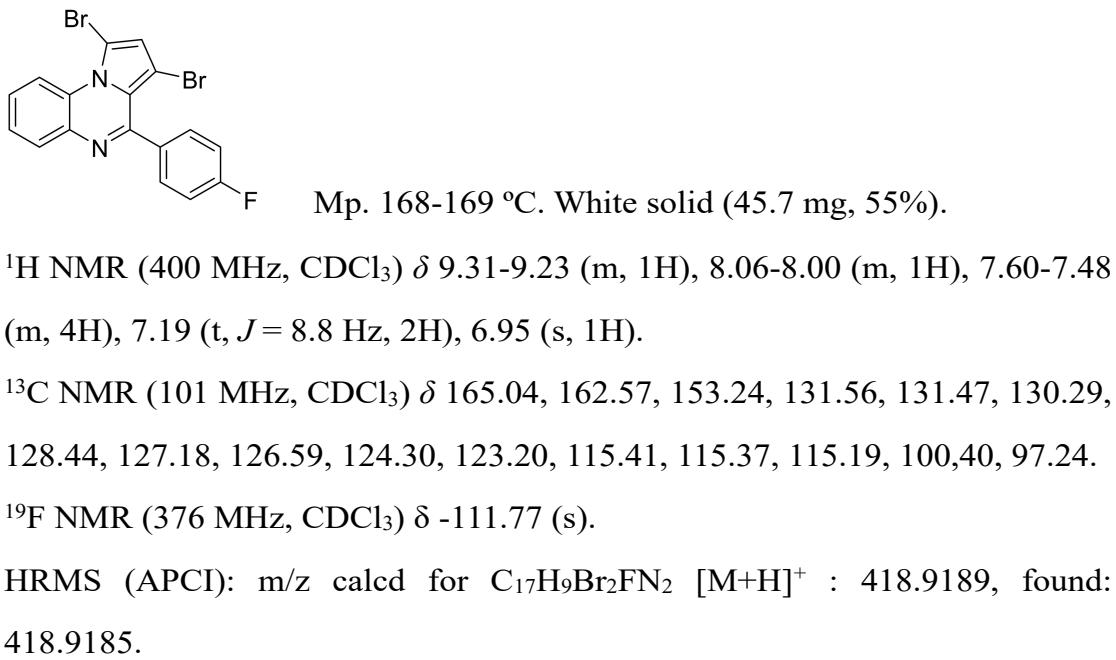
Mp. 207-211°C. White solid (49.5 mg, 53%).

¹H NMR (400 MHz, CDCl₃) δ 9.32-9.23 (m, 1H), 8.06-7.96 (m, 1H), 7.63 (d, *J* = 8.0 Hz, 2H), 7.58-7.49 (m, 2H), 7.46 (d, *J* = 8.0 Hz, 2H), 6.95 (s, 1H).

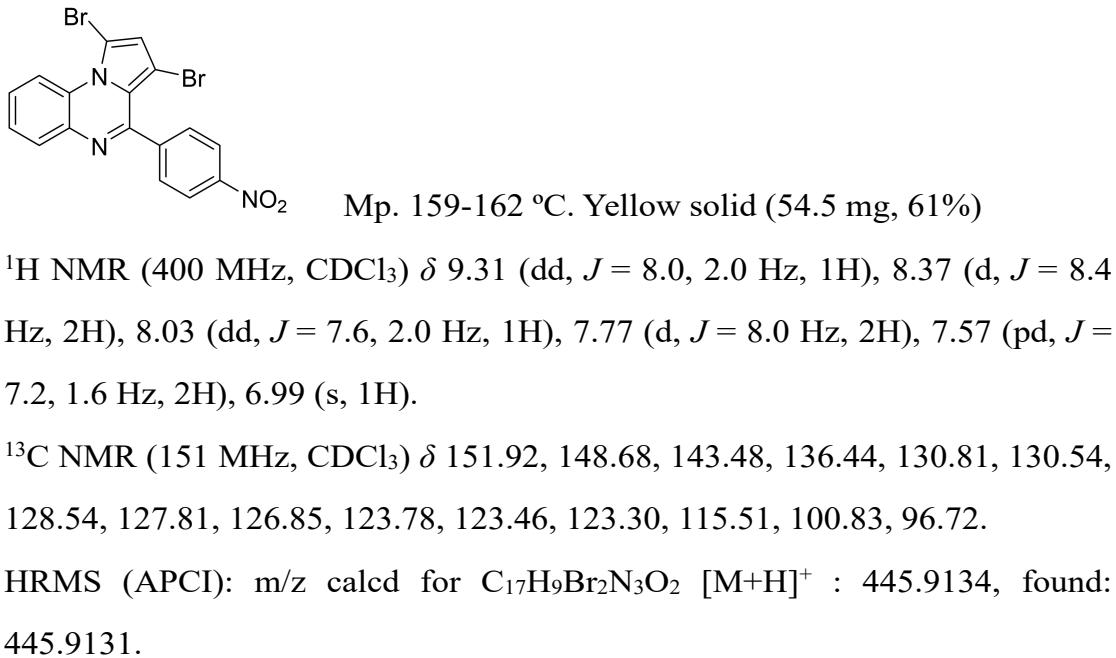
¹³C NMR (101 MHz, CDCl₃) δ 153.06, 136.53, 136.06, 131.39, 131.21, 130.33, 128.41, 127.25, 126.60, 124.08, 124.00, 123.17, 115.37, 100.33, 96.87.

HRMS (APCI): m/z calcd for C₁₇H₉Br₃N₂ [M+H]⁺ : 478.8389, found: 478.8386.

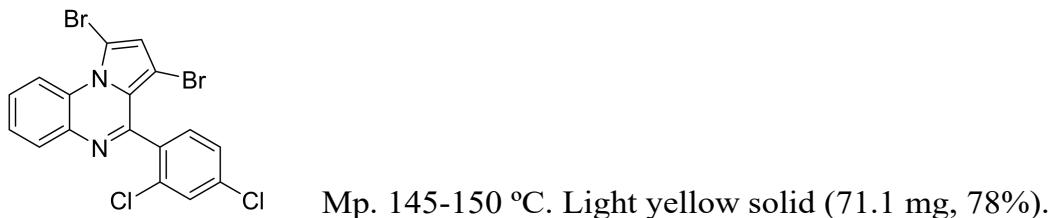
1,3-dibromo-4-(4-fluorophenyl)pyrrolo[1,2-*a*]quinoxaline [4g]



1,3-dibromo-4-(4-nitrophenyl)pyrrolo[1,2-a]quinoxaline [4h]



1,3-dibromo-4-(2,4-dichlorophenyl)pyrrolo[1,2-a]quinoxaline [4i]



Hz, 1H), 6.92 (s, 1H).

^{13}C NMR (151 MHz, CDCl_3) δ 150.72, 136.48, 136.00, 134.96, 131.62, 130.53, 130.17, 129.42, 128.76, 127.56, 127.38, 126.58, 124.15, 122.79, 115.36, 100.16, 96.64.

HRMS (APCI): m/z calcd for $\text{C}_{17}\text{H}_8\text{Br}_2\text{Cl}_2\text{N}_2$ $[\text{M}+\text{H}]^+$: 468.8504, found: 468.8500.

1,3-dibromo-7-methylpyrrolo[1,2-*a*]quinoxaline [4j]



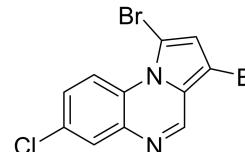
Mp. 149-151 °C. Light yellow solid (59.1 mg, 87%).

^1H NMR (400 MHz, CDCl_3) δ 8.99 (d, $J = 8.8$ Hz, 1H), 8.63 (s, 1H), 7.74-7.70 (m, 1H), 7.28 (dd, $J = 8.8, 2.0$ Hz, 1H), 6.84 (s, 1H), 2.47 (s, 3H).

^{13}C NMR (101 MHz, CDCl_3) δ 143.18, 136.95, 136.42, 130.14, 128.38, 126.58, 125.79, 120.43, 114.94, 98.99, 95.75, 77.48, 77.16, 76.84, 21.06.

HRMS (APCI): m/z calcd for $\text{C}_{12}\text{H}_8\text{Br}_2\text{N}_2$ $[\text{M}+\text{H}]^+$: 338.9127, found: 338.9124.

1,3-dibromo-7-chloropyrrolo[1,2-*a*]quinoxaline [4k]



Mp. 187-190 °C. Light yellow solid (36.3 mg, 51%).

^1H NMR (600 MHz, CDCl_3) δ 9.22 (d, $J = 2.2$ Hz, 1H), 9.22 (d, $J = 2.4$ Hz, 1H), 8.69 (s, 1H), 7.92 (d, $J = 8.4$ Hz, 1H), 7.47 (dd, $J = 9.0, 2.4$ Hz, 1H), 6.95 (s, 1H).

^{13}C NMR (151 MHz, CDCl_3) δ 143.53, 135.79, 132.81, 131.48, 129.24, 126.93, 125.93, 121.33, 115.42, 99.76, 96.69.

HRMS (APCI): m/z calcd for $\text{C}_{11}\text{H}_5\text{Br}_2\text{ClN}_2$ $[\text{M}+\text{H}]^+$: 358.8581, found: 358.8587.

1,3-dibromo-8-methylpyrrolo[1,2-*a*]quinoxaline [4l]



Mp. 166-171 °C. Light yellow solid (48.1 mg, 71%).

¹H NMR (400 MHz, CDCl₃) δ 8.64 (s, 1H), 8.64 (s, 1H), 7.86 (d, *J* = 8.4 Hz, 1H), 7.31 (dd, *J* = 8.4, 1.6 Hz, 1H), 6.88 (s, 1H), 2.53 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 142.25, 138.01, 134.75, 129.90, 128.62, 127.71, 125.97, 120.85, 115.33, 99.09, 95.95, 22.28.

HRMS (APCI): m/z calcd for C₁₂H₈Br₂N₂ [M+H]⁺ : 338.9127, found: 338.9125.

1,3-dibromo-8-chloropyrrolo[1,2-*a*]quinoxaline [4m]



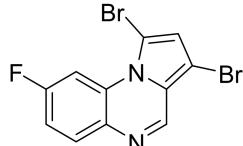
Mp. 197-199 °C. Light yellow solid (62.0 mg, 86%).

¹H NMR (600 MHz, CDCl₃) δ 9.15 (d, *J* = 9.0 Hz, 1H), 8.72 (s, 1H), 7.99 (d, *J* = 2.4 Hz, 1H), 7.49 (dd, *J* = 9.6, 2.4 Hz, 1H), 6.95 (s, 1H).

¹³C NMR (151 MHz, CDCl₃) δ 144.50, 138.36, 131.79, 129.79, 127.43, 127.31, 125.91, 121.12, 116.41, 99.77, 96.78.

HRMS (APCI): m/z calcd for C₁₁H₅Br₂ClN₂ [M+H]⁺ : 358.8581, found: 358.8586.

1,3-dibromo-8-fluoropyrrolo[1,2-*a*]quinoxaline [4n]



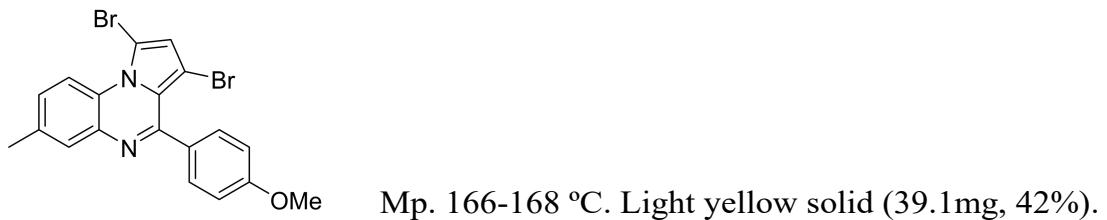
Mp. 184-186 °C. Light yellow solid (55.1 mg, 81%).

¹H NMR (400 MHz, CDCl₃) δ 8.93 (dd, *J* = 10.8, 2.4 Hz, 1H), 8.66 (s, 1H), 7.97 (dd, *J* = 8.8, 6.0 Hz, 1H), 7.27-7.21 (m, 1H), 6.93 (s, 1H).

¹³C NMR (151 MHz, CDCl₃) δ 160.53 (d, *J* = 248.6 Hz), 142.47, 133.51, 131.86 (d, *J* = 9.8 Hz), 125.59, 121.30, 114.48, 114.33, 102.57 (d, *J* = 29.8 Hz), 99.59, 96.57.

HRMS (APCI): m/z calcd for C₁₁H₅Br₂FN₂ [M+H]⁺ : 342.8876, found: 342.8872.

1,3-dibromo-4-(4-methoxyphenyl)-7-methylpyrrolo[1,2-*a*]quinoxaline [4o]

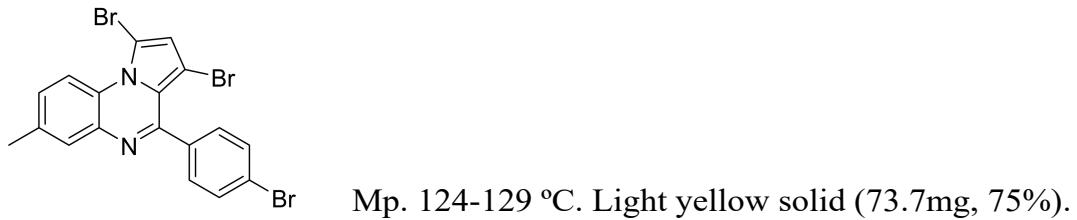


¹H NMR (400 MHz, CDCl₃) δ 7.89 (d, *J* = 2.8 Hz, 1H), 7.82 (s, 1H), 7.71 (d, *J* = 8.4 Hz, 1H), 7.60 (d, *J* = 2.0 Hz, 1H), 7.58 (d, *J* = 2.0 Hz, 1H), 7.32 (dd, *J* = 8.4, 2.0 Hz, 1H), 7.03 (d, *J* = 2.0 Hz, 1H), 7.02 (d, *J* = 1.8 Hz, 1H), 6.88 (d, *J* = 3.2 Hz, 1H), 3.89 (s, 3H), 2.48 (s, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 160.79, 153.92, 136.67, 136.36, 131.02, 129.97, 127.88, 126.15, 124.40, 122.73, 115.00, 113.54, 99.66, 96.52, 55.50, 21.05.

HRMS (APCI): m/z calcd for C₁₉H₁₄Br₂N₂O [M+H]⁺ : 444.9546, found: 444.9542.

1,3-dibromo-4-(4-bromophenyl)-7-methylpyrrolo[1,2-a]quinoxaline [4q]

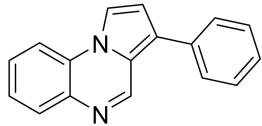


¹H NMR (400 MHz, CDCl₃) δ 9.14 (d, *J* = 8.8 Hz, 1H), 7.83 (s, 1H), 7.64 (d, *J* = 2.0 Hz, 1H), 7.62 (d, *J* = 2.0 Hz, 1H), 7.46 (d, *J* = 2.0 Hz, 1H), 7.44 (d, *J* = 2.0 Hz, 1H), 7.35 (dd, *J* = 8.8, 1.6 Hz, 1H), 6.93 (s, 1H), 2.49 (s, 3H).

¹³C NMR (151 MHz, CDCl₃) δ 153.04, 136.59, 136.33, 131.36, 131.22, 130.21, 128.35, 126.29, 124.01, 123.89, 122.80, 115.11, 99.98, 96.44, 21.06.

HRMS (APCI): m/z calcd for C₁₈H₁₁Br₃N₂ [M+H]⁺ : 492.8545, found: 492.8542.

3-phenylpyrrolo[1,2-a]quinoxaline [3aa]

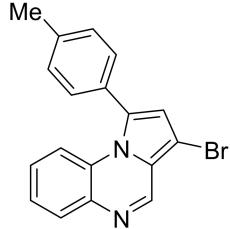


Mp. 178-180 °C. Light yellow solid (42.7 mg, 88%).

¹H NMR (400 MHz, CDCl₃) δ 9.04 (s, 1H), 7.97 (dd, *J* = 8.0, 1.6 Hz, 1H), 7.93 (d, *J* = 2.8 Hz, 1H), 7.84 (dd, *J* = 8.0, 1.2 Hz, 1H), 7.65 (d, *J* = 7.2 Hz, 2H),

7.55-7.42 (m, 4H), 7.37 (t, J = 7.2 Hz, 1H), 7.02 (d, J = 2.8 Hz, 1H).
 ^{13}C NMR (101 MHz, CDCl_3) δ 145.16, 136.15, 134.49, 130.08, 129.15, 128.36, 128.08, 128.03, 127.19, 125.49, 123.88, 122.75, 114.28, 113.84, 113.80.
HRMS (APCI): m/z calcd for $\text{C}_{17}\text{H}_{12}\text{N}_2$ [$\text{M}+\text{H}]^+$: 245.1073, found: 245.1071.

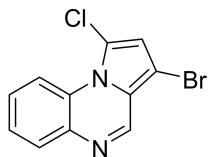
3-bromo-1-(p-tolyl)pyrrolo[1,2-a]quinoxaline [3ab]



Mp. 118-123 °C. Light yellow solid (56.9 mg, 85%).

^1H NMR (400 MHz, CDCl_3) δ 8.82 (s, 1H), 7.98 (d, J = 8.4 Hz, 1H), 7.35-7.41 (m, 4H), 7.32 (d, J = 8.0 Hz, 2H), 7.18-7.13 (m, 1H), 6.79 (s, 1H), 2.49 (s, 3H).
 ^{13}C NMR (101 MHz, CDCl_3) δ 144.00, 139.55, 136.70, 133.10, 129.95, 129.88, 129.76, 129.69, 128.59, 127.22, 125.74, 124.65, 118.93, 116.57, 96.56, 21.61.
HRMS (APCI): m/z calcd for $\text{C}_{18}\text{H}_{13}\text{BrN}_2$ [$\text{M}+\text{H}]^+$: 337.0335, found: 337.0331.

3-bromo-1-chloropyrrolo[1,2-a]quinoxaline [3ac]

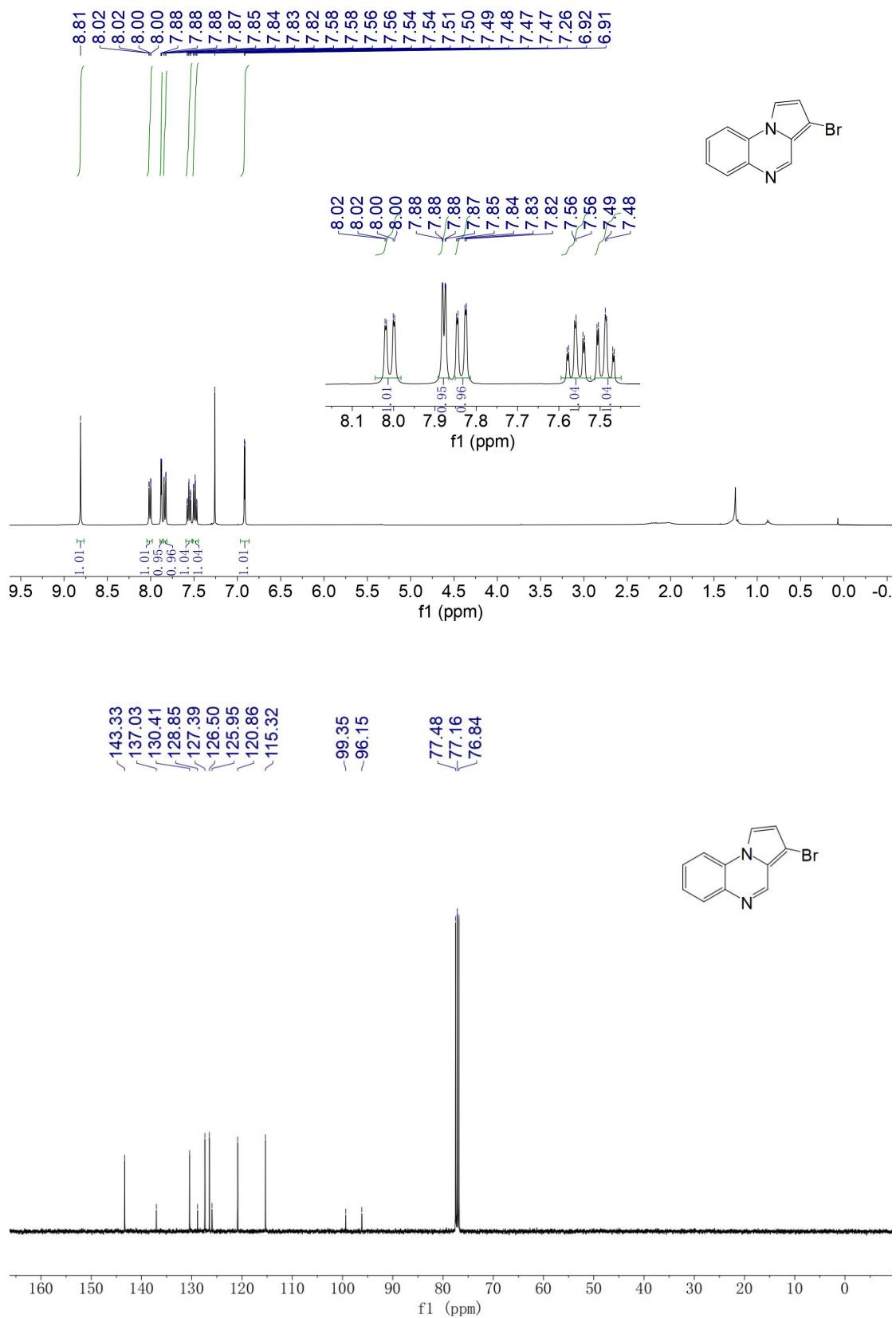


Mp 136-138 °C. Light yellow solid (53.9 mg, 96%).

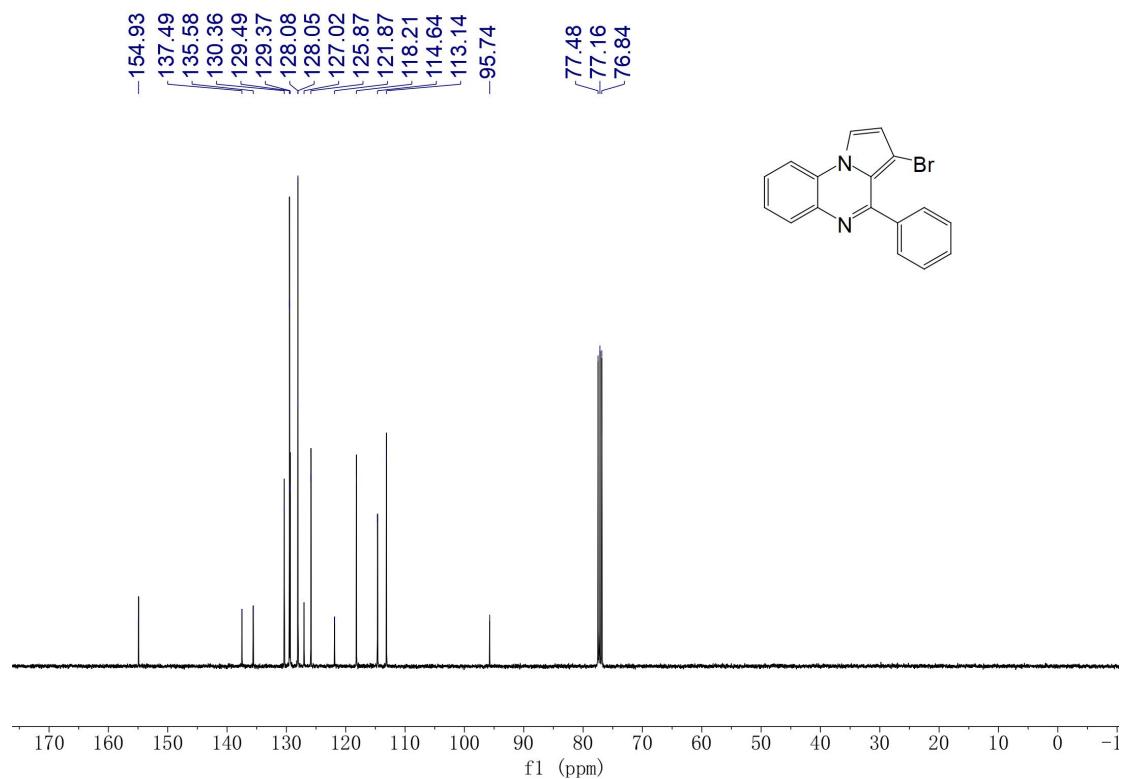
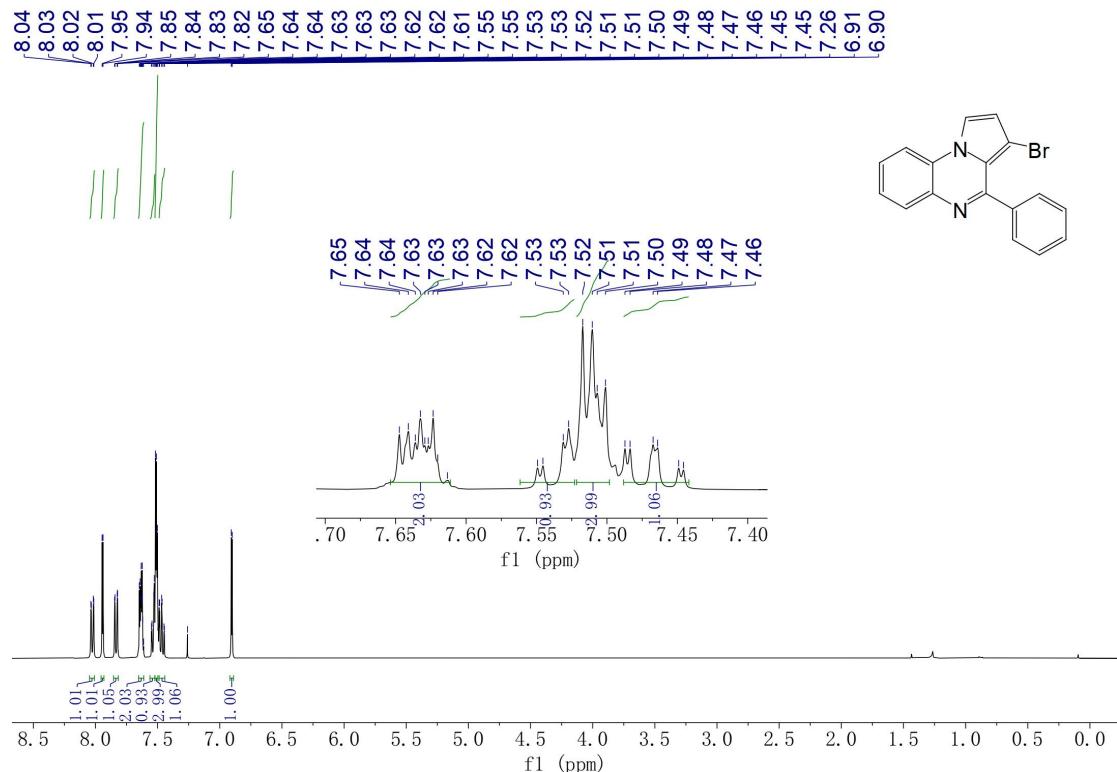
^1H NMR (400 MHz, CDCl_3) δ 8.93 (dd, J = 6.4, 3.6 Hz, 1H), 8.68 (s, 1H), 7.95 (dd, J = 6.4, 3.2 Hz, 1H), 7.48 (dd, J = 6.4, 3.2 Hz, 2H), 6.78 (s, 1H).
 ^{13}C NMR (101 MHz, CDCl_3) δ 143.51, 137.12, 130.42, 128.48, 127.57, 126.41, 124.31, 116.69, 115.76, 115.67, 95.02.
HRMS (APCI): m/z calcd for $\text{C}_{11}\text{H}_6\text{BrClN}_2$ [$\text{M}+\text{H}]^+$: 280.9476, found: 280.9473.

4. Copies of Spectra

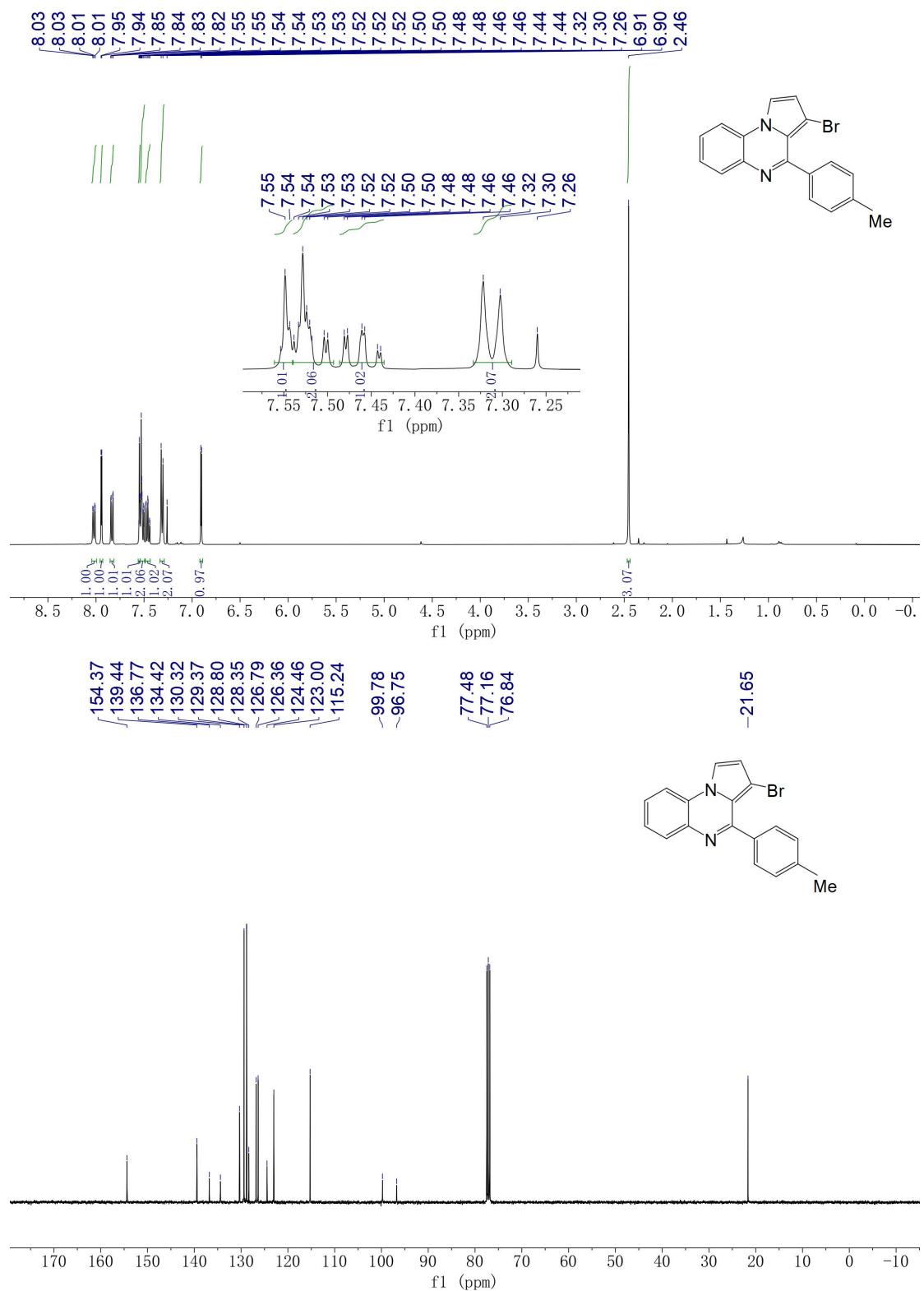
3-bromopyrrolo[1,2-a]quinoxaline [3a]



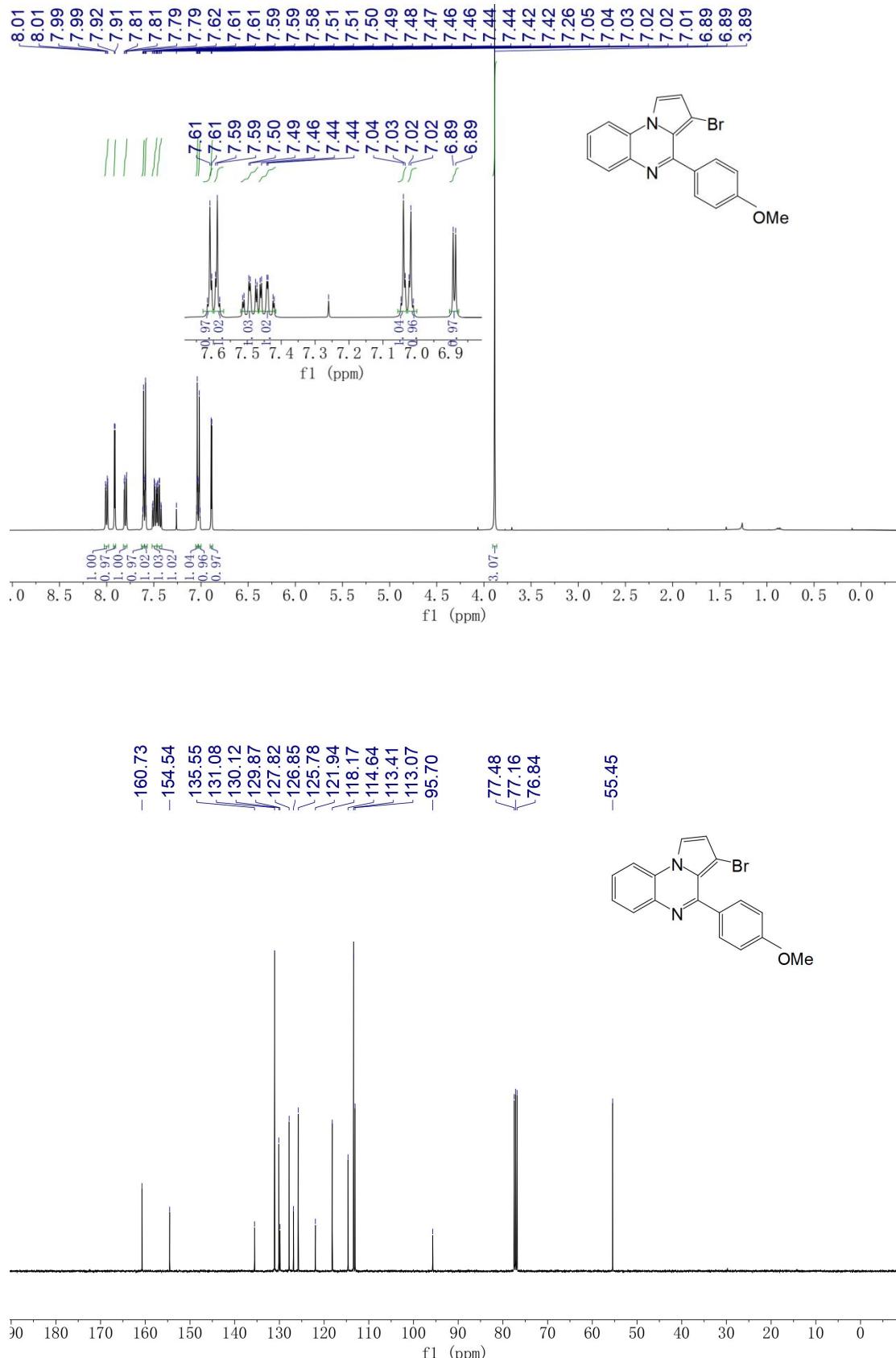
3-bromo-4-phenylpyrrolo[1,2-*a*]quinoxaline [3b]



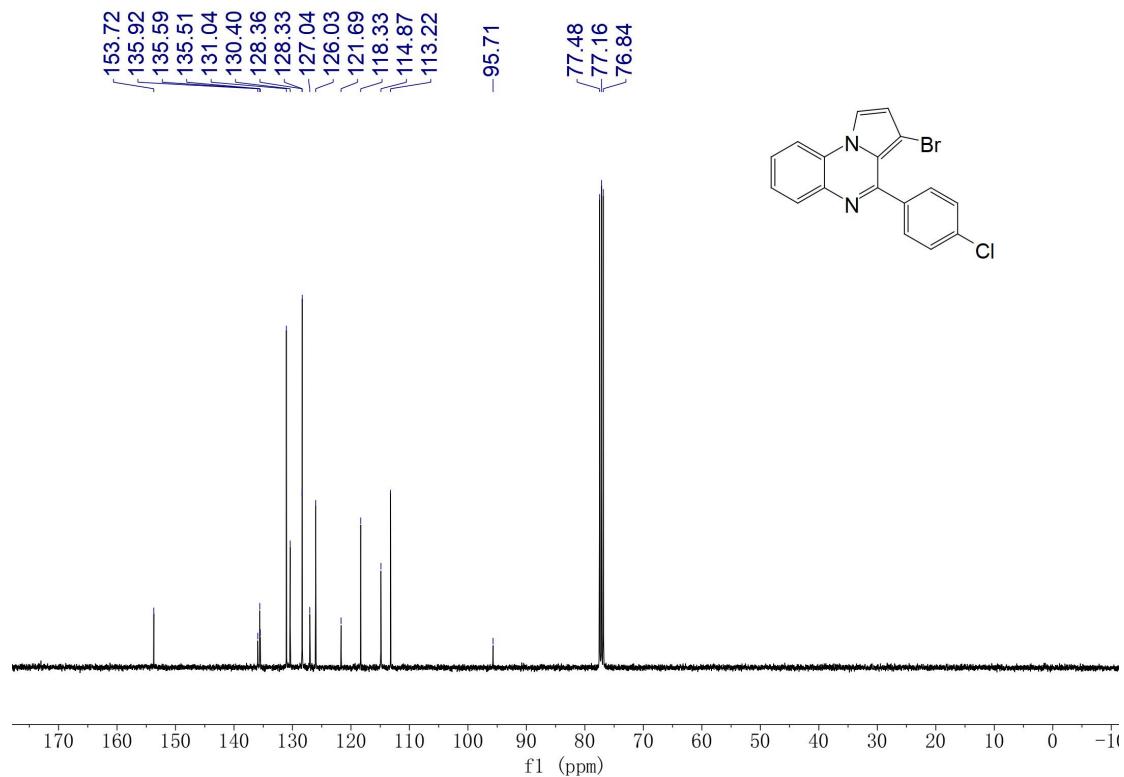
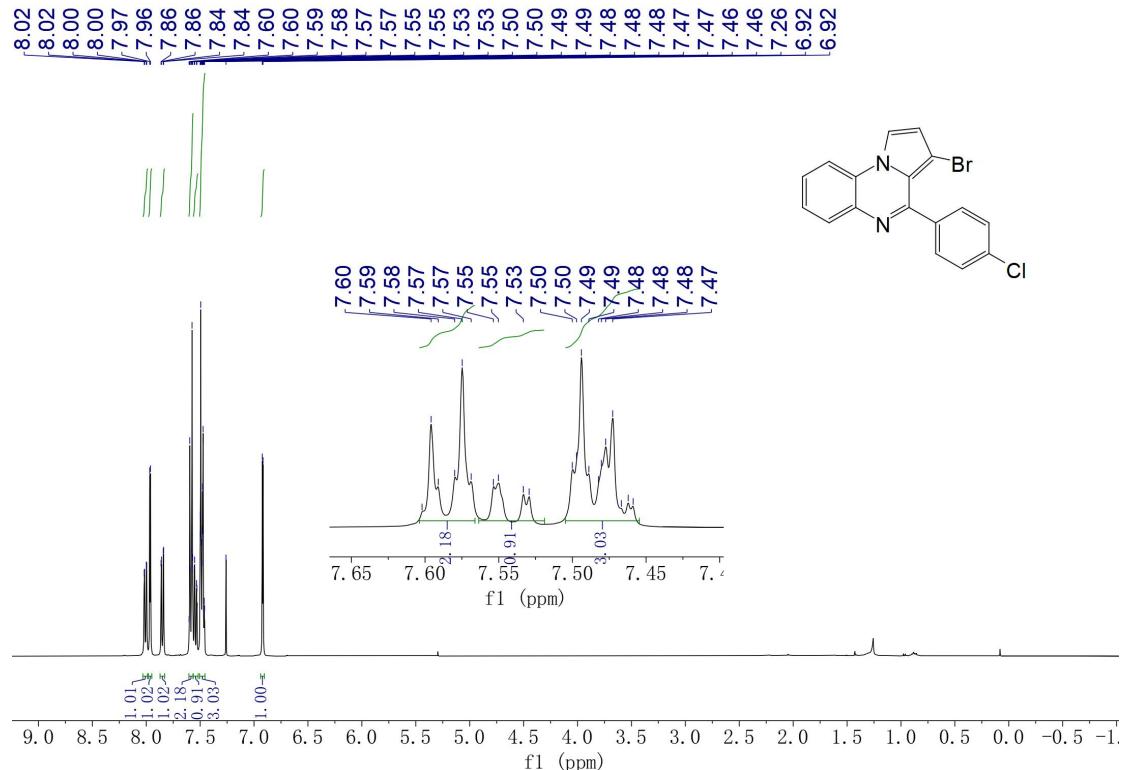
3-bromo-4-(p-tolyl)pyrrolo[1,2-*a*]quinoxaline [3c]



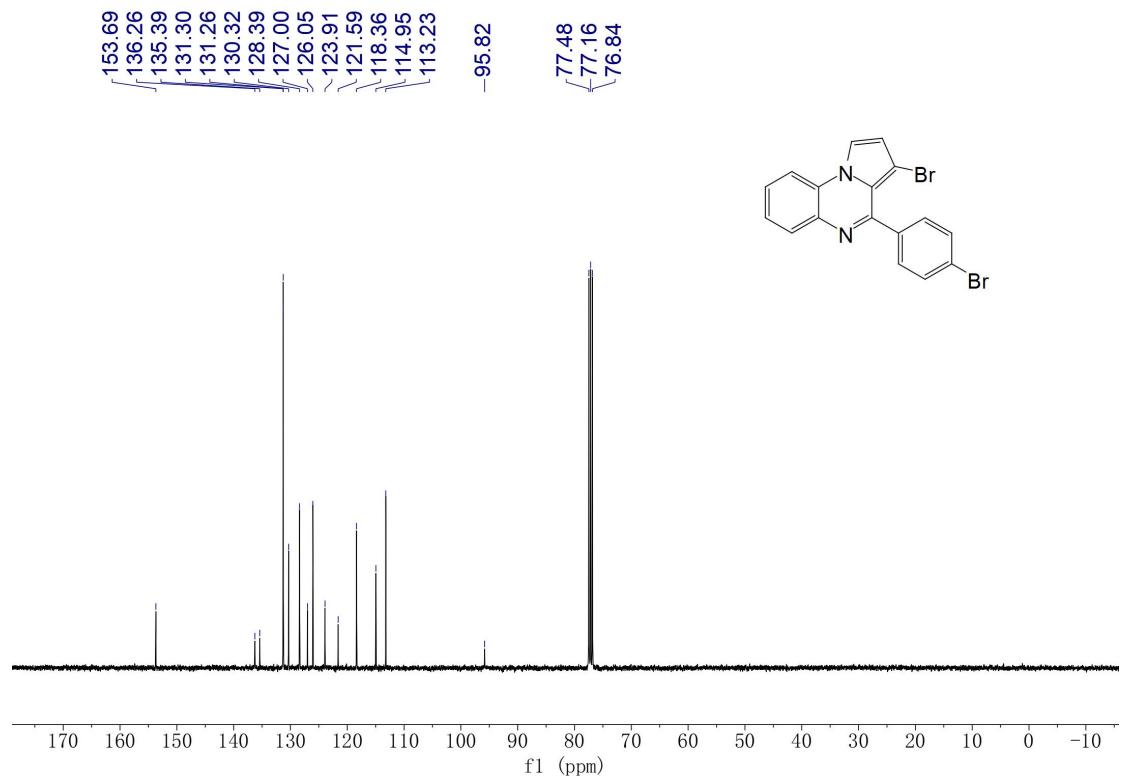
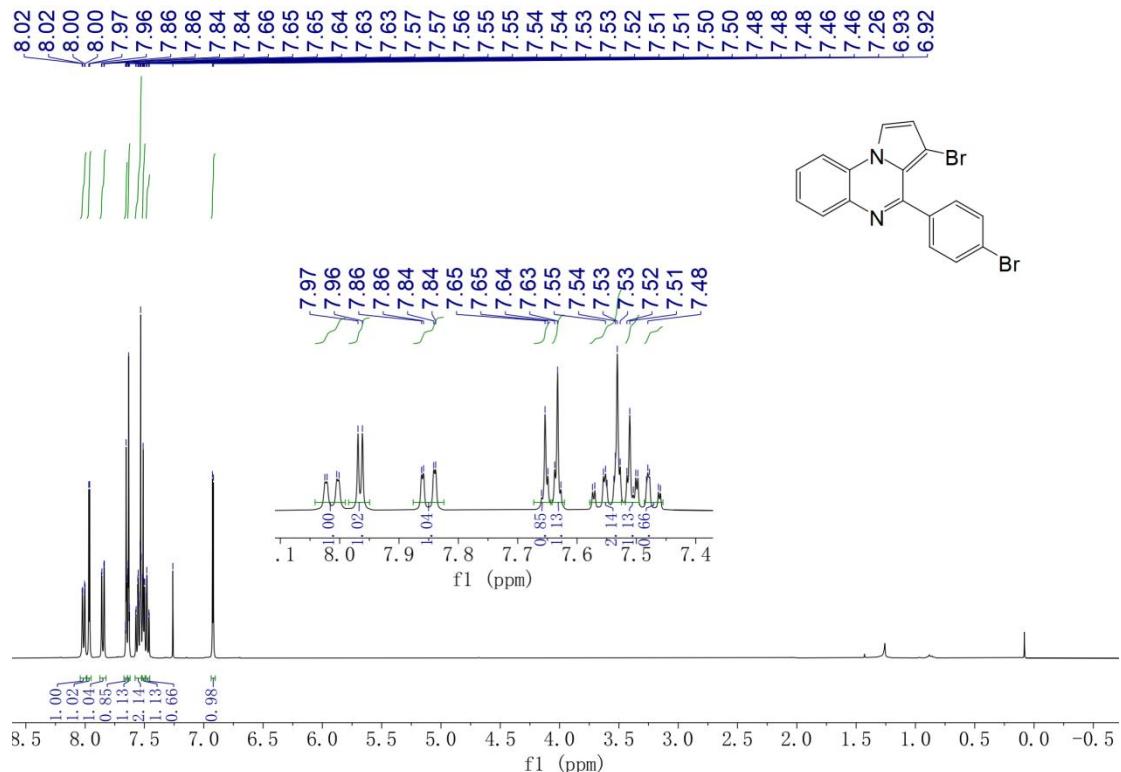
3-bromo-4-(4-methoxyphenyl)pyrrolo[1,2-*a*]quinoxaline [3d]



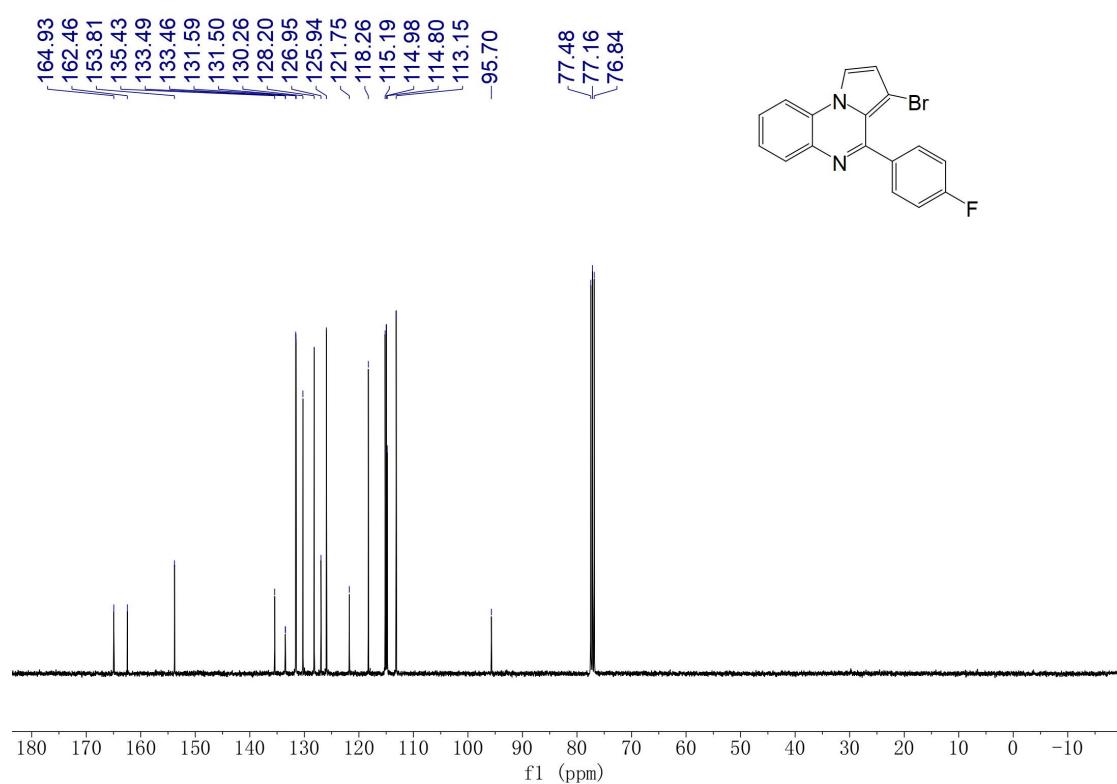
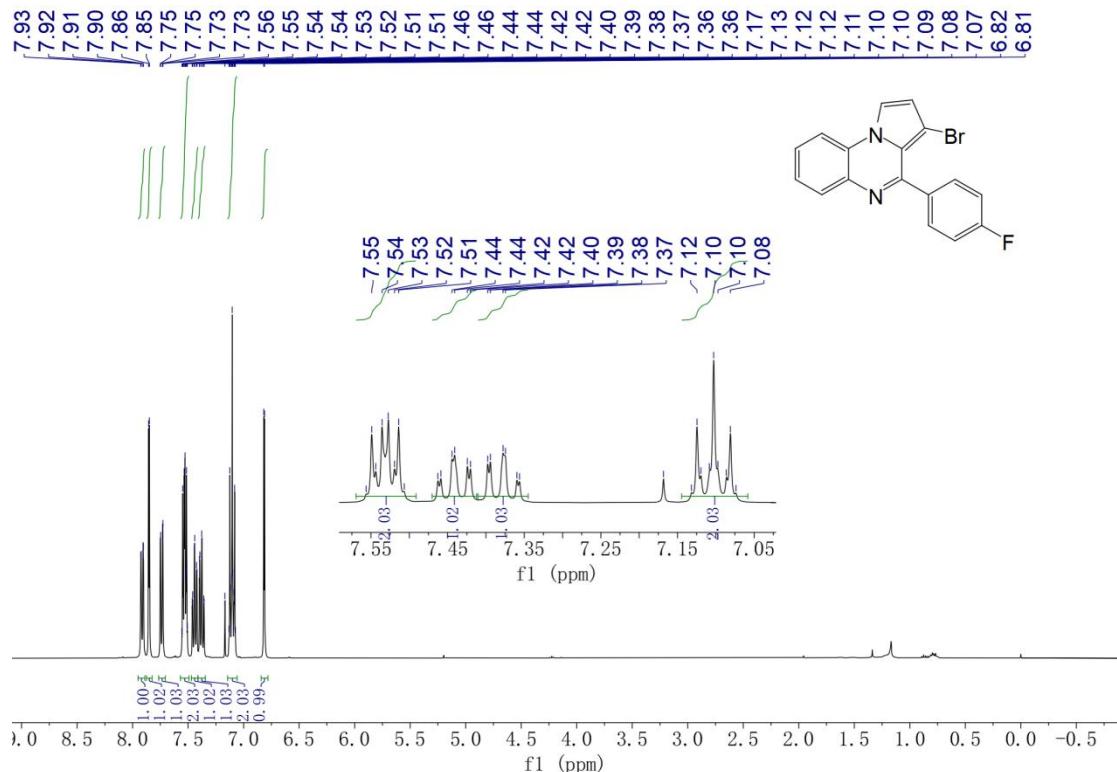
3-bromo-4-(4-chlorophenyl)pyrrolo[1,2-*a*]quinoxaline [3e]

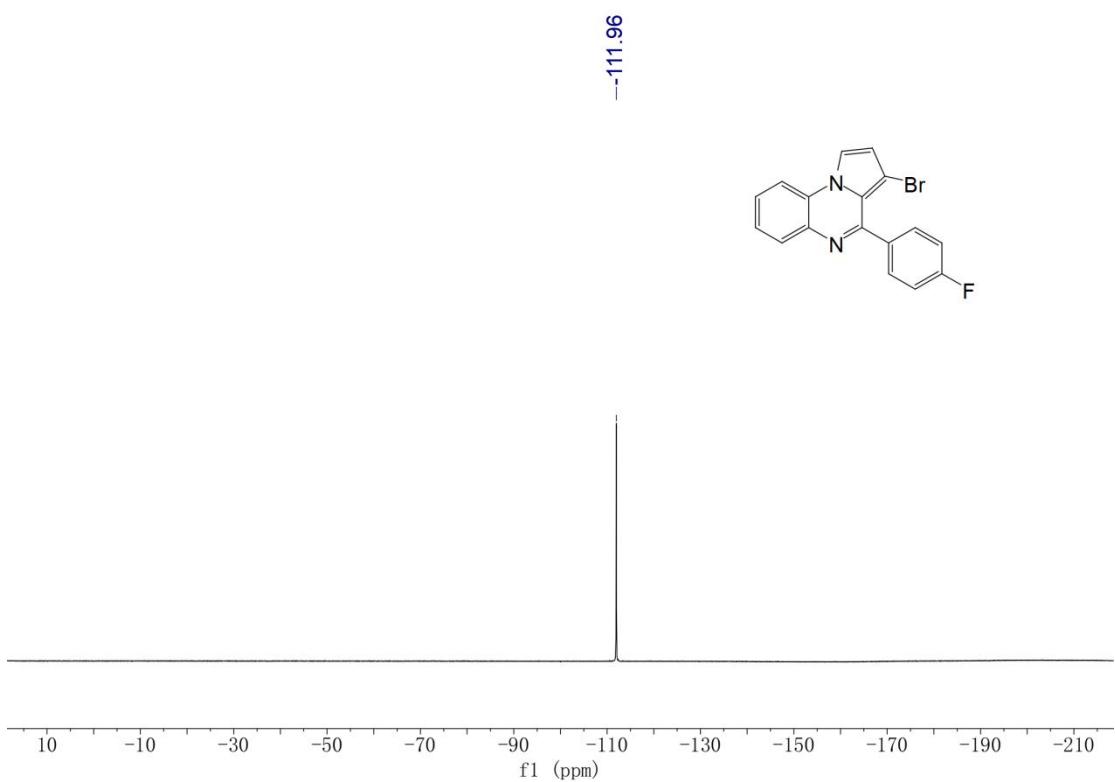


3-bromo-4-(4-bromophenyl)pyrrolo[1,2-*a*]quinoxaline [3f]

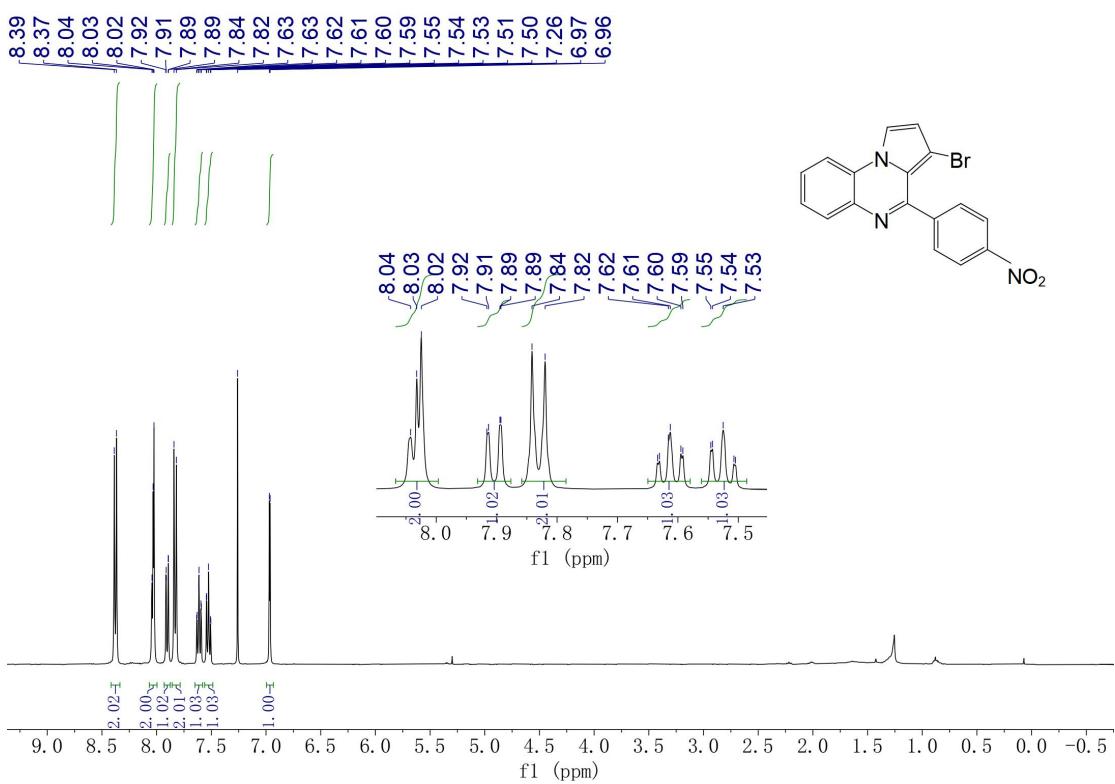


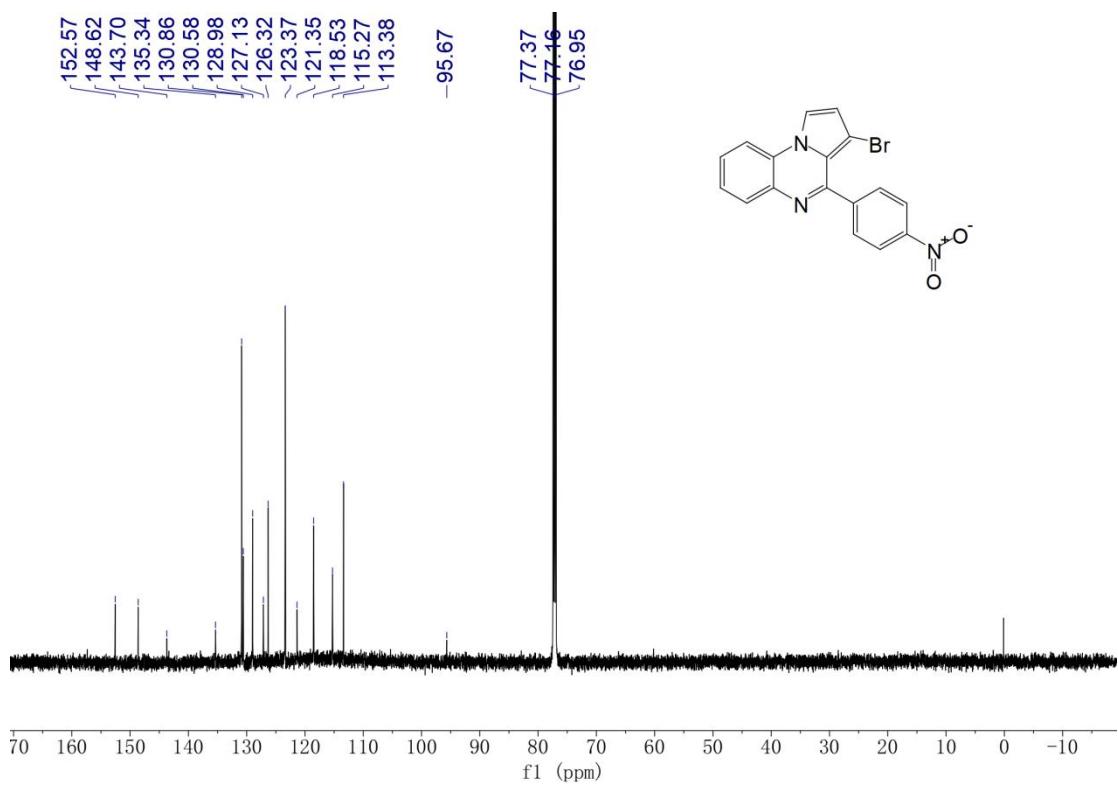
3-bromo-4-(4-fluorophenyl)pyrrolo[1,2-*a*]quinoxaline [3g]



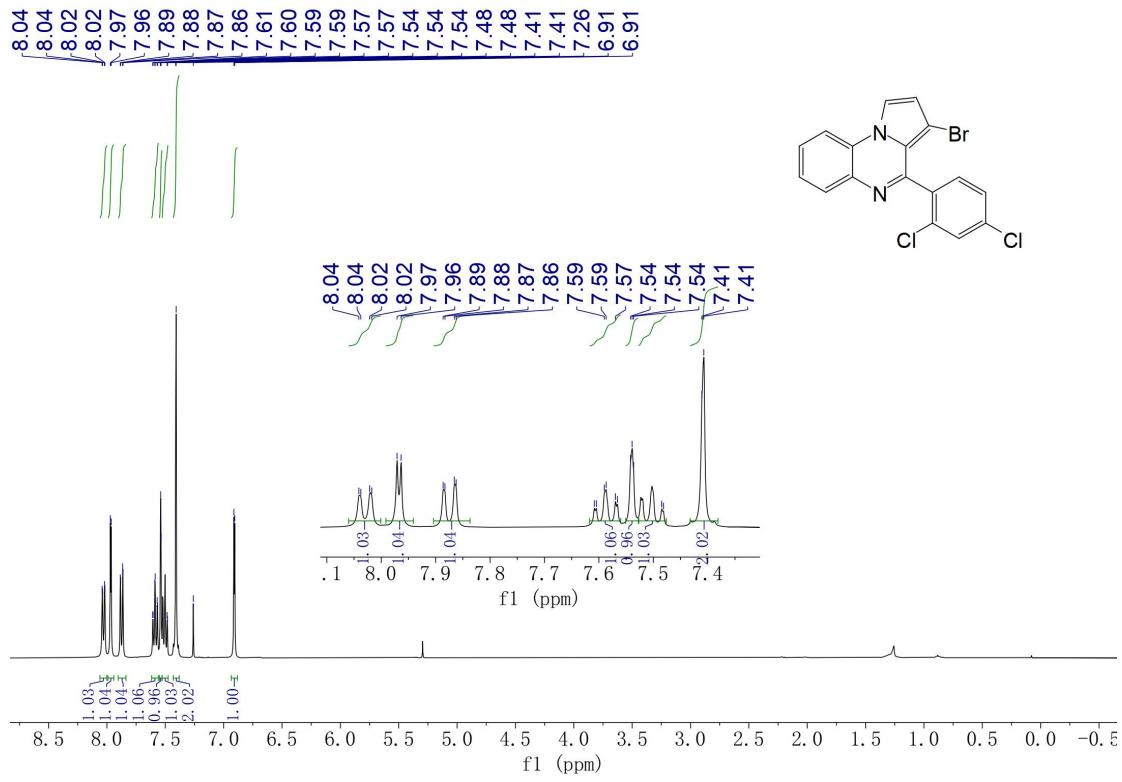


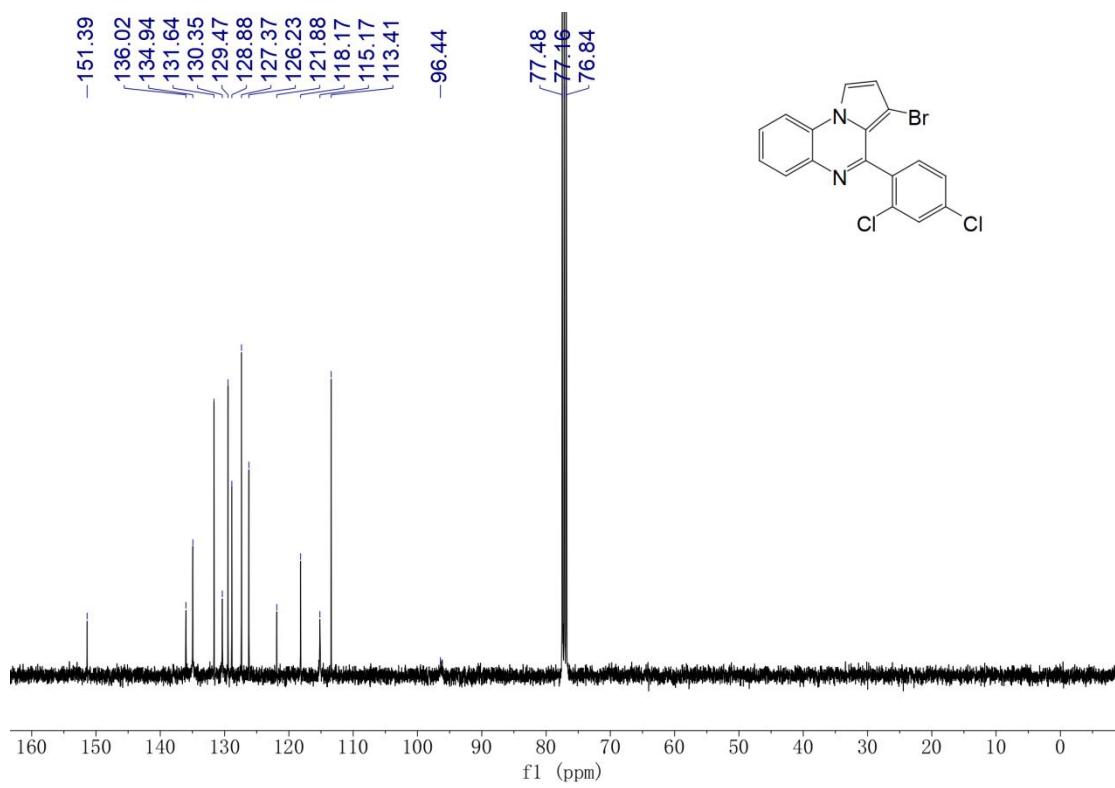
3-bromo-4-(4-nitrophenyl)pyrrolo[1,2-a]quinoxaline [3h]



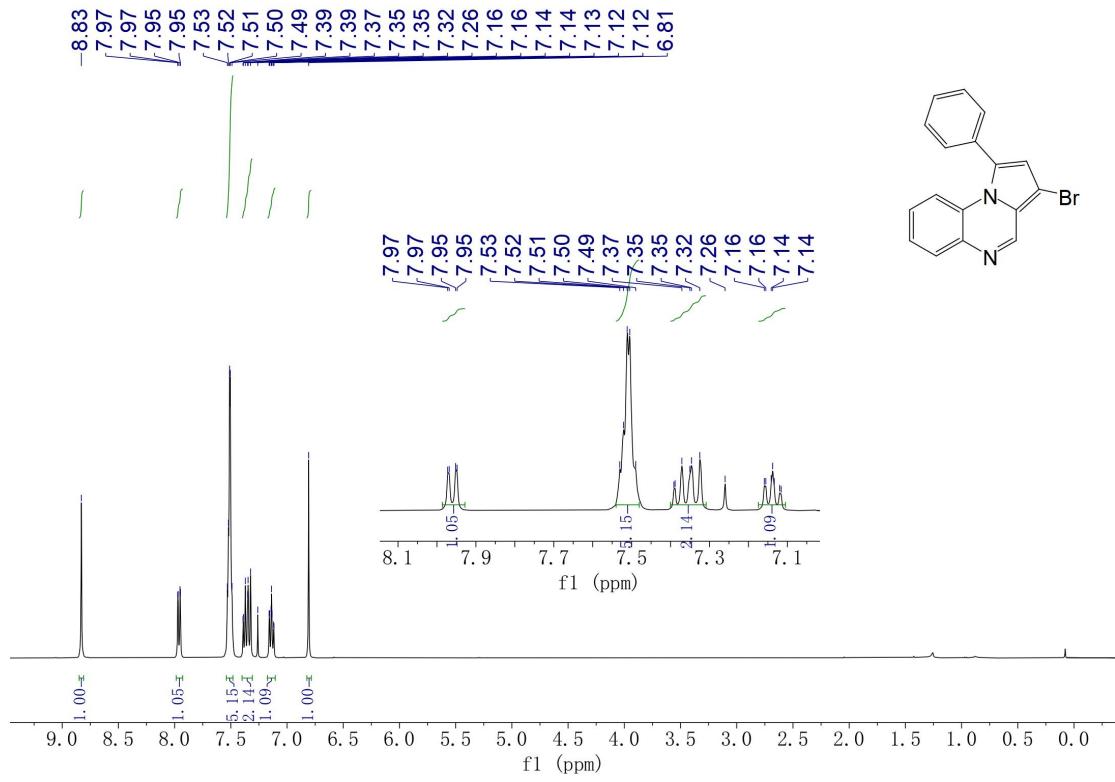


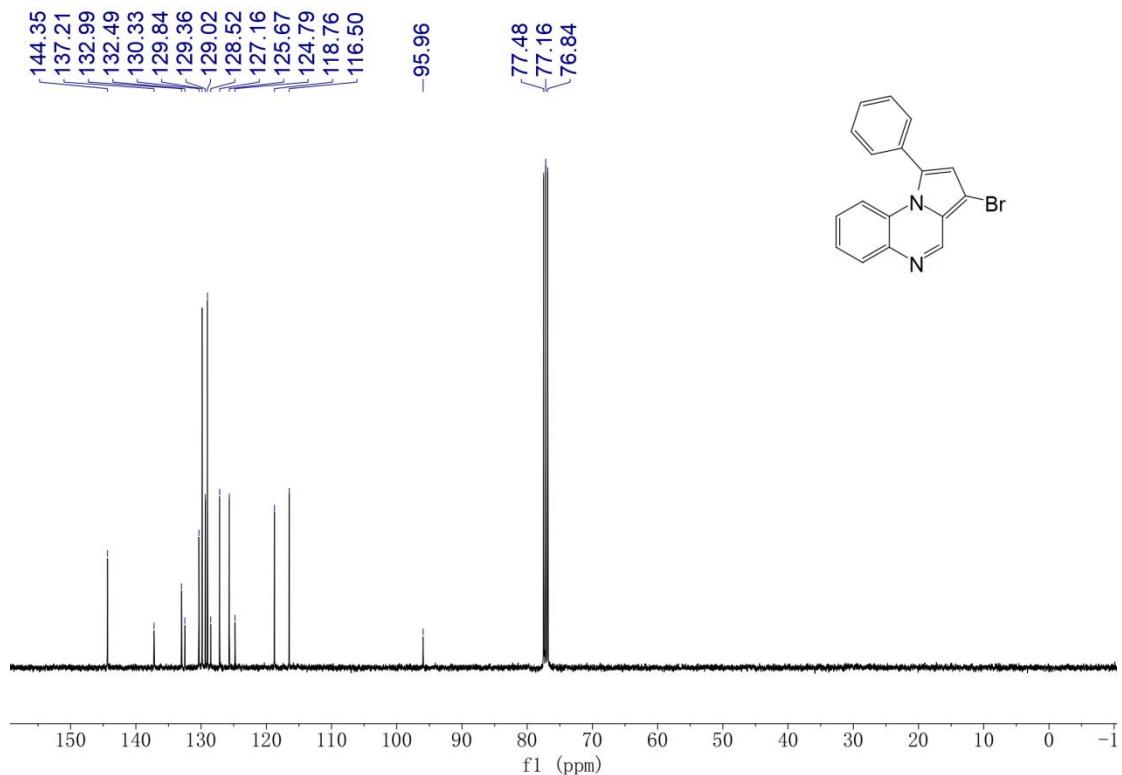
3-bromo-4-(2,4-dichlorophenyl)pyrrolo[1,2-a]quinoxaline [3i]



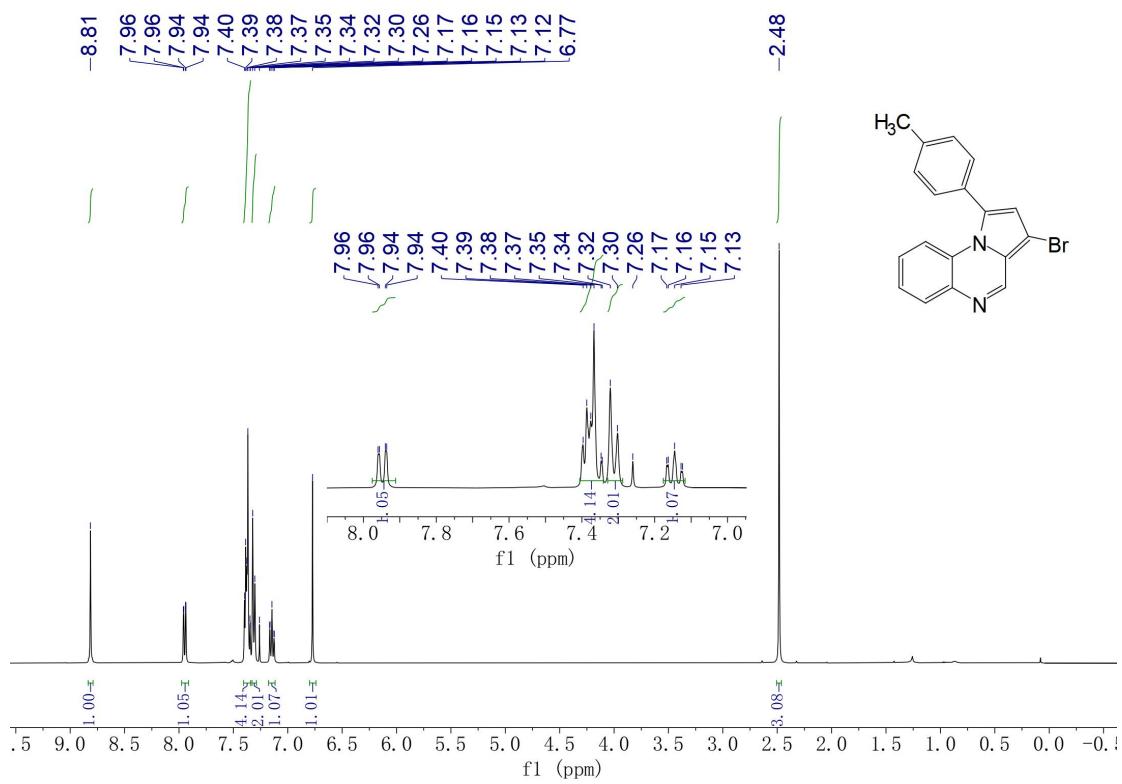


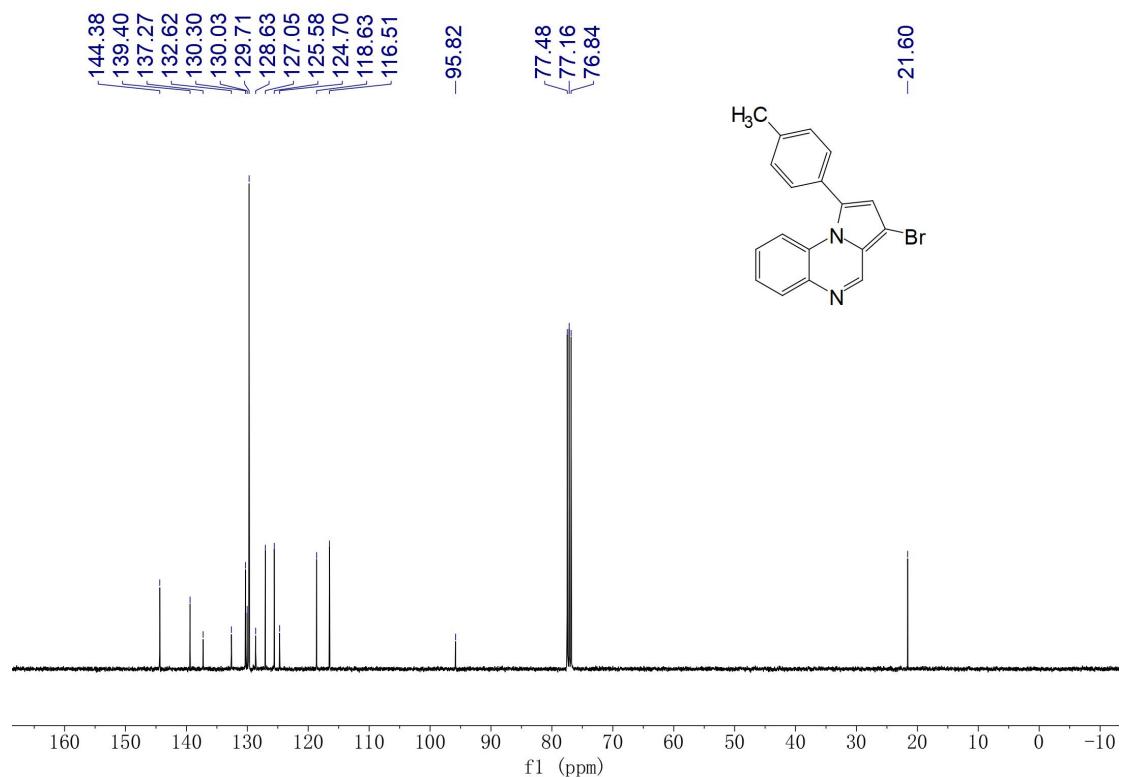
3-bromo-1-phenylpyrrolo[1,2-*a*]quinoxaline [3j]



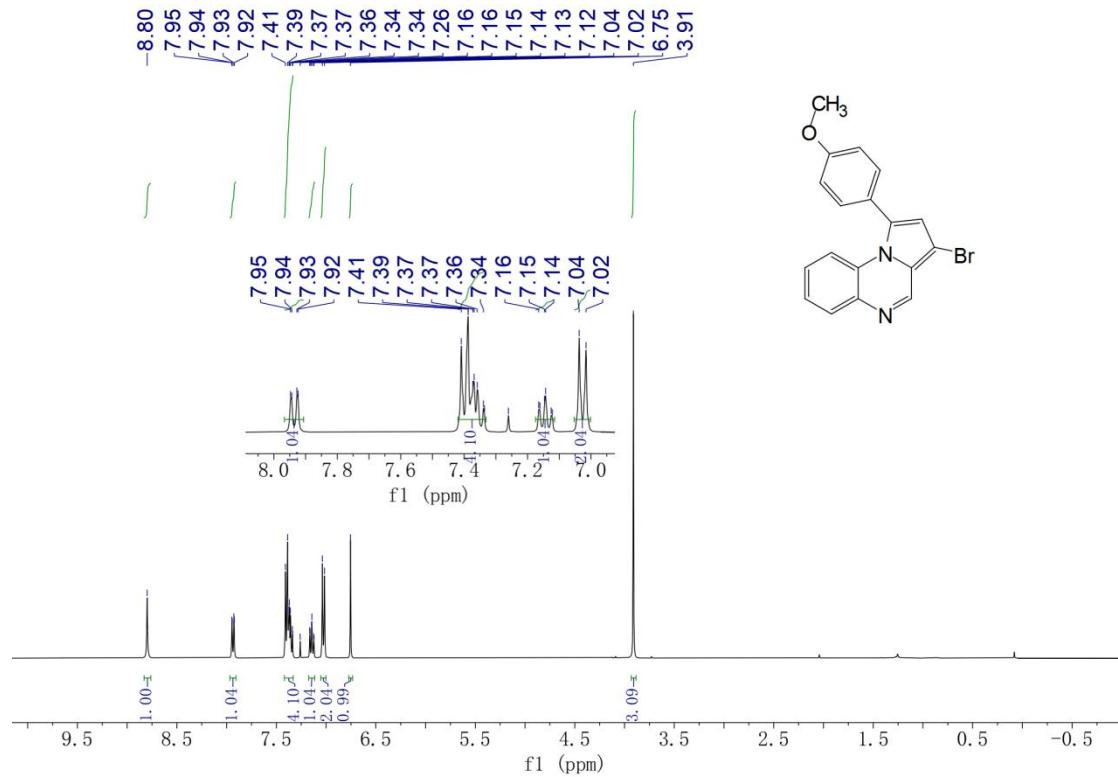


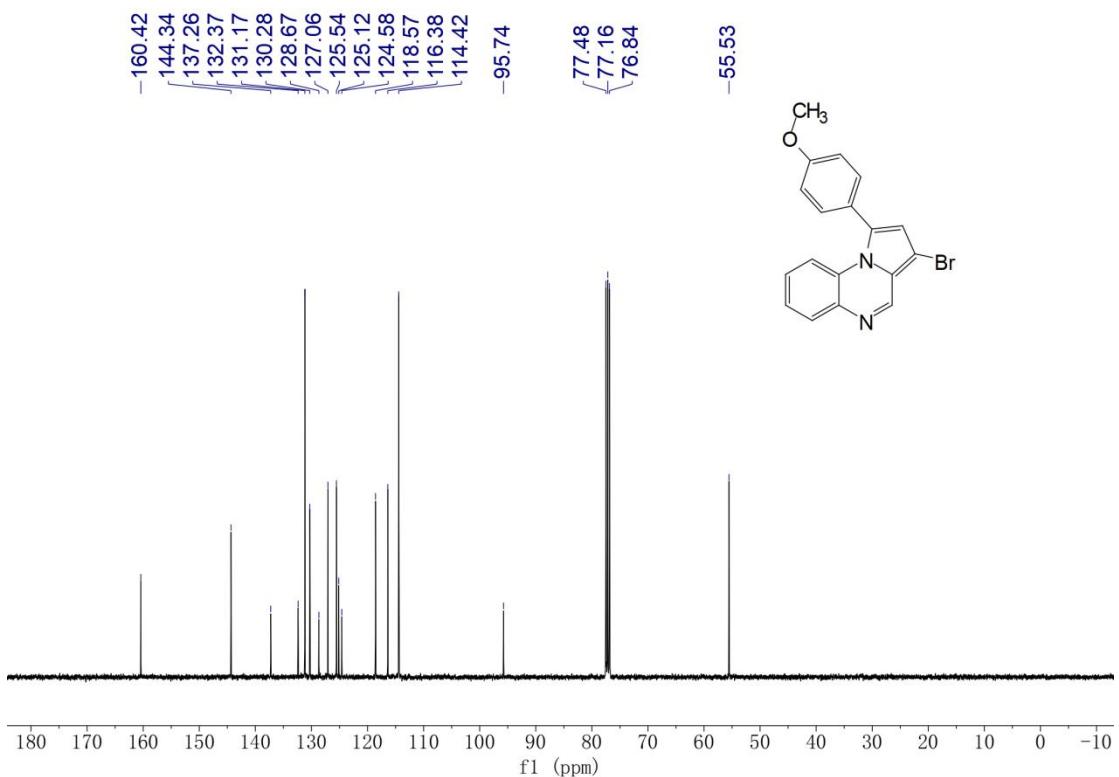
3-bromo-1-(p-tolyl)pyrrolo[1,2-a]quinoxaline [3k]



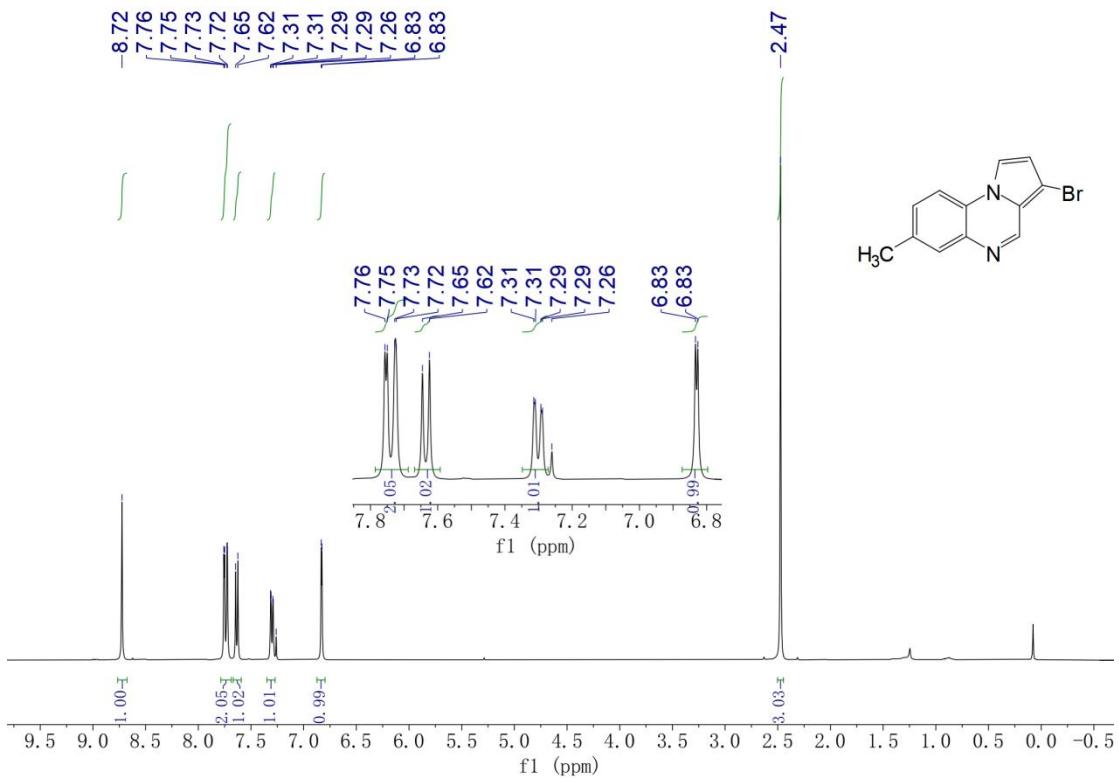


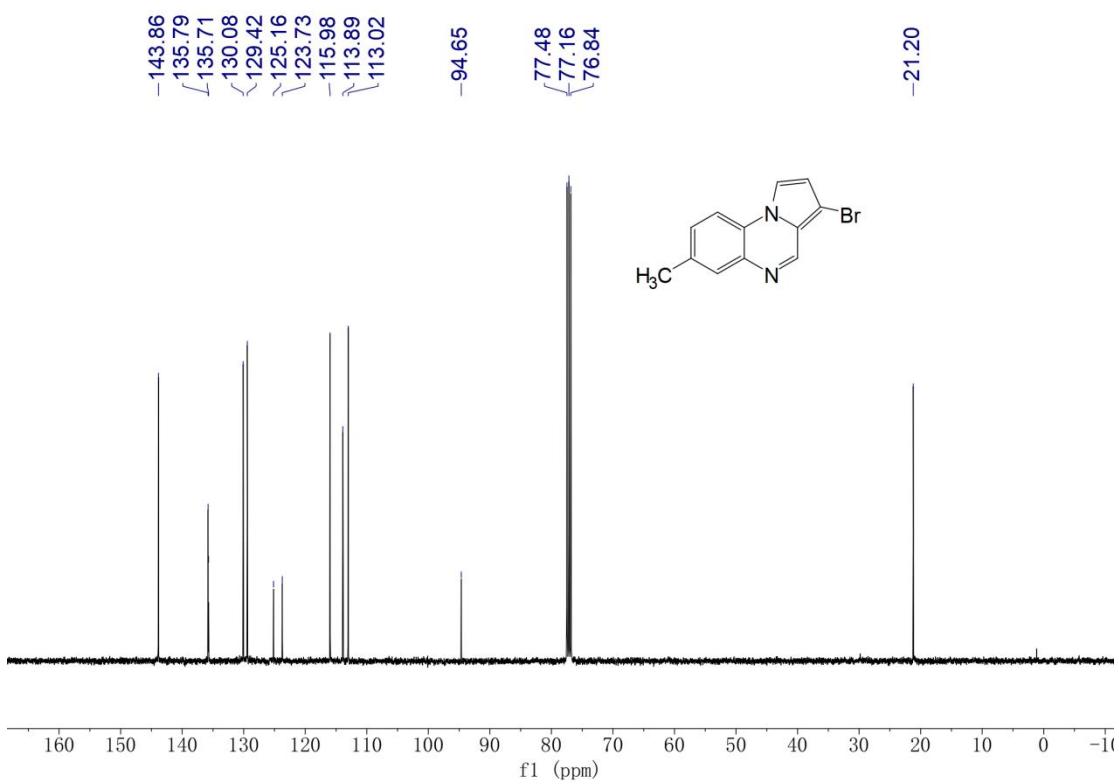
3-bromo-1-(4-methoxyphenyl)pyrrolo[1,2-a]quinoxaline [3l]



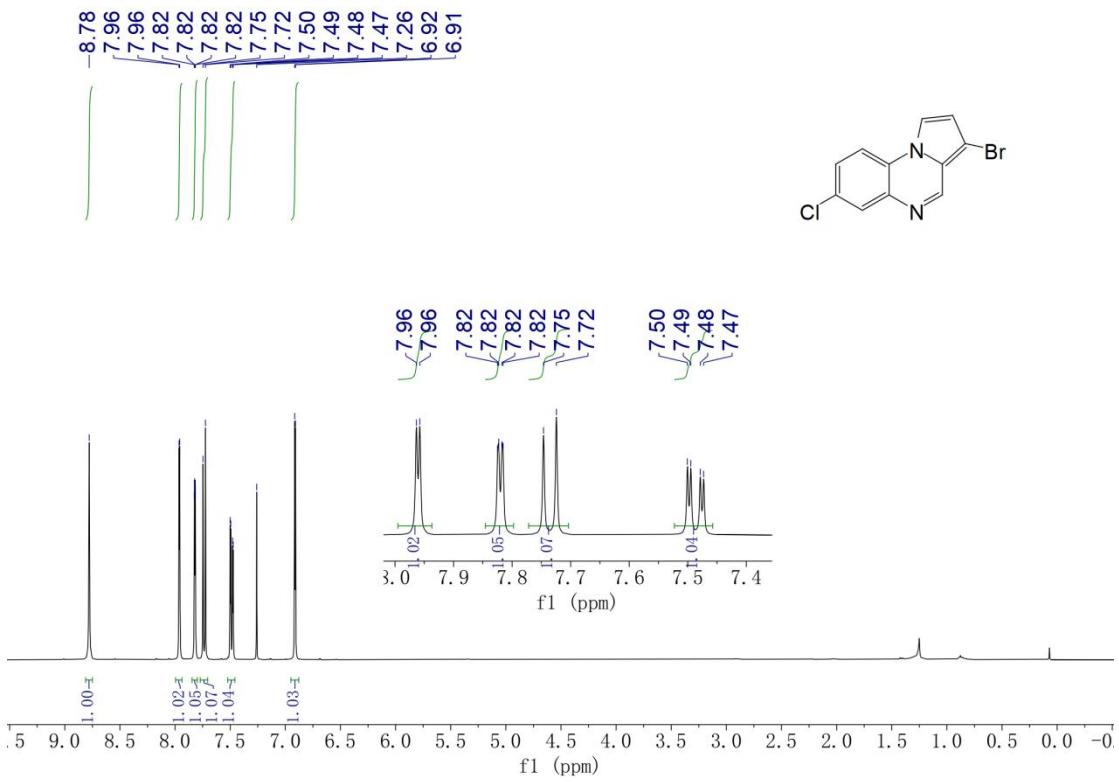


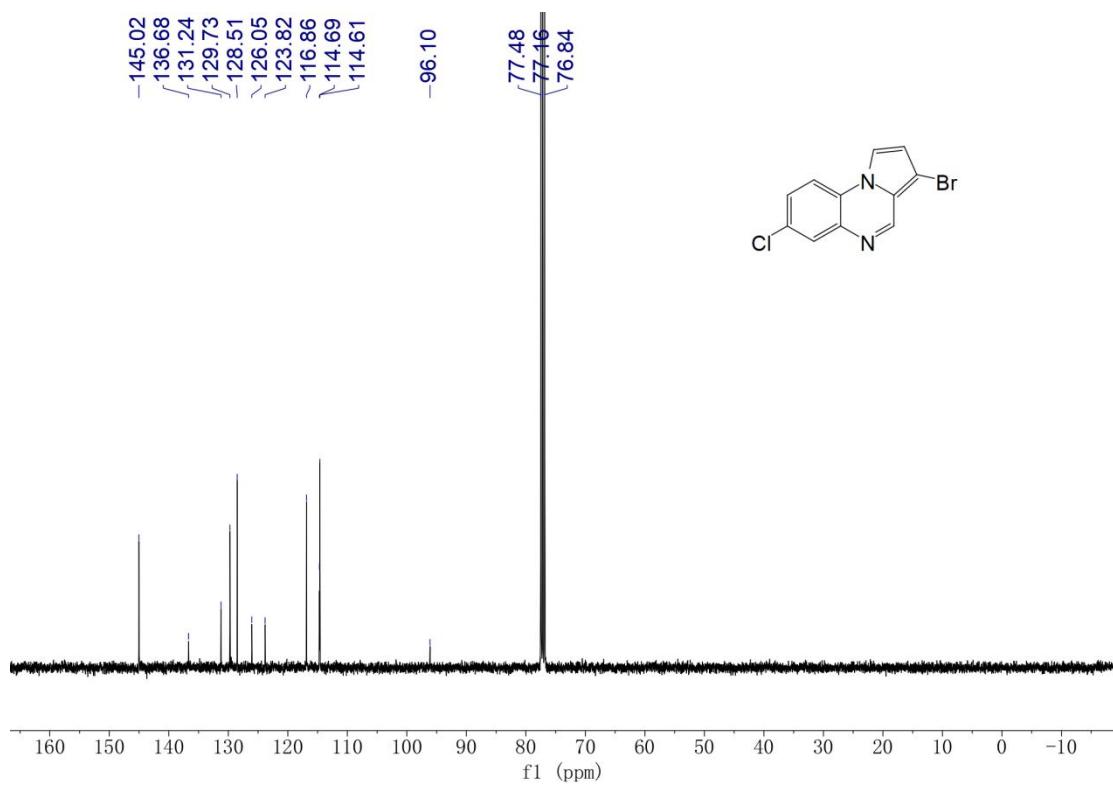
3-bromo-7-methylpyrrolo[1,2-*a*]quinoxaline [3m]



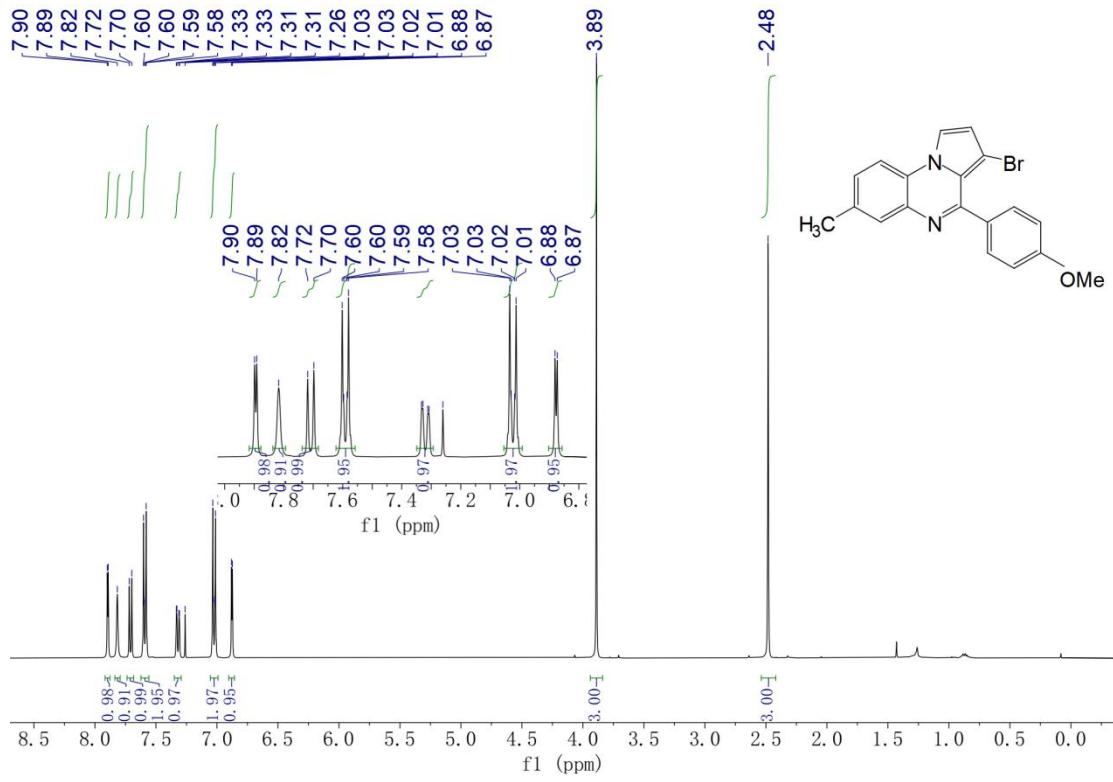


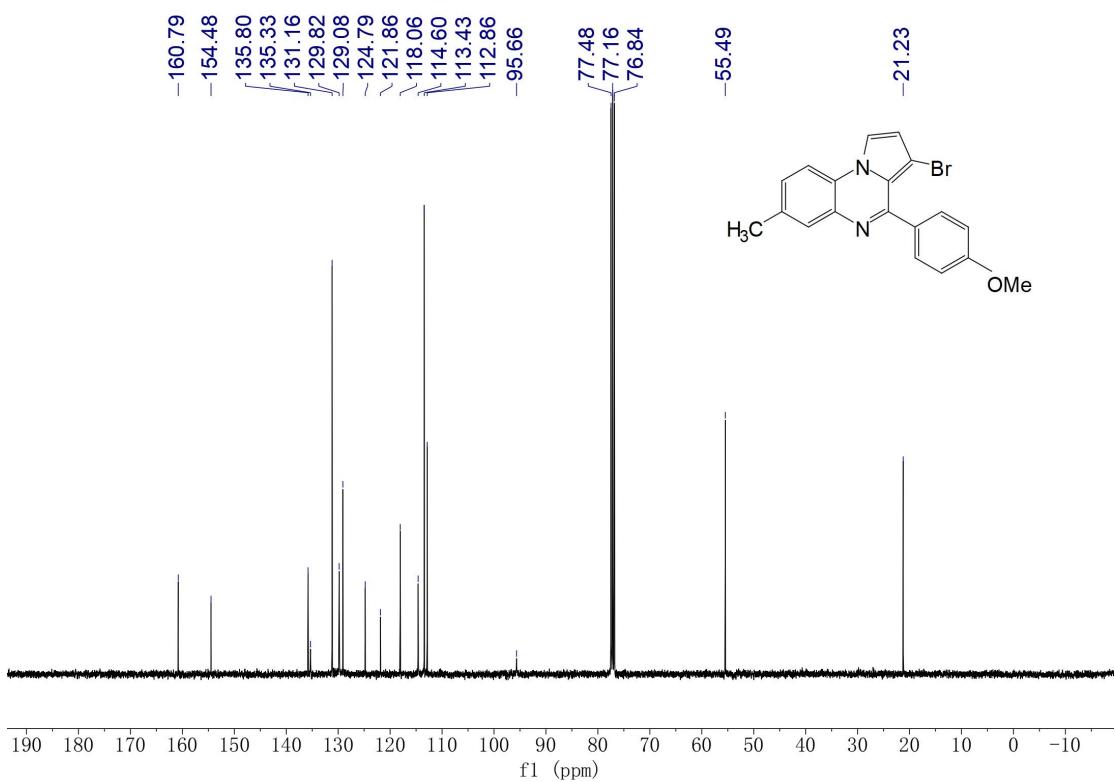
3-bromo-7-chloropyrrolo[1,2-*a*]quinoxaline [3n]



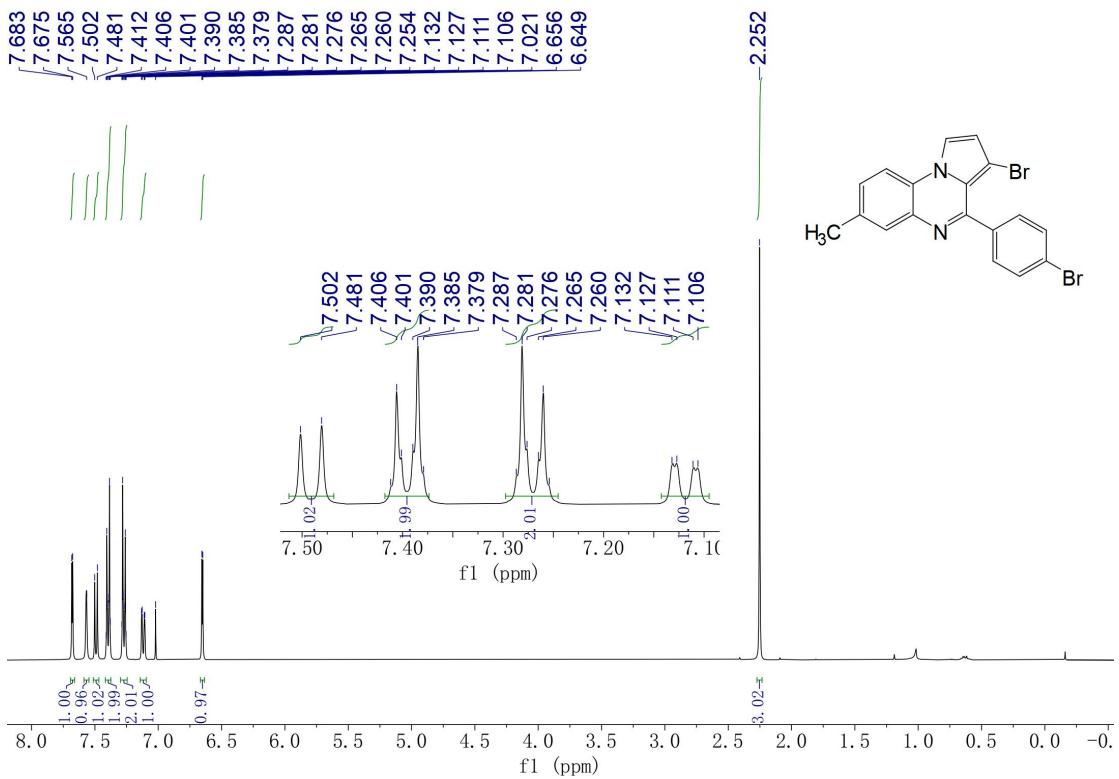


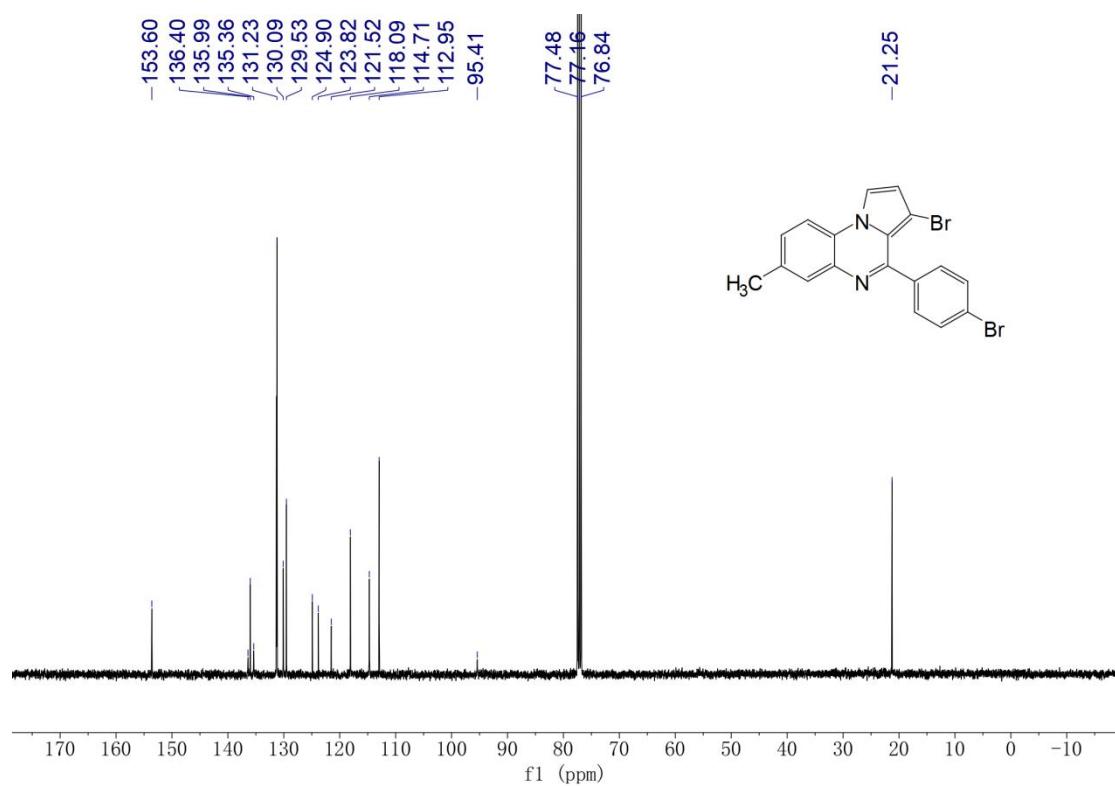
3-bromo-4-(4-methoxyphenyl)-7-methylpyrrolo[1,2-*a*]quinoxaline [3o]



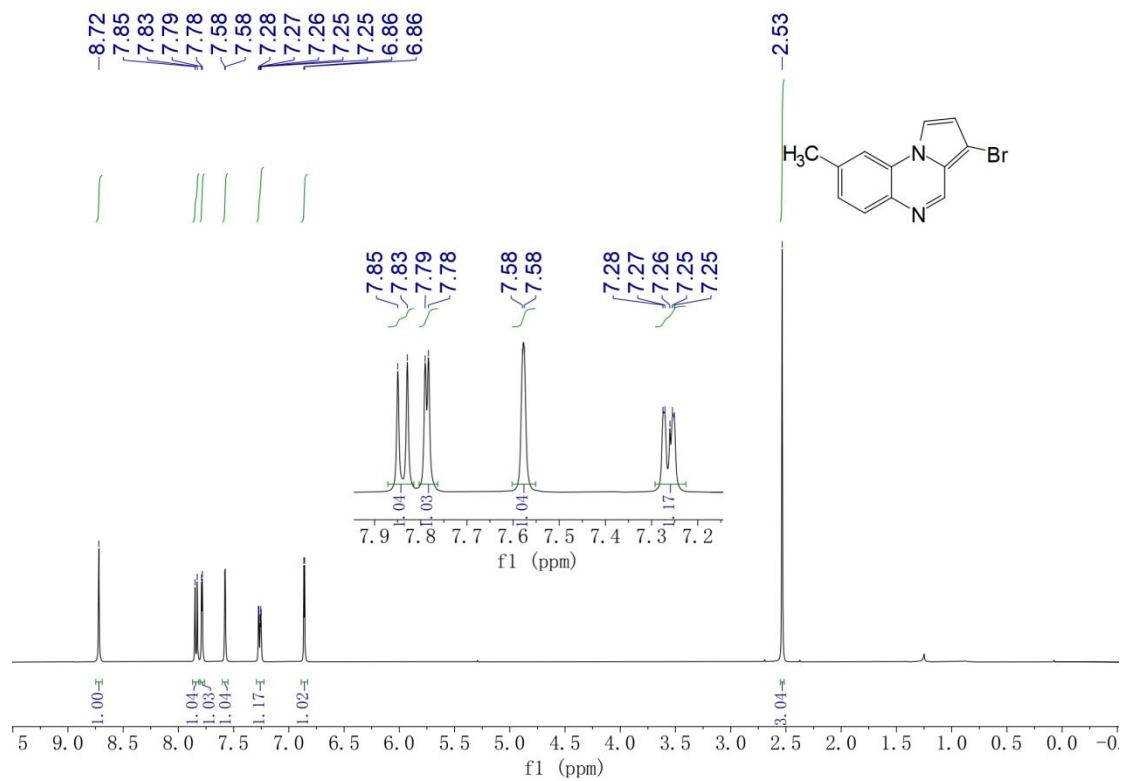


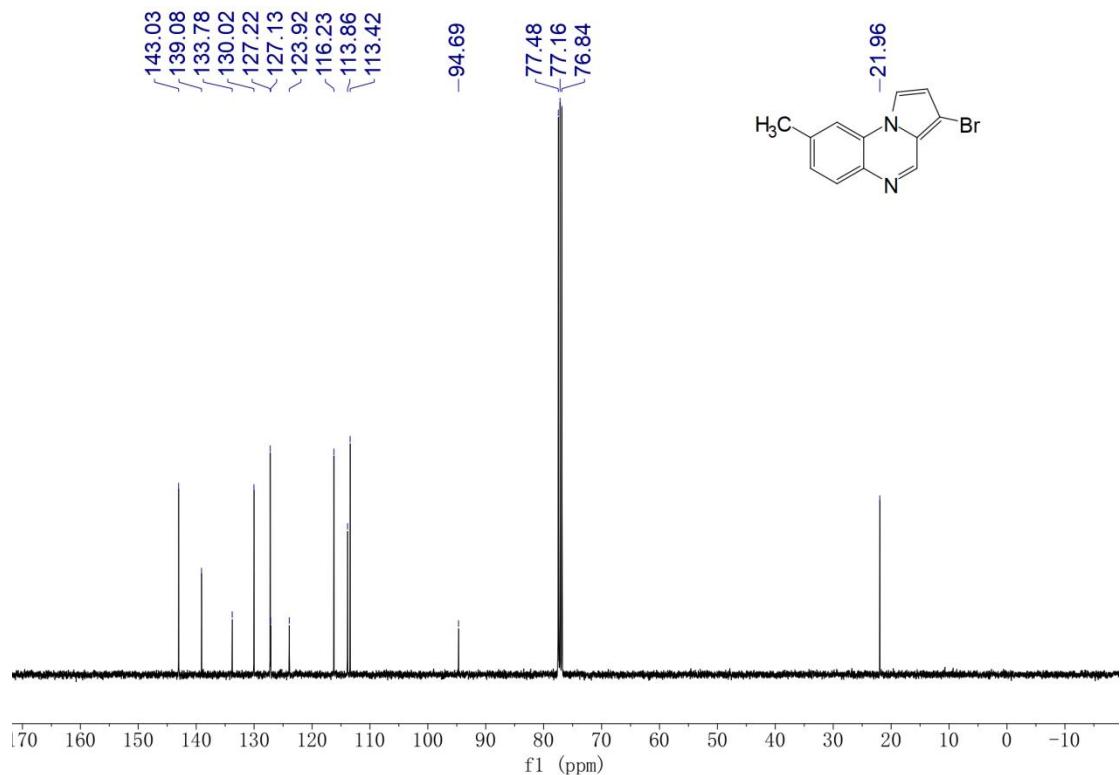
3-bromo-4-(4-bromophenyl)-7-methylpyrrolo[1,2-a]quinoxaline [3p]



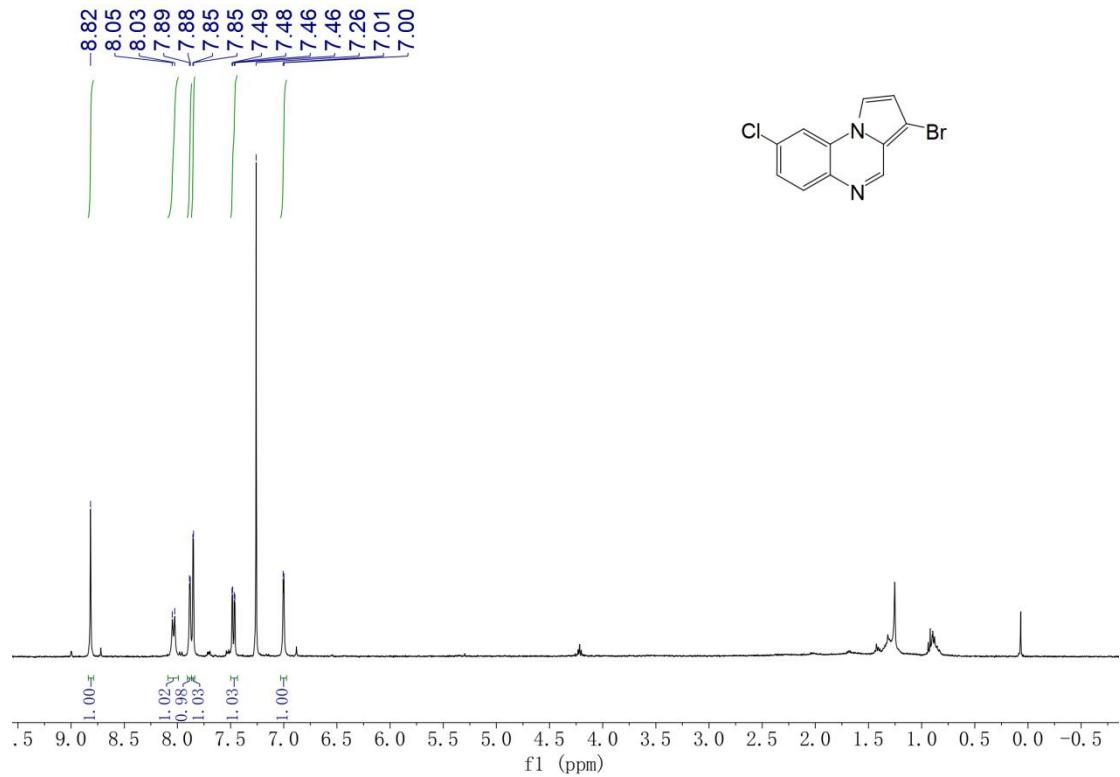


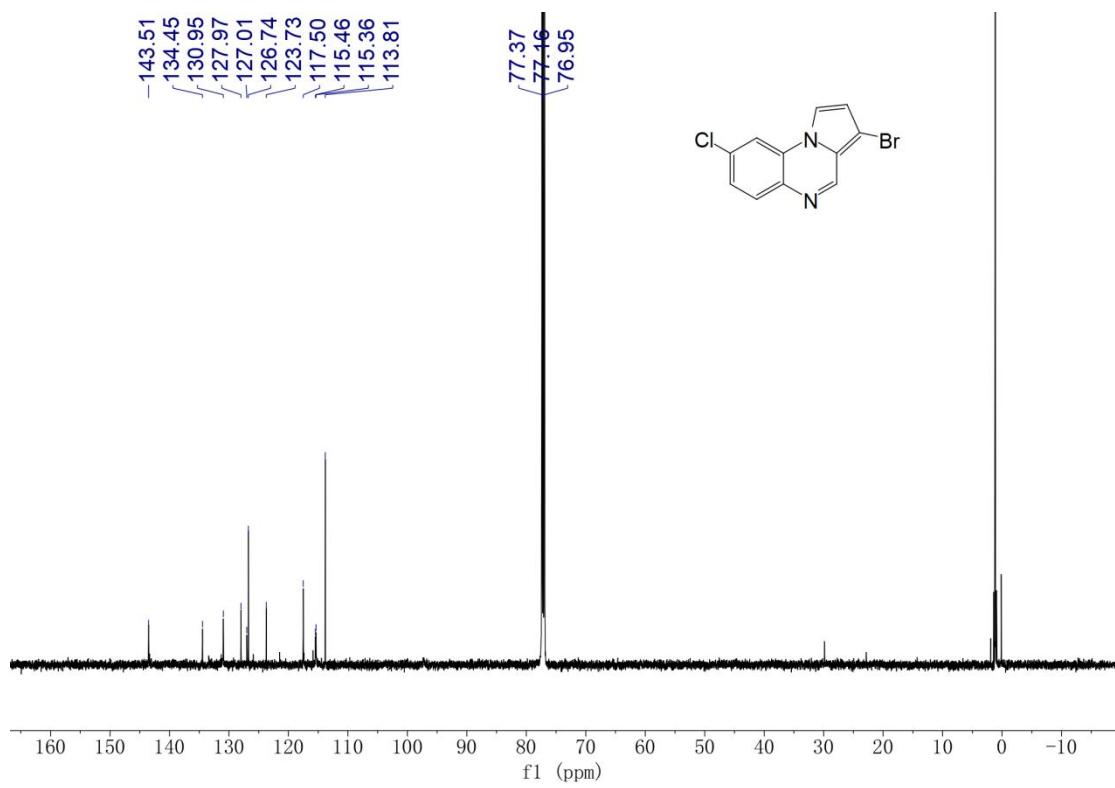
3-bromo-8-methylpyrrolo[1,2-*a*]quinoxaline [3q]



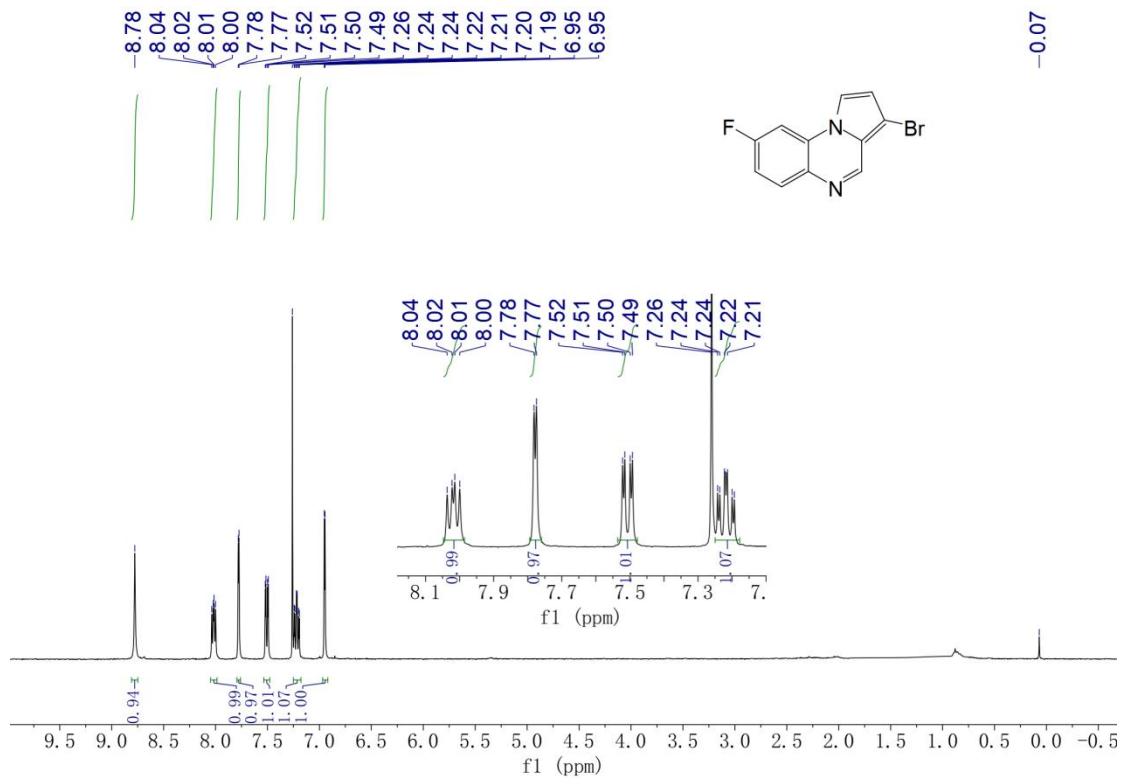


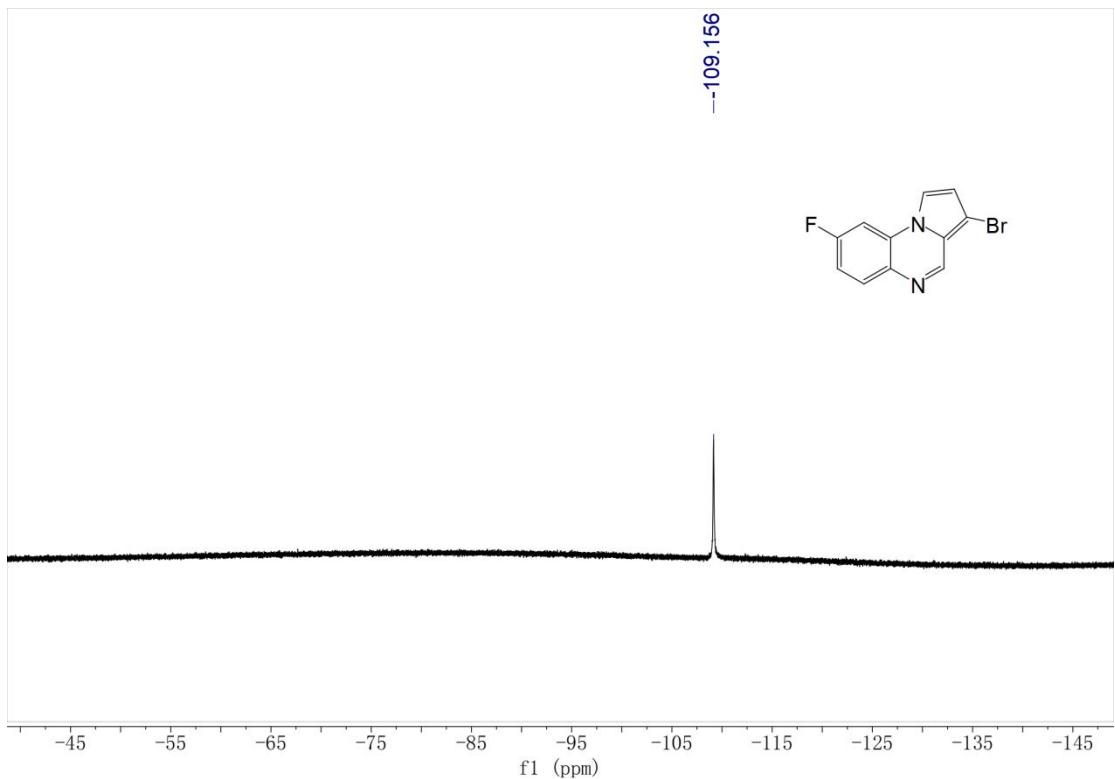
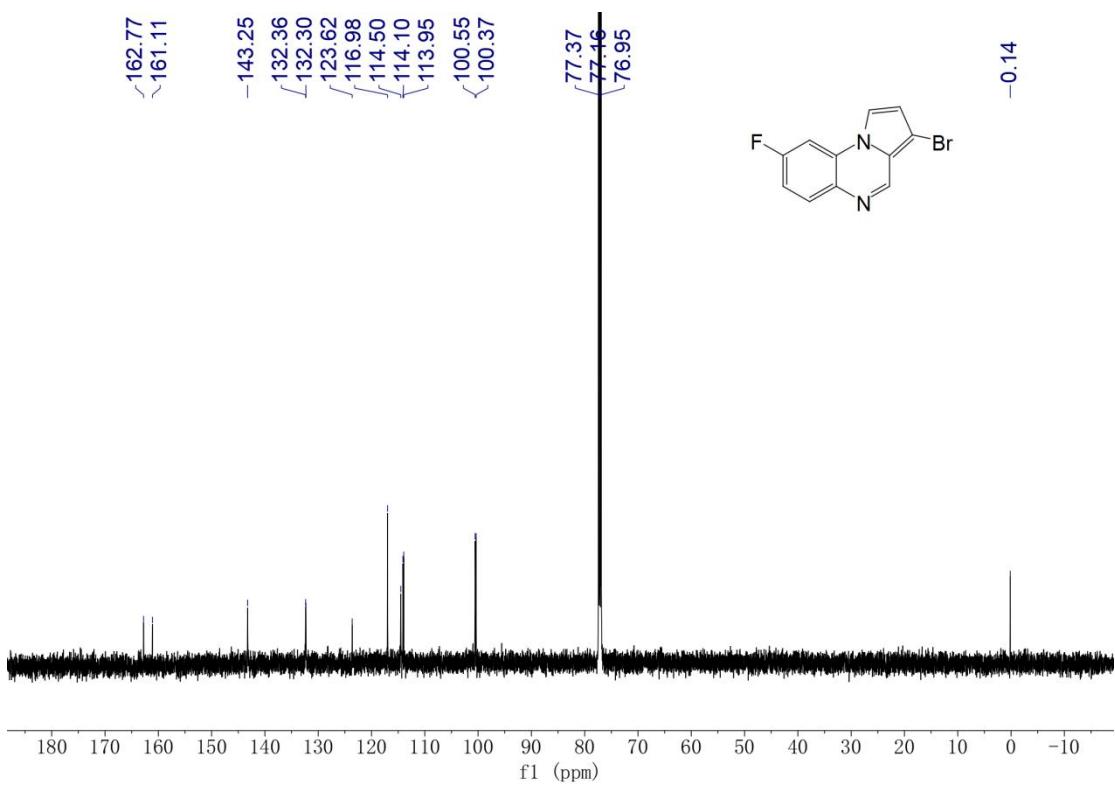
3-bromo-8-chloropyrrolo[1,2-a]quinoxaline [3r]



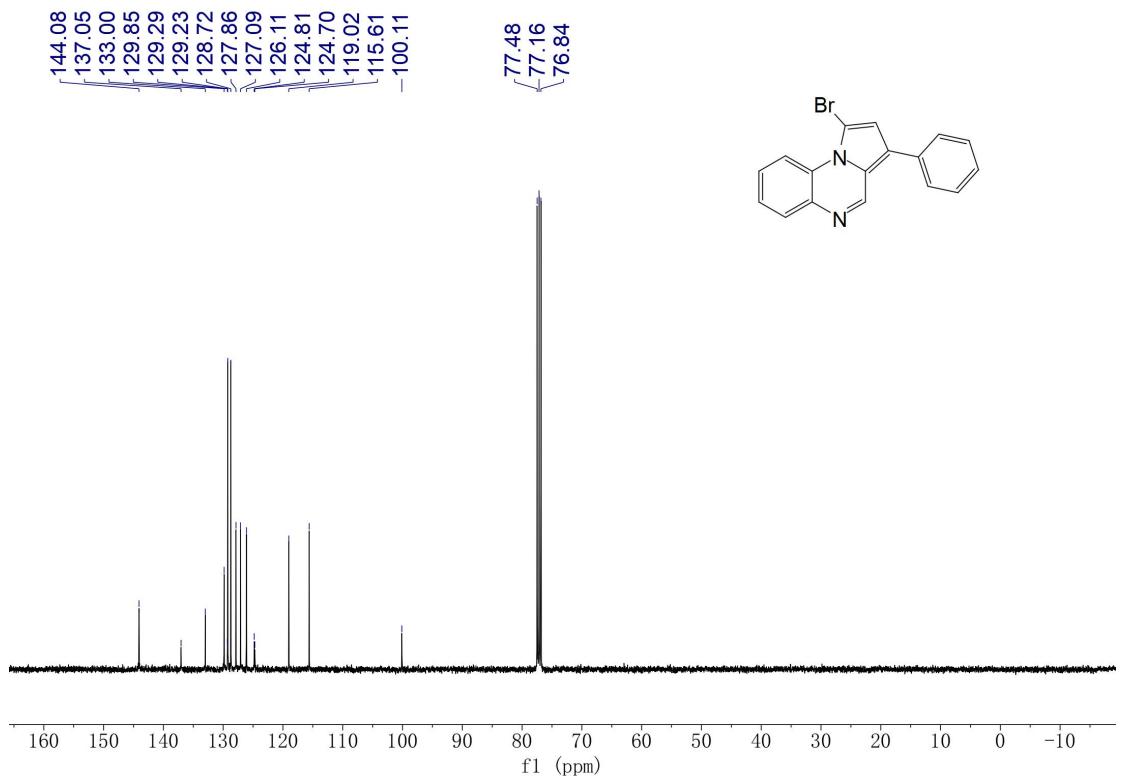
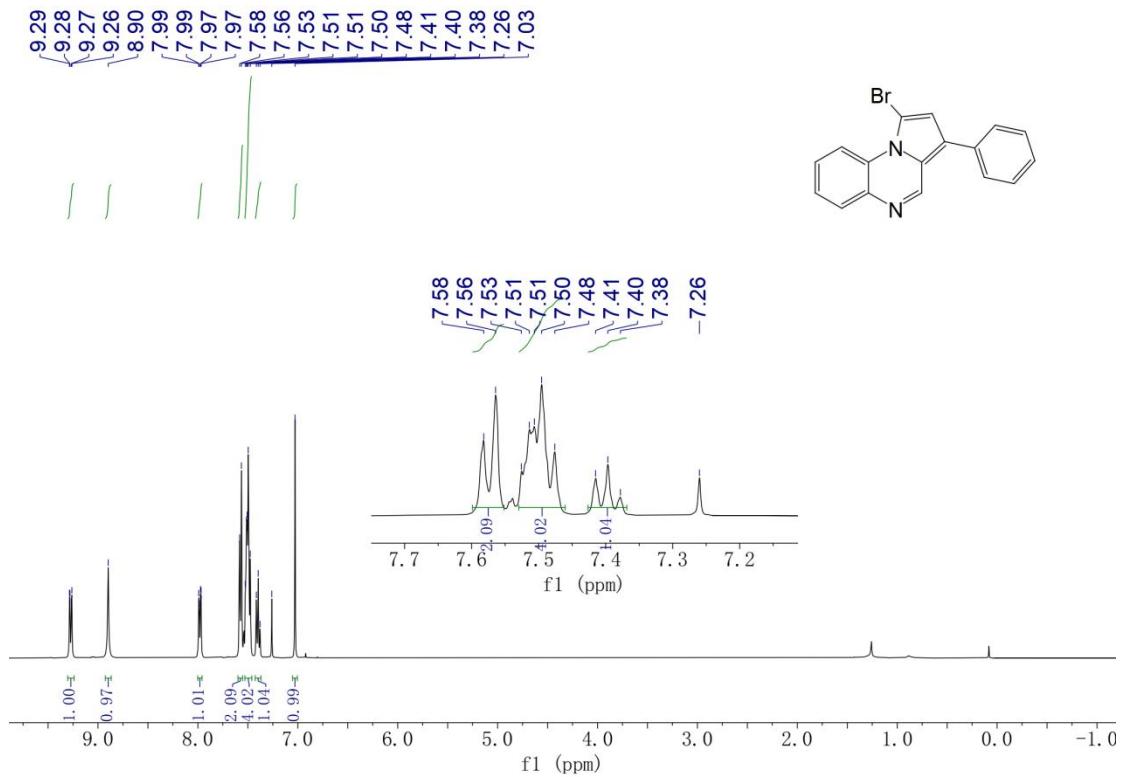


3-bromo-8-fluoropyrrolo[1,2-*a*]quinoxaline [3s]

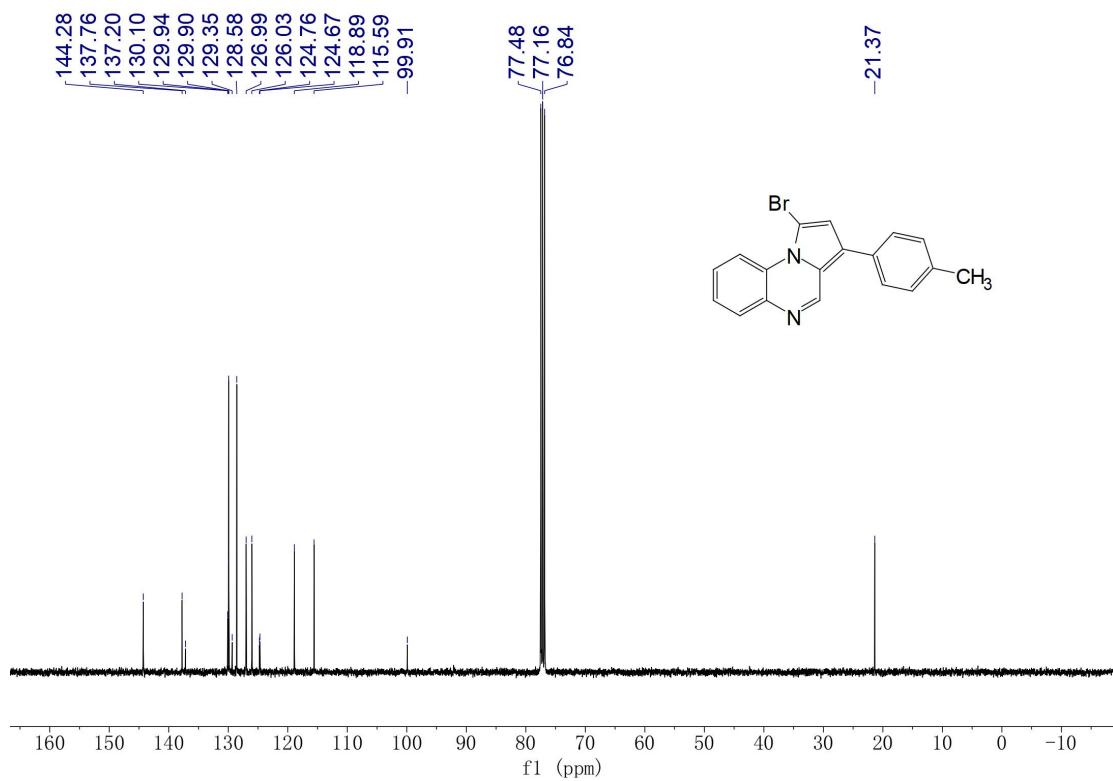
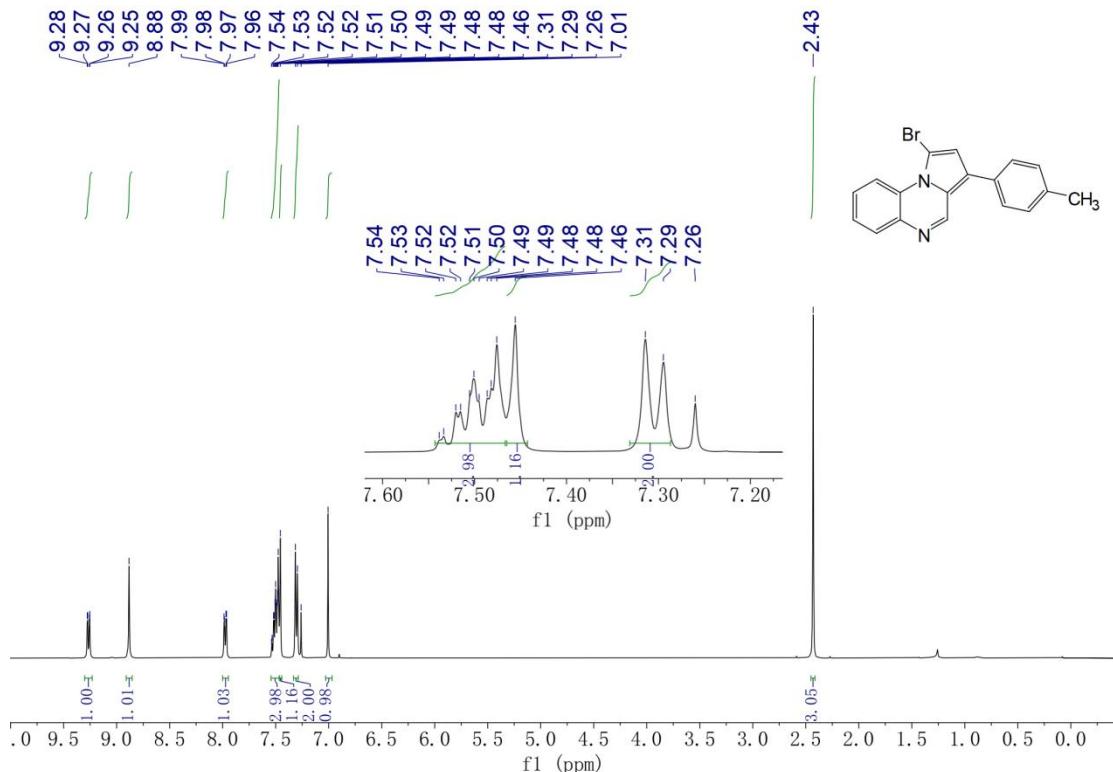




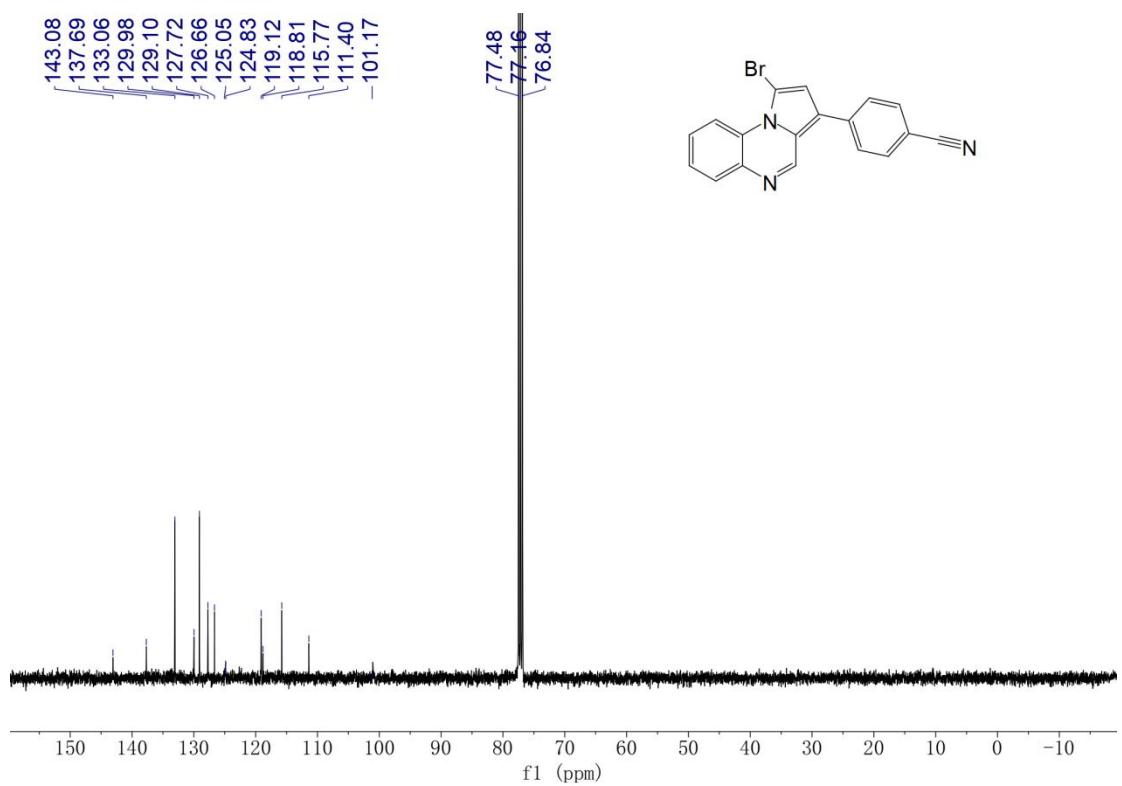
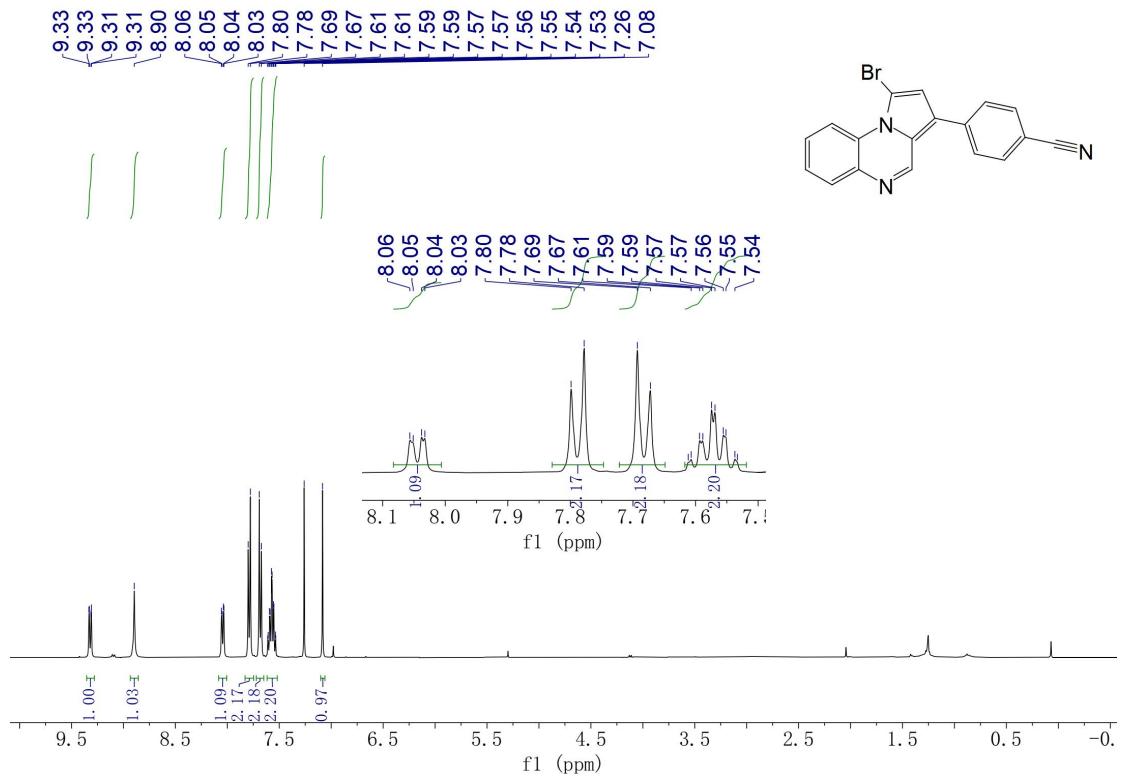
1-bromo-3-phenylpyrrolo[1,2-*a*]quinoxaline [3t]



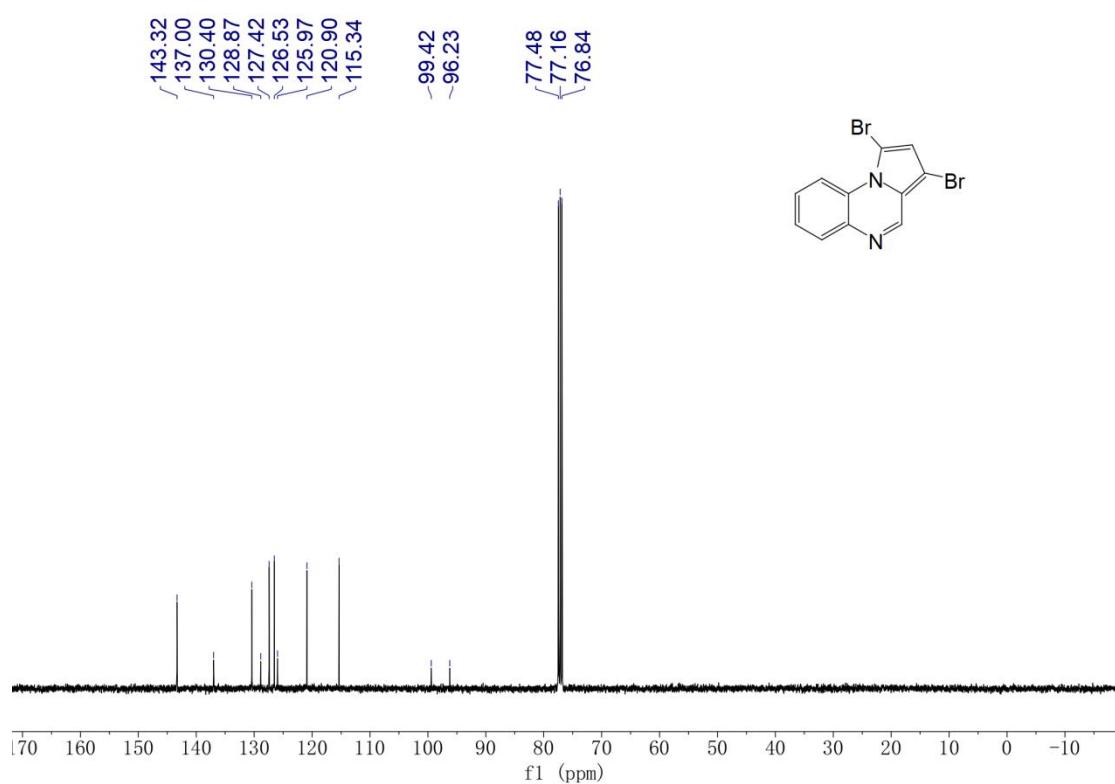
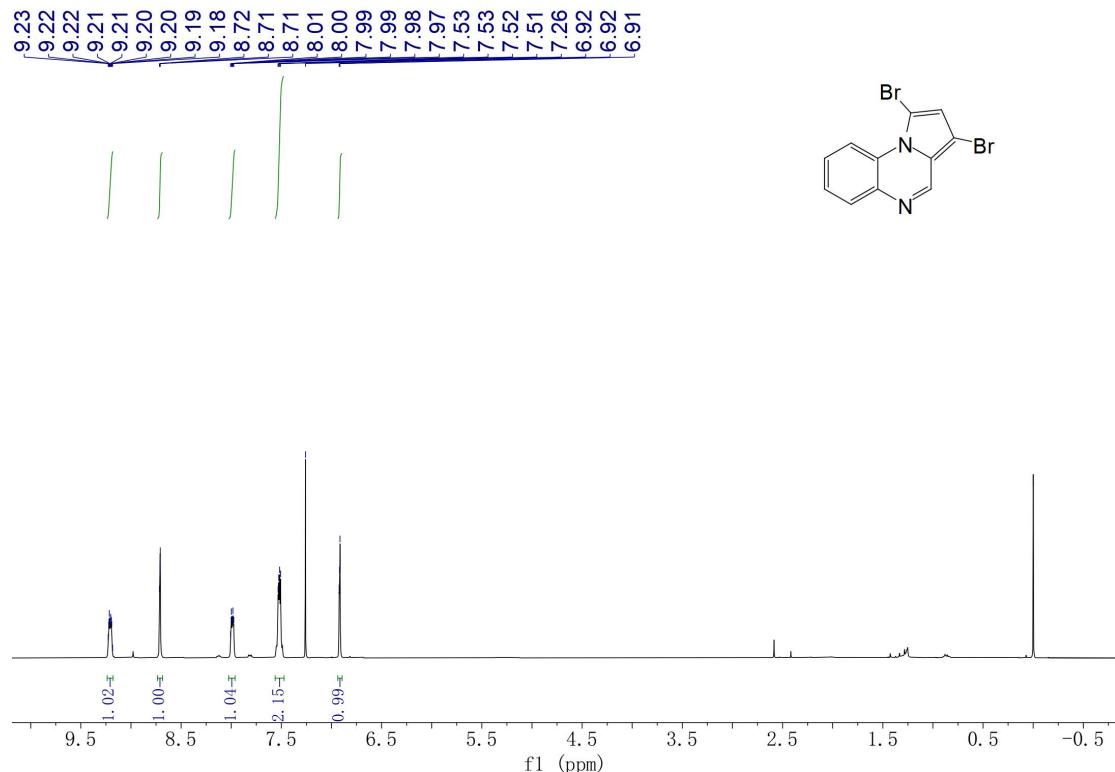
1-bromo-3-(p-tolyl)pyrrolo[1,2-*a*]quinoxaline [3u]



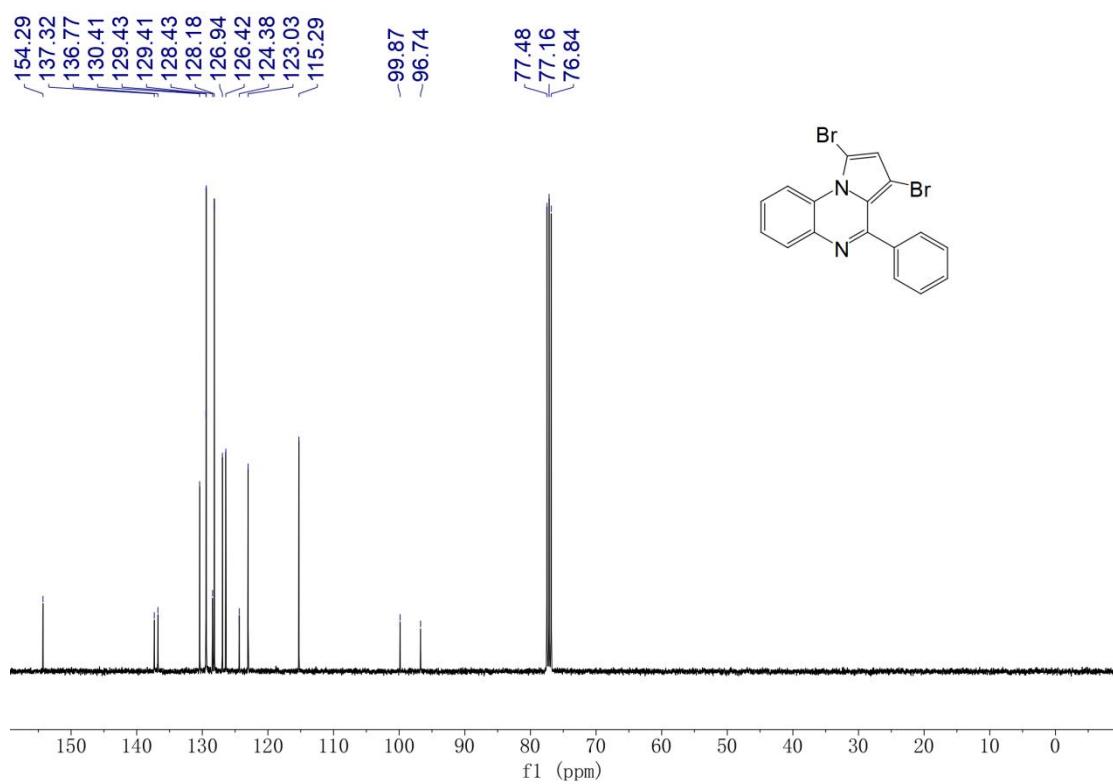
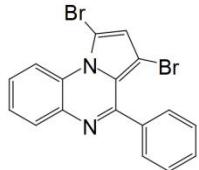
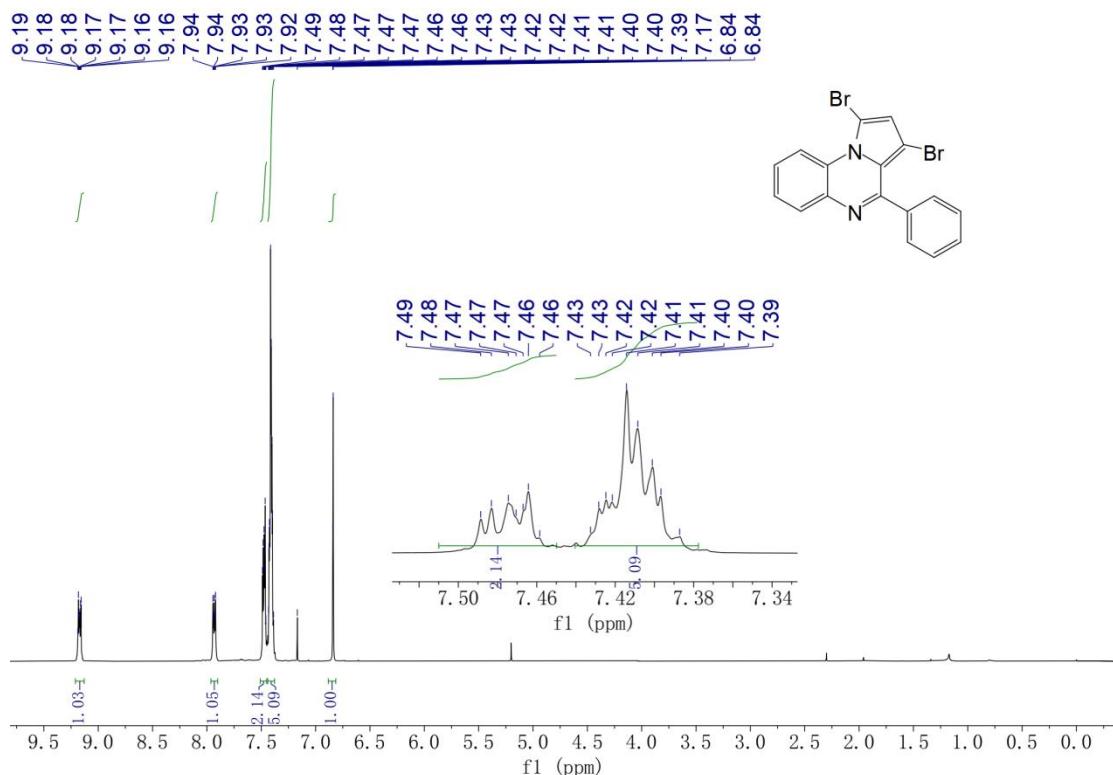
4-(1-bromopyrrolo[1,2-*a*]quinoxalin-3-yl)benzonitrile [3v]



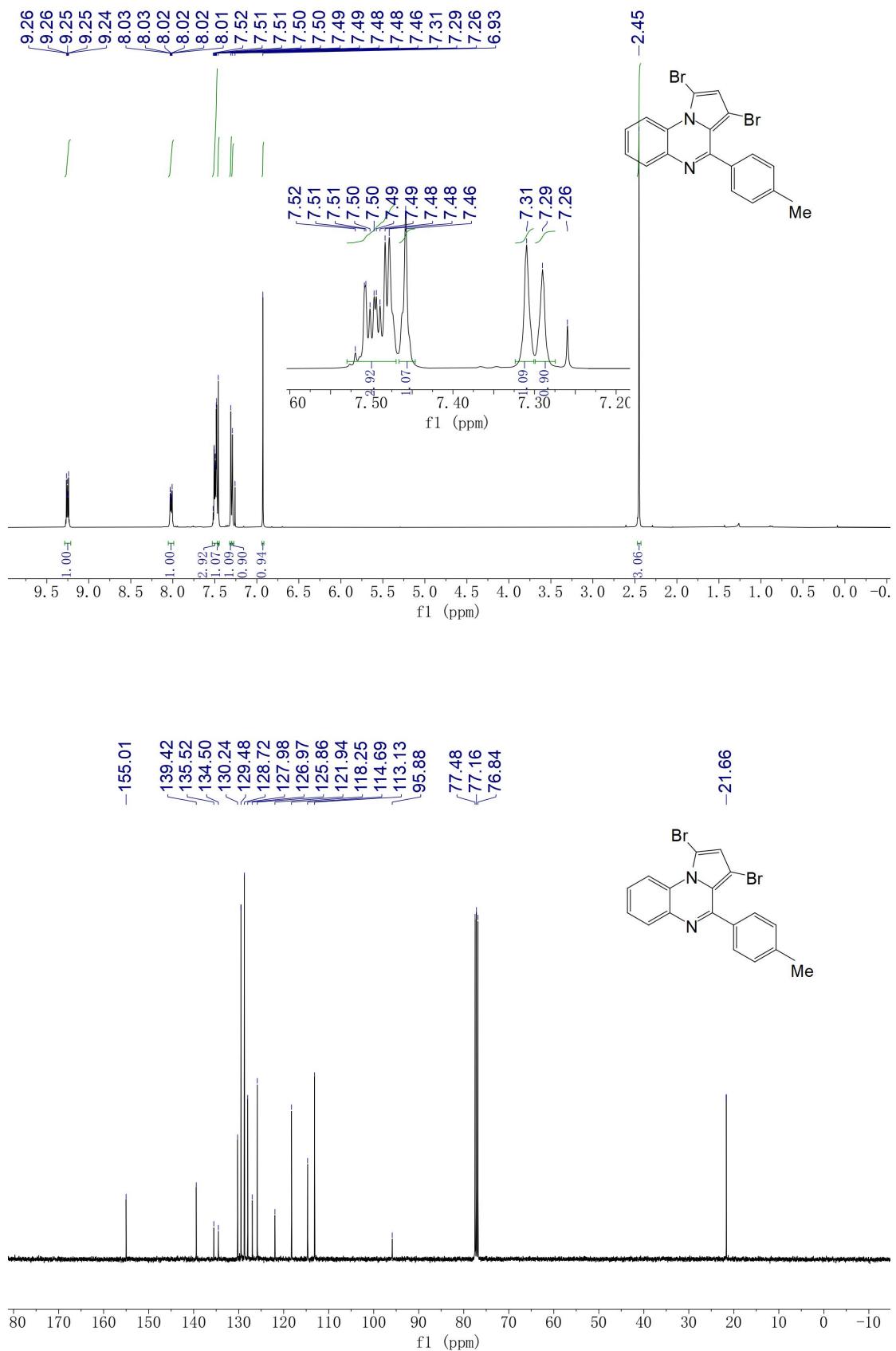
1,3-dibromopyrrolo[1,2-*a*]quinoxaline [4a]



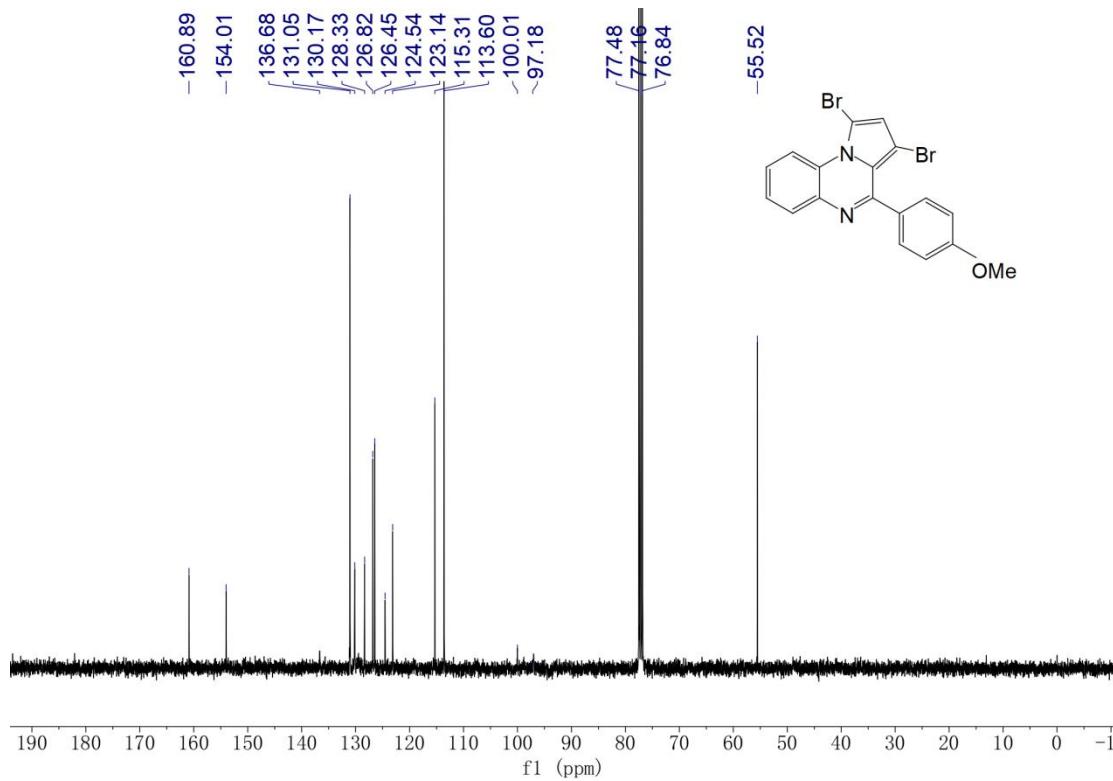
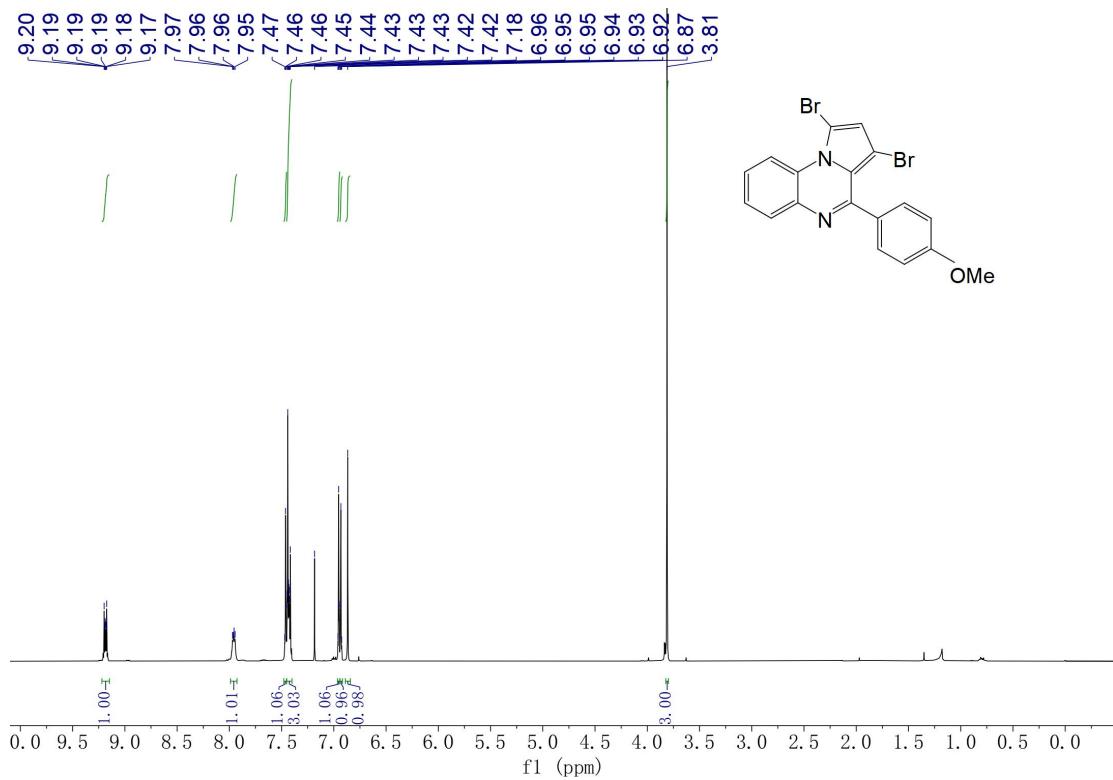
1,3-dibromo-4-phenylpyrrolo[1,2-a]quinoxaline [4b]



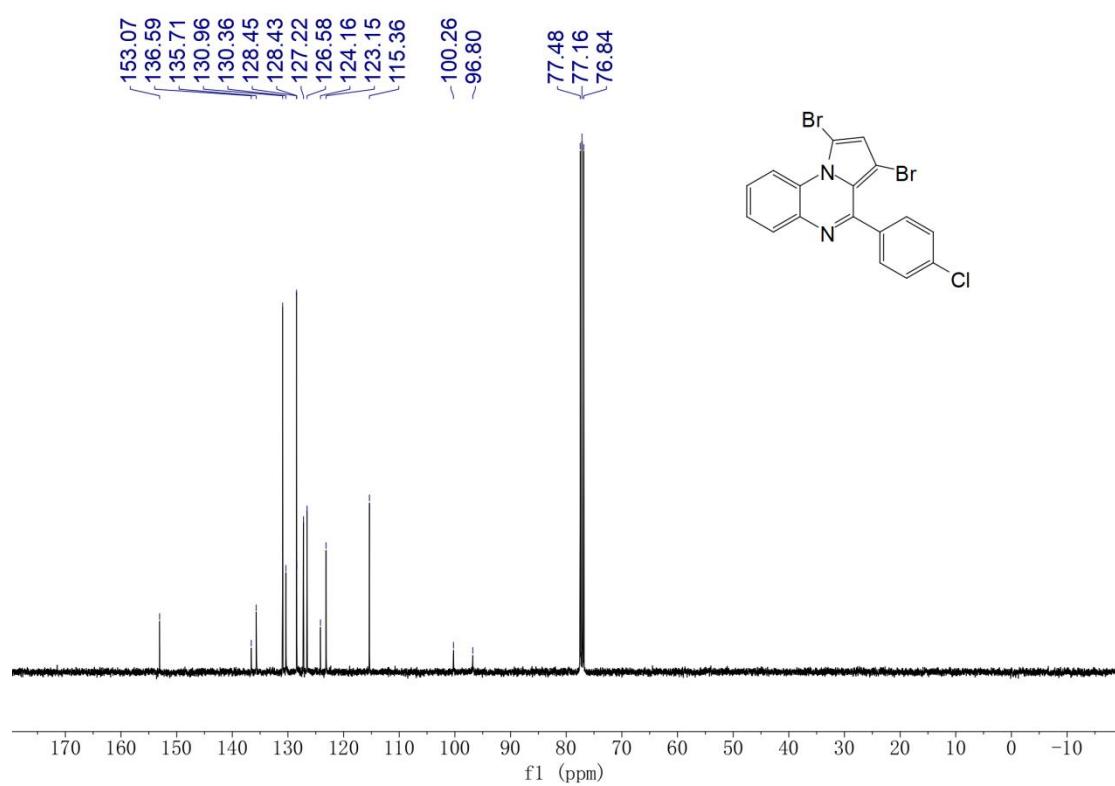
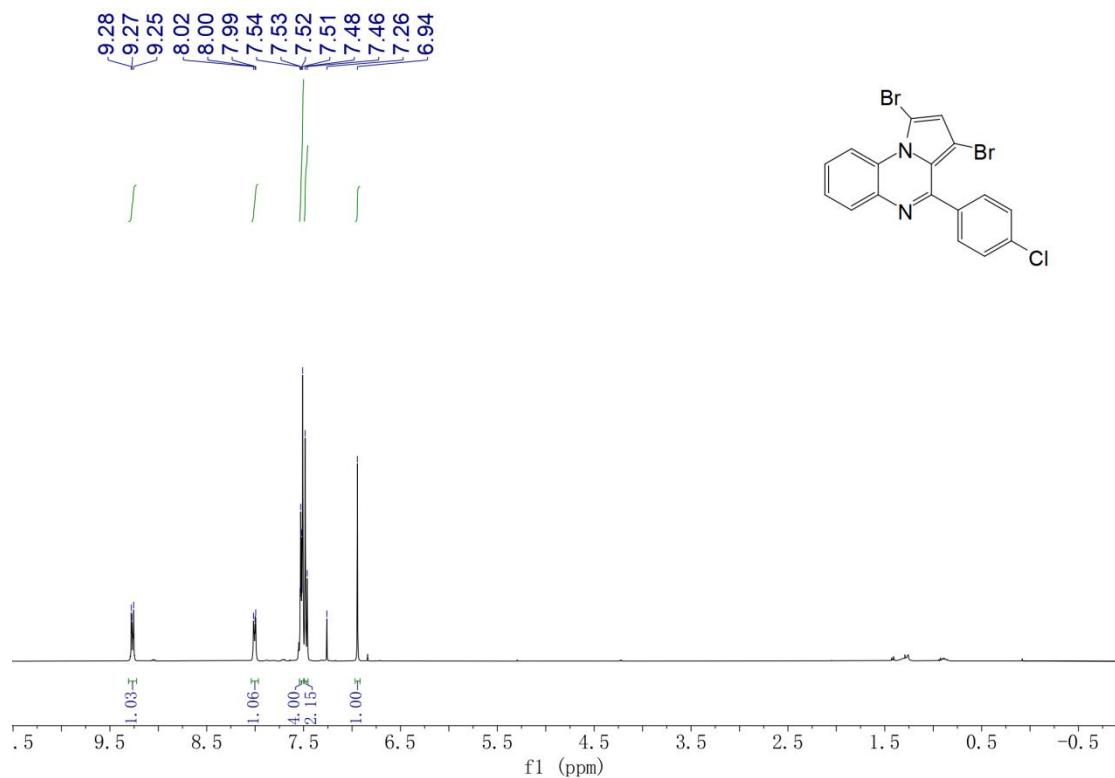
1,3-dibromo-4-(p-tolyl)pyrrolo[1,2-a]quinoxaline [4c]



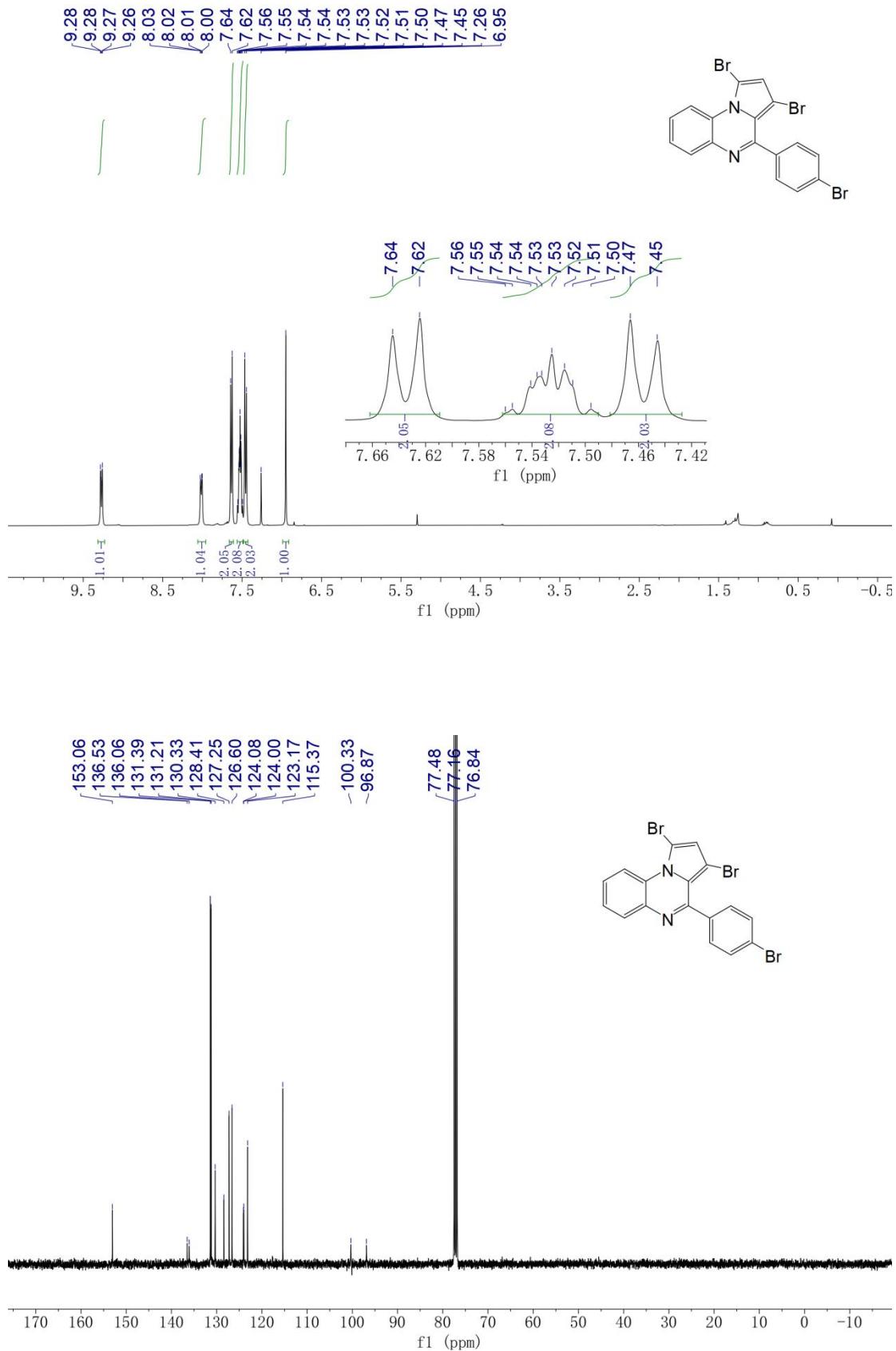
1,3-dibromo-4-(4-methoxyphenyl)pyrrolo[1,2-*a*]quinoxaline [4d]



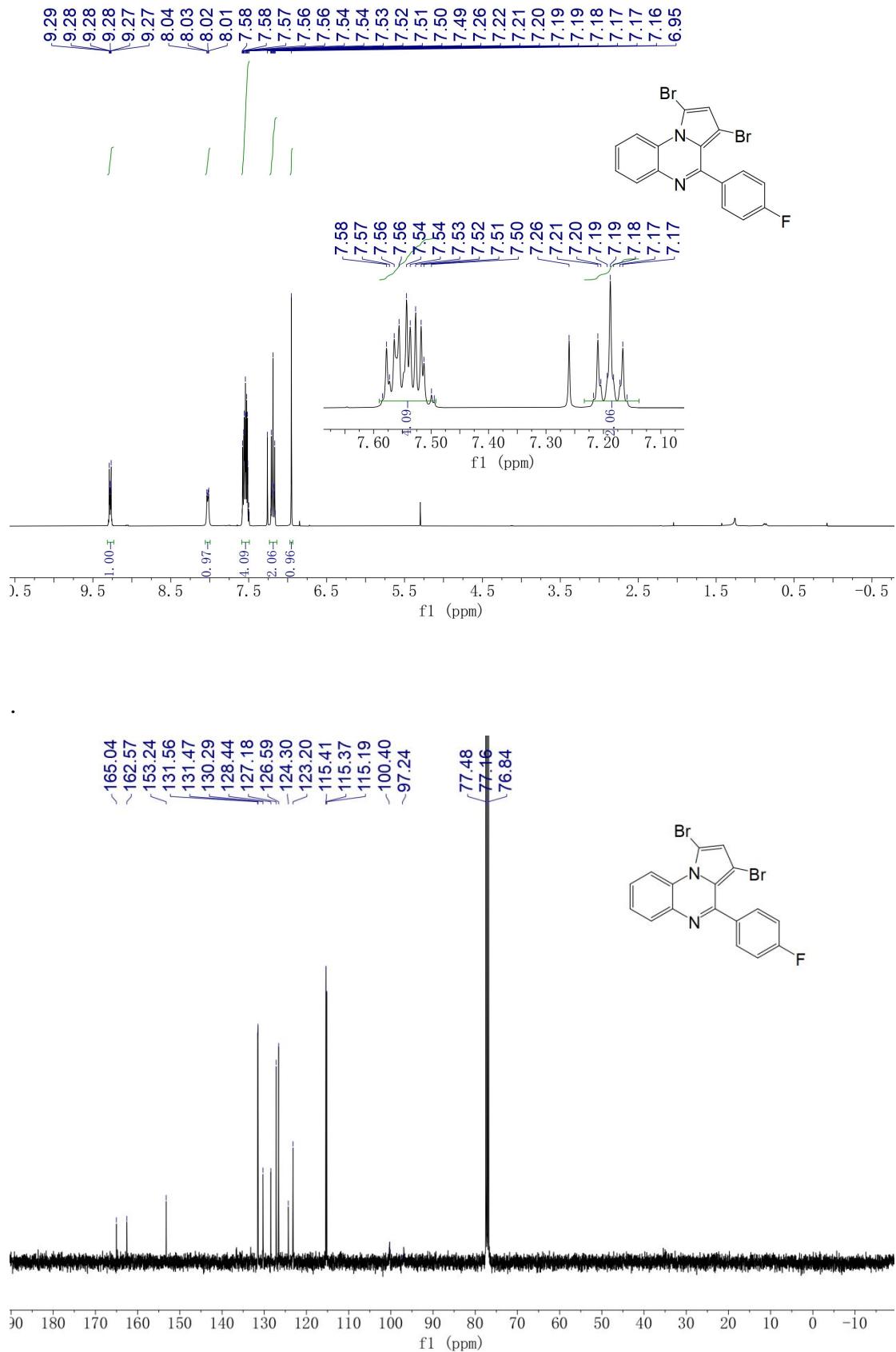
1,3-dibromo-4-(4-chlorophenyl)pyrrolo[1,2-*a*]quinoxaline [4e]

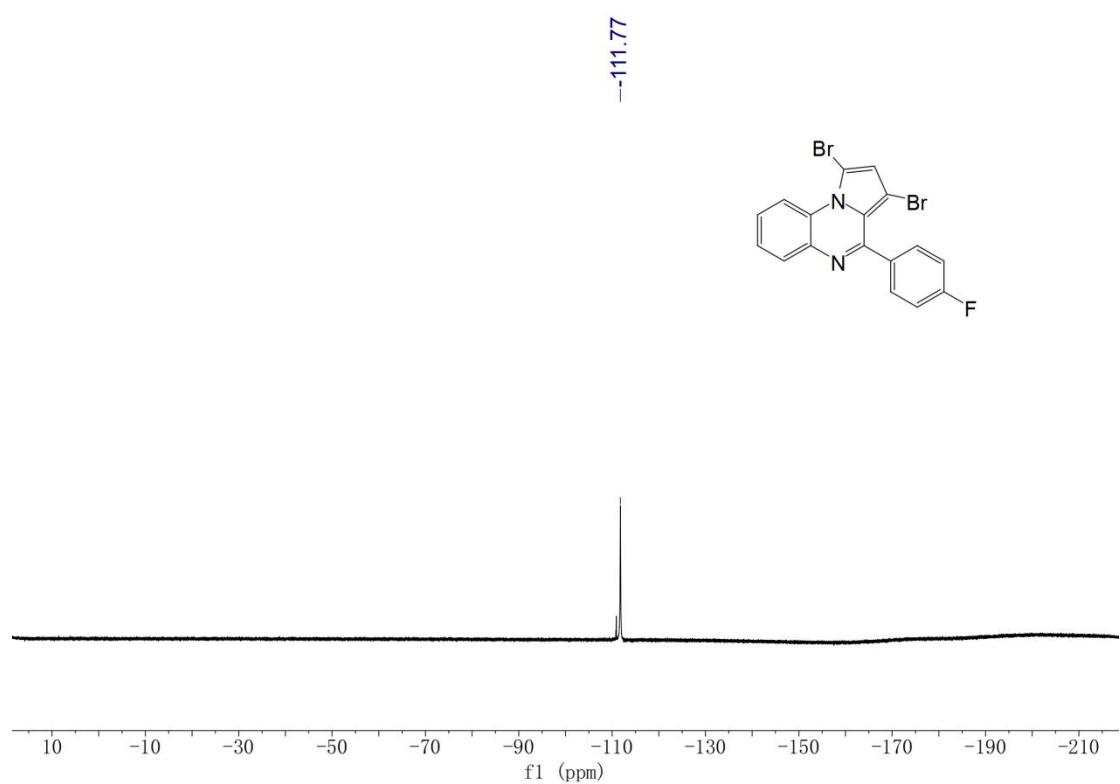


1,3-dibromo-4-(4-bromophenyl)pyrrolo[1,2-a]quinoxaline [4f]

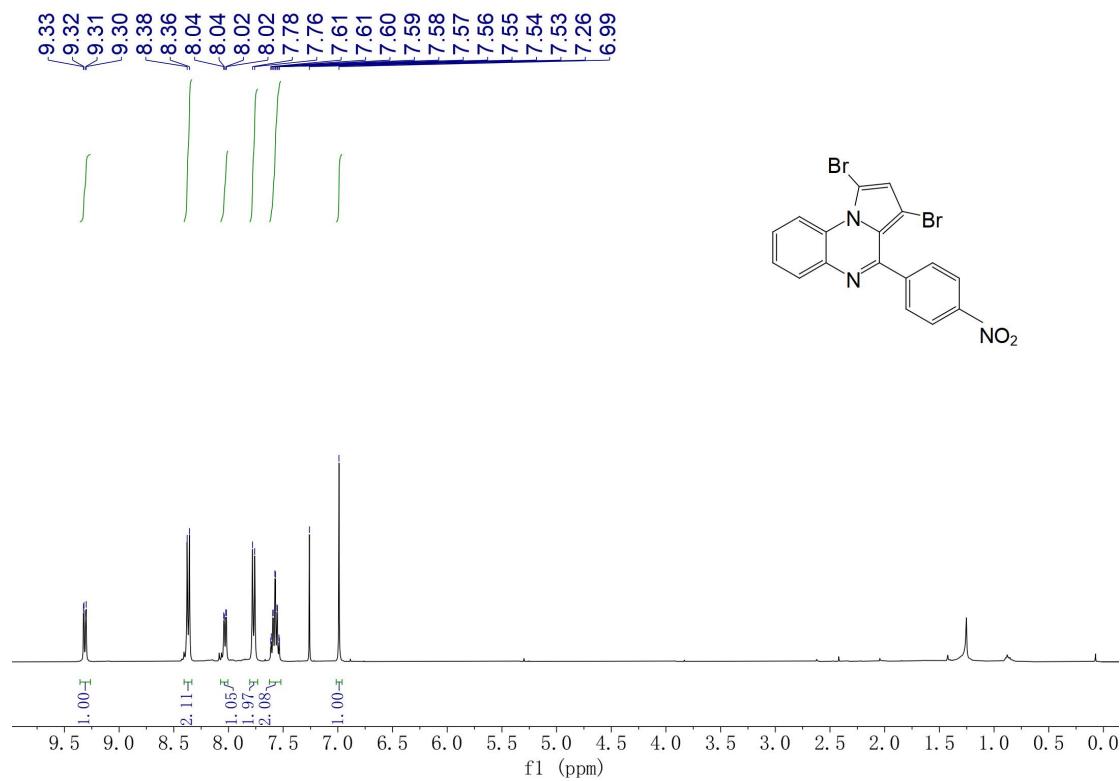


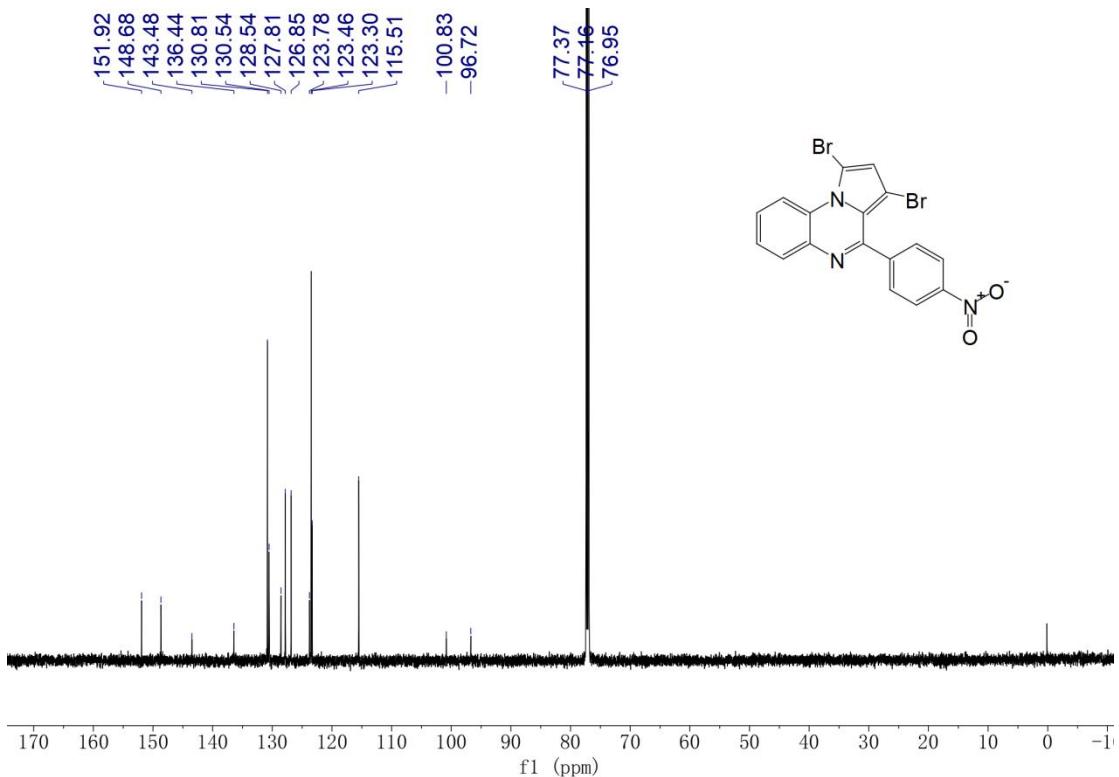
1,3-dibromo-4-(4-fluorophenyl)pyrrolo[1,2-*a*]quinoxaline [4g]



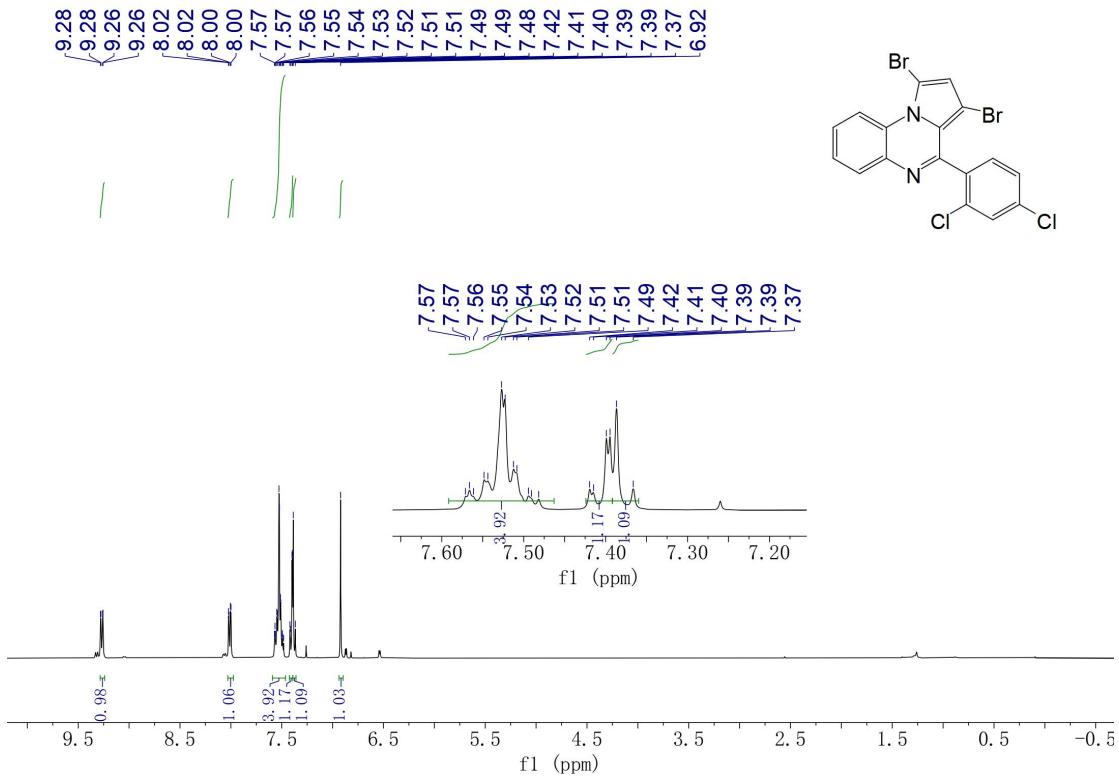


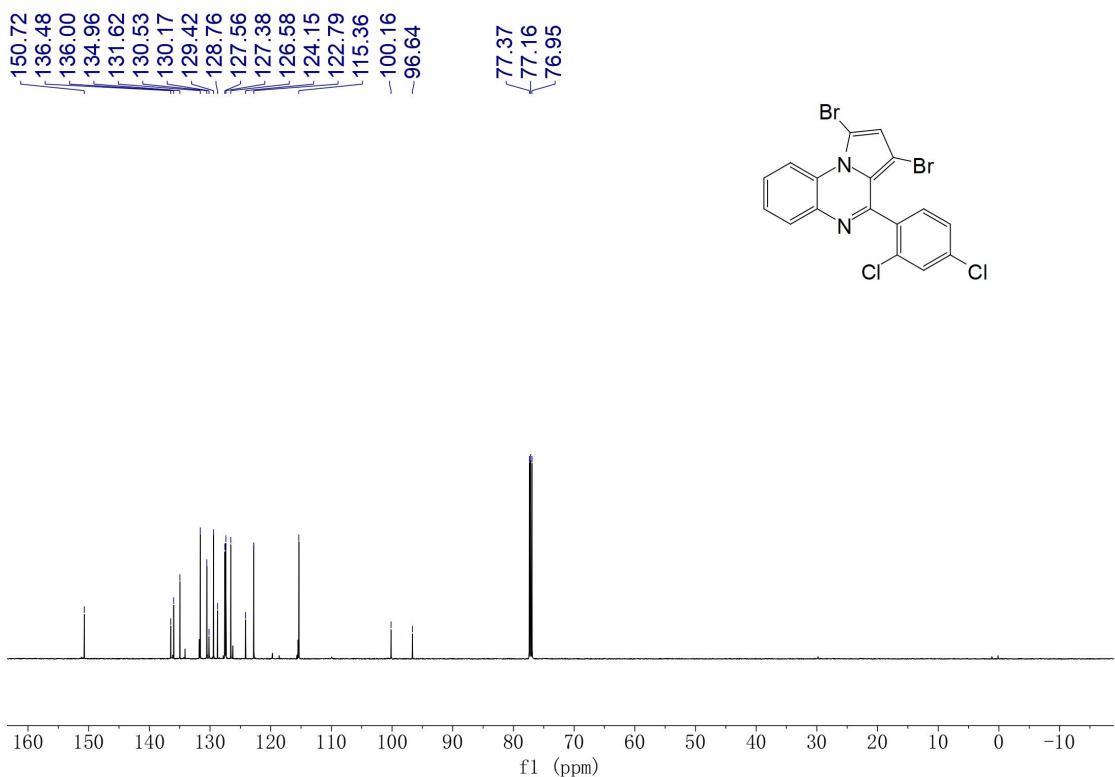
1,3-dibromo-4-(4-nitrophenyl)pyrrolo[1,2-a]quinoxaline [4h]



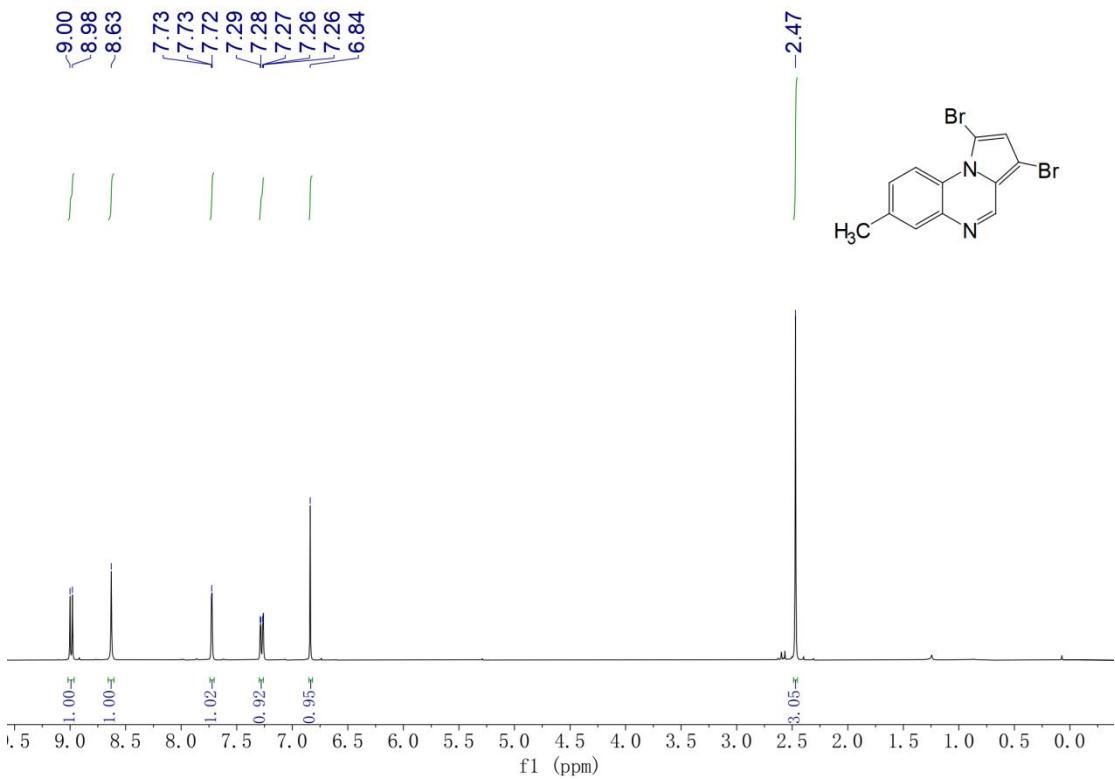


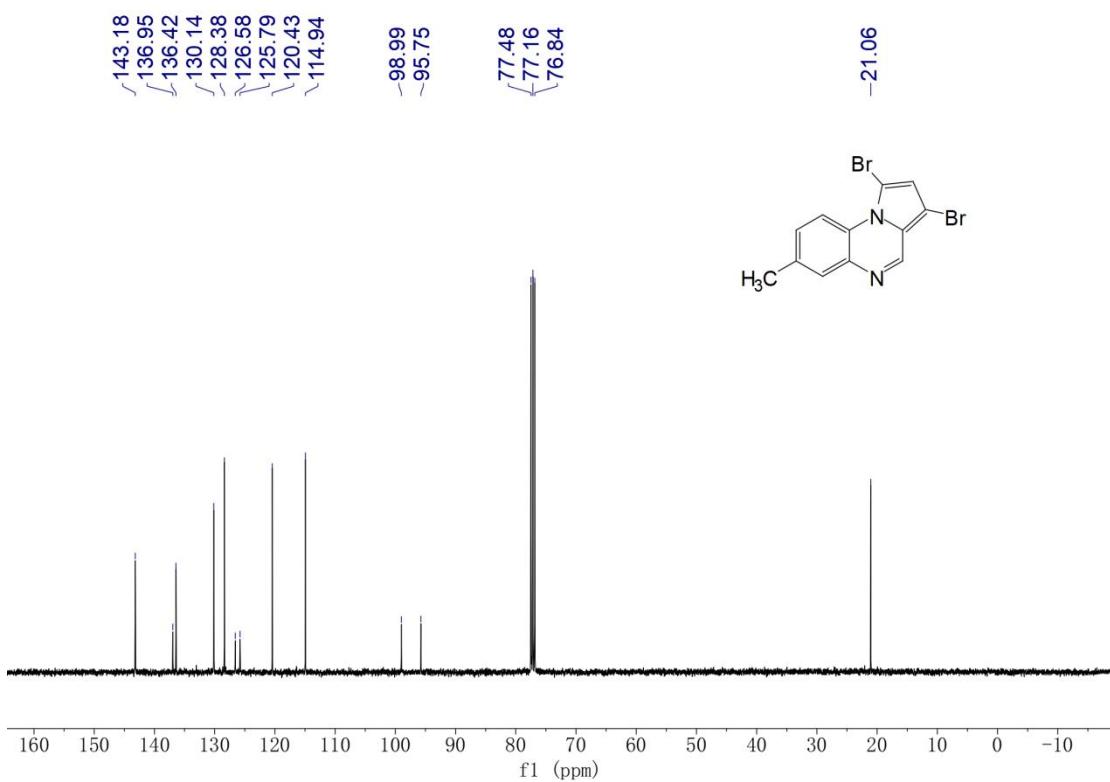
1,3-dibromo-4-(2,4-dichlorophenyl)pyrrolo[1,2-*a*]quinoxaline [4i]



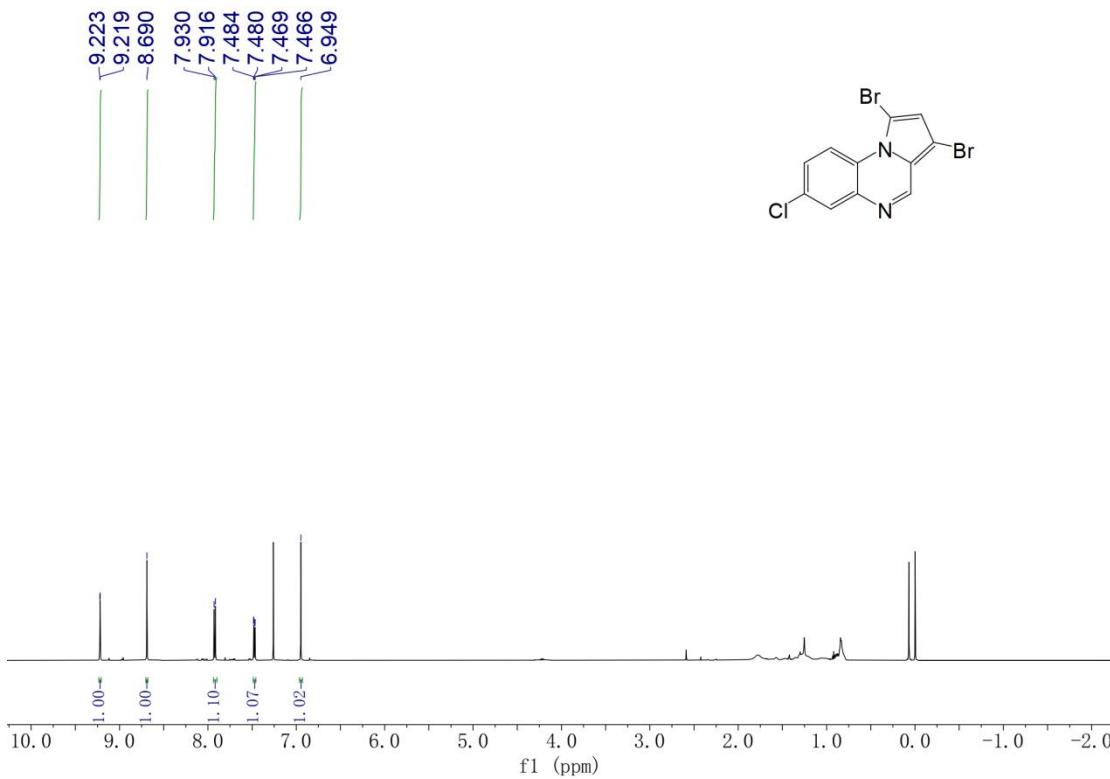


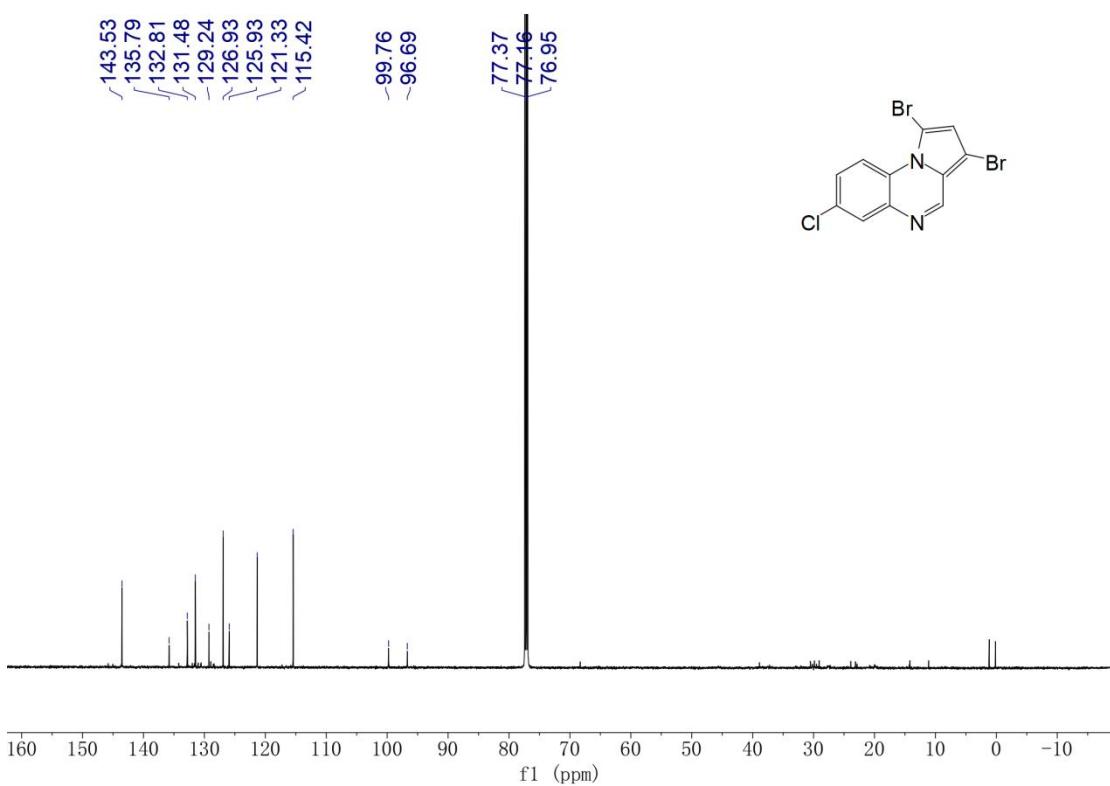
1,3-dibromo-7-methylpyrrolo[1,2-a]quinoxaline [4j]



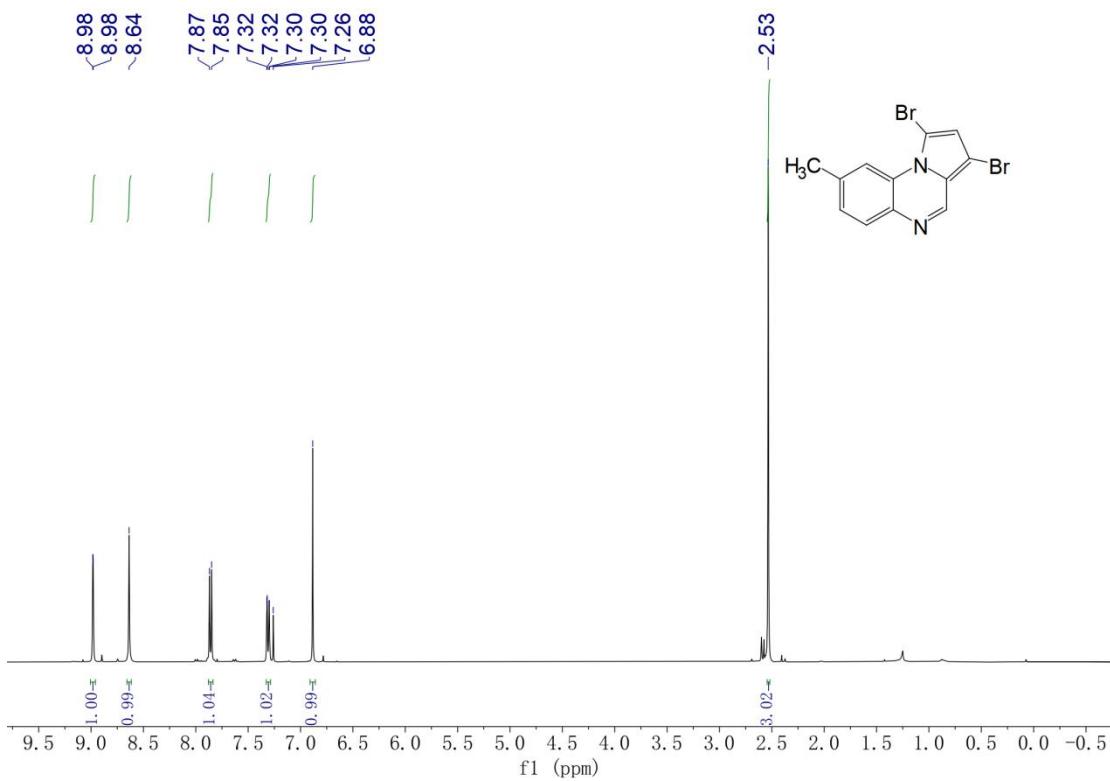


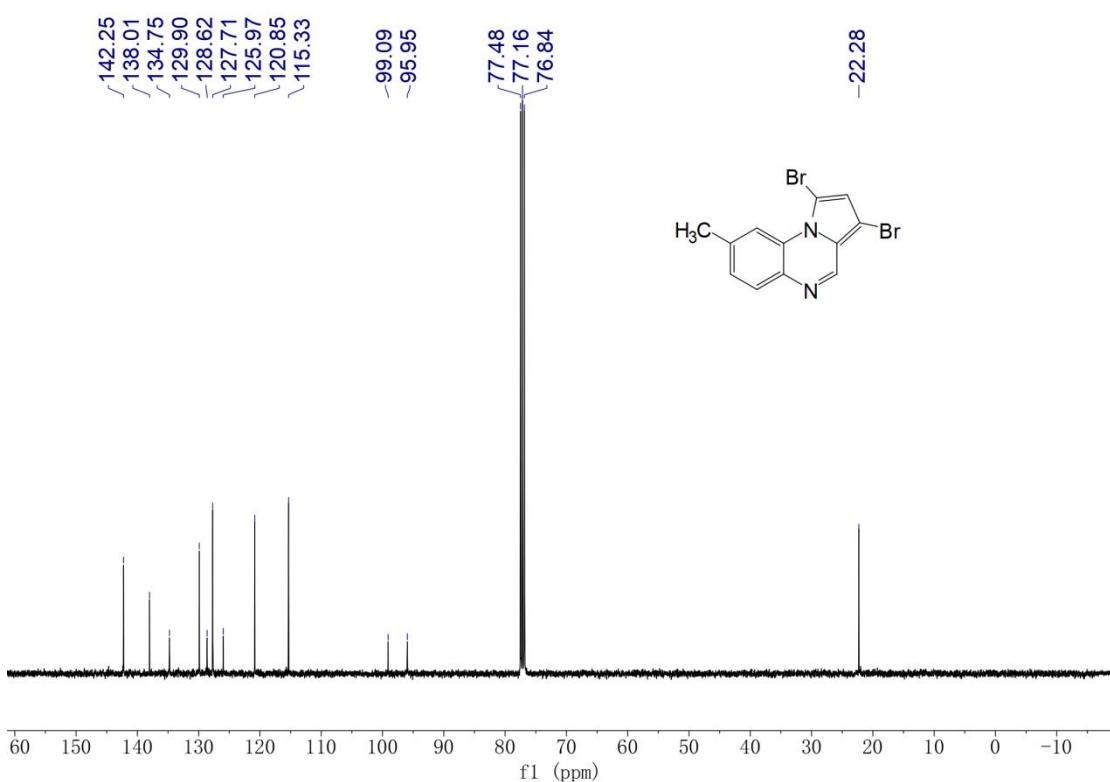
1,3-dibromo-7-chloropyrrolo[1,2-a]quinoxaline [4k]



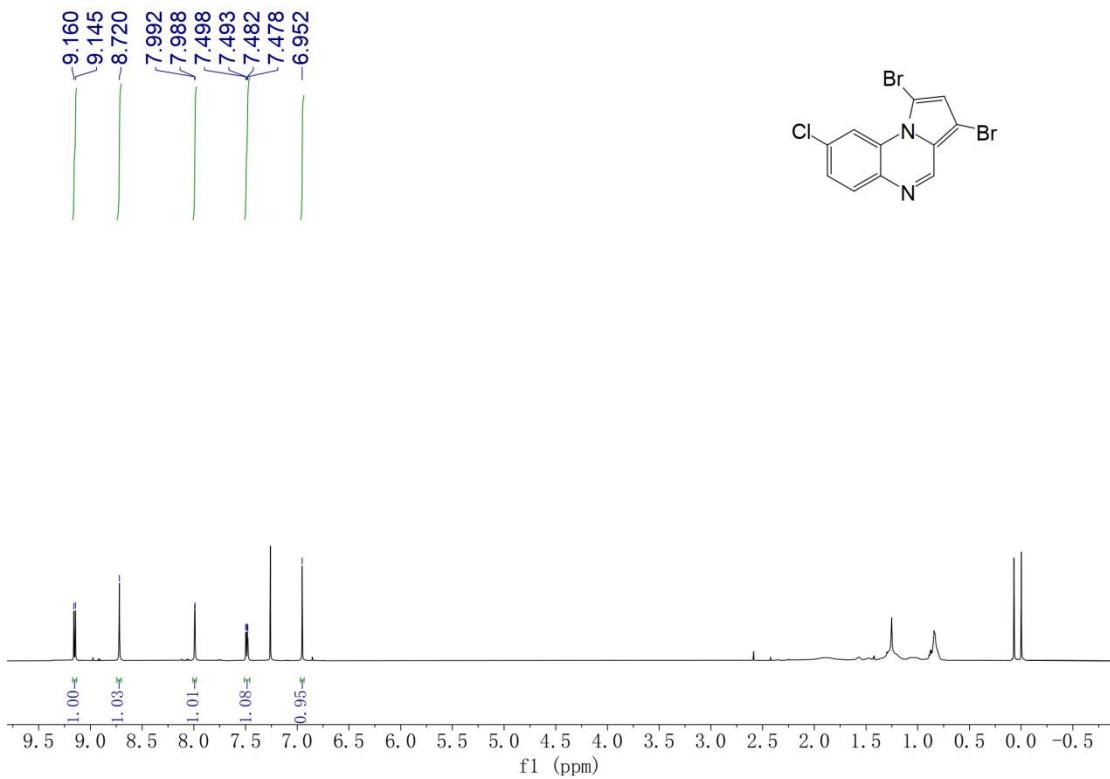


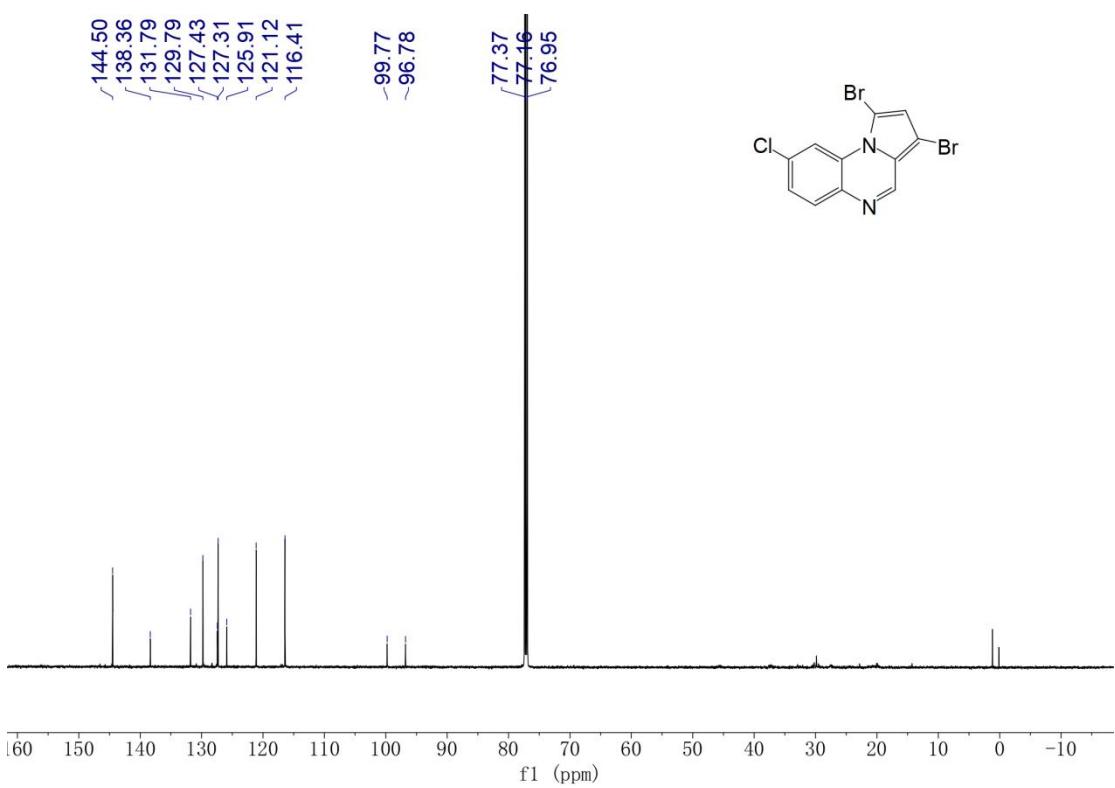
1,3-dibromo-8-methylpyrrolo[1,2-*a*]quinoxaline [4l]



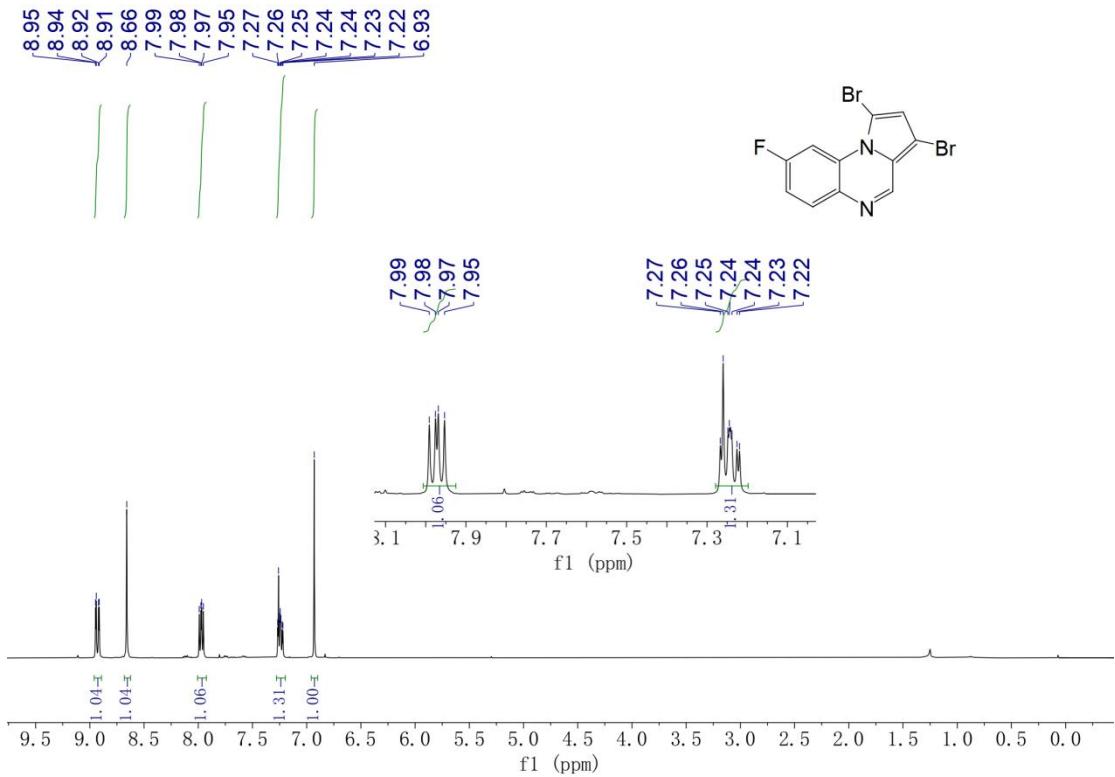


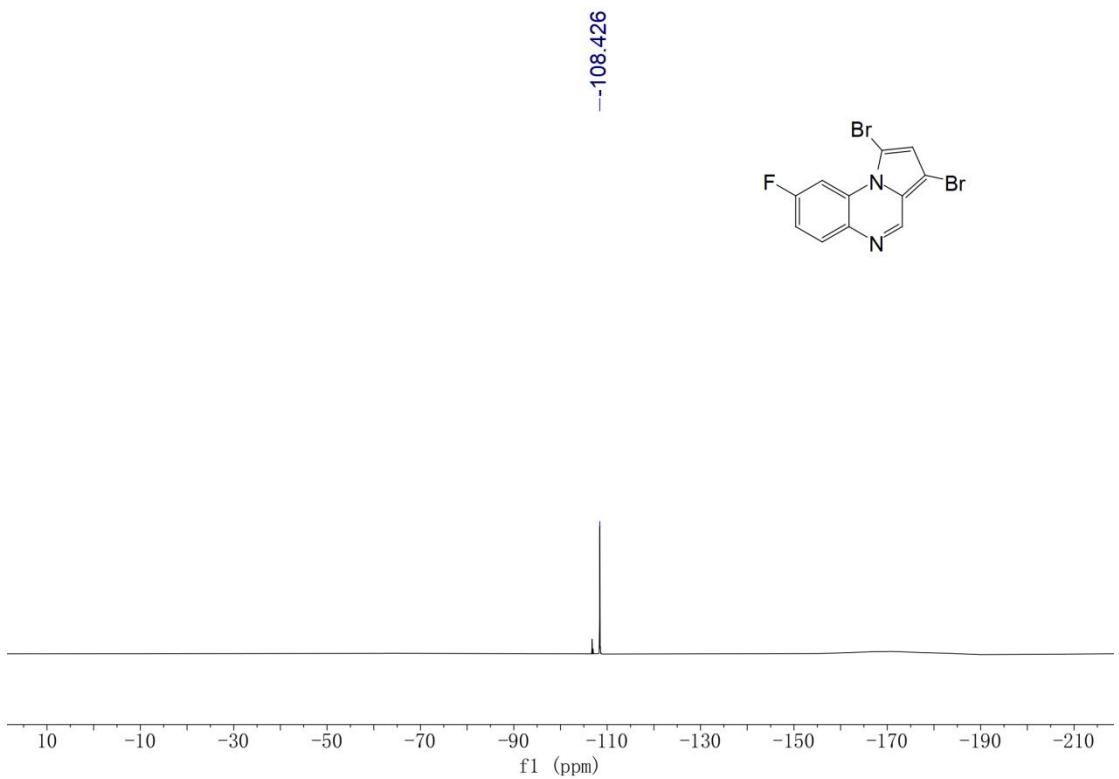
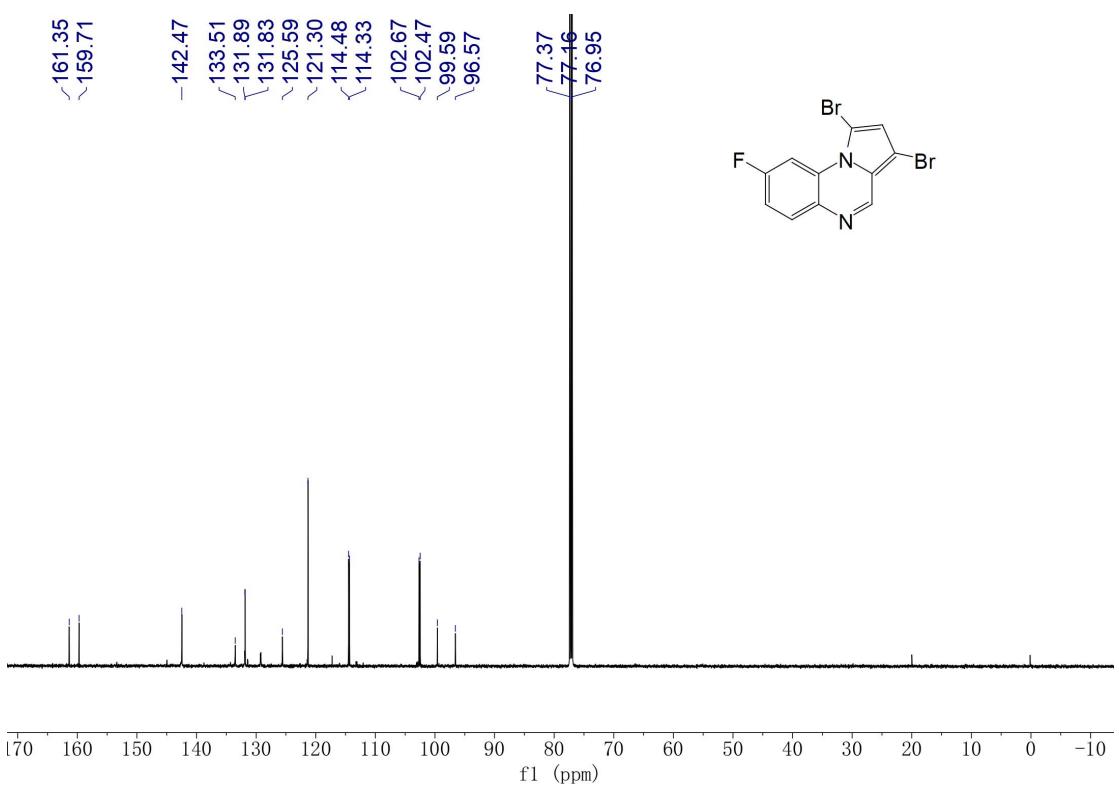
1,3-dibromo-8-chloropyrrolo[1,2-a]quinoxaline [4m]



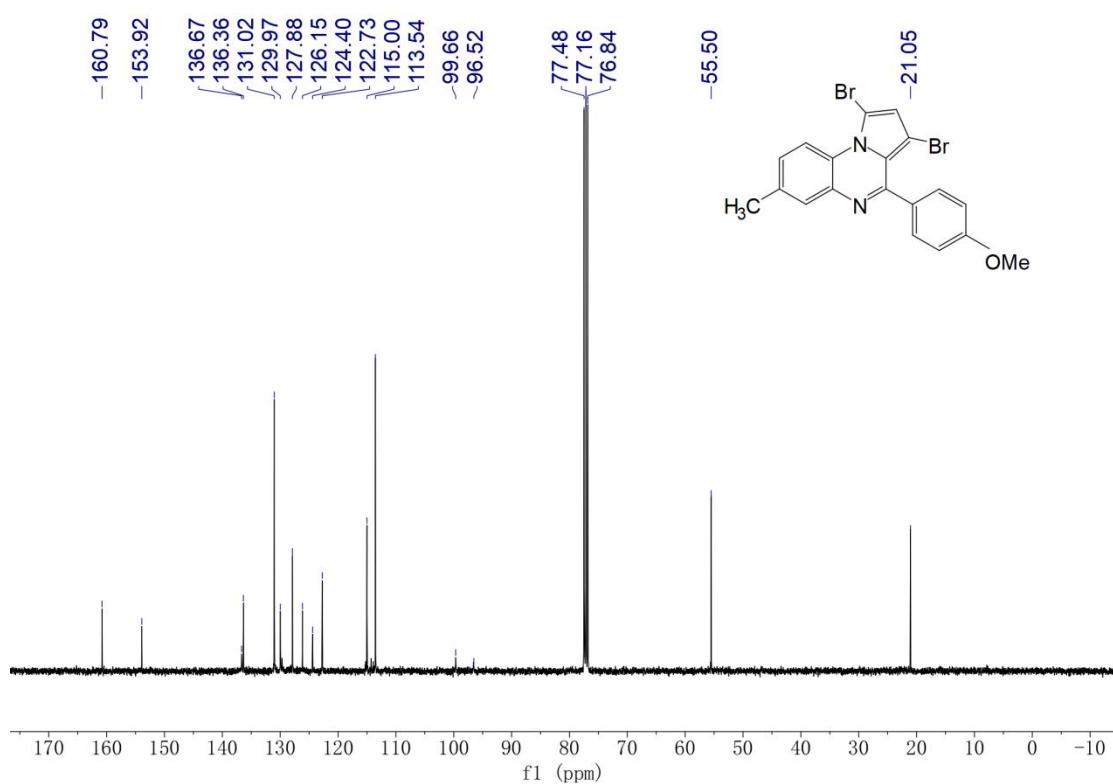
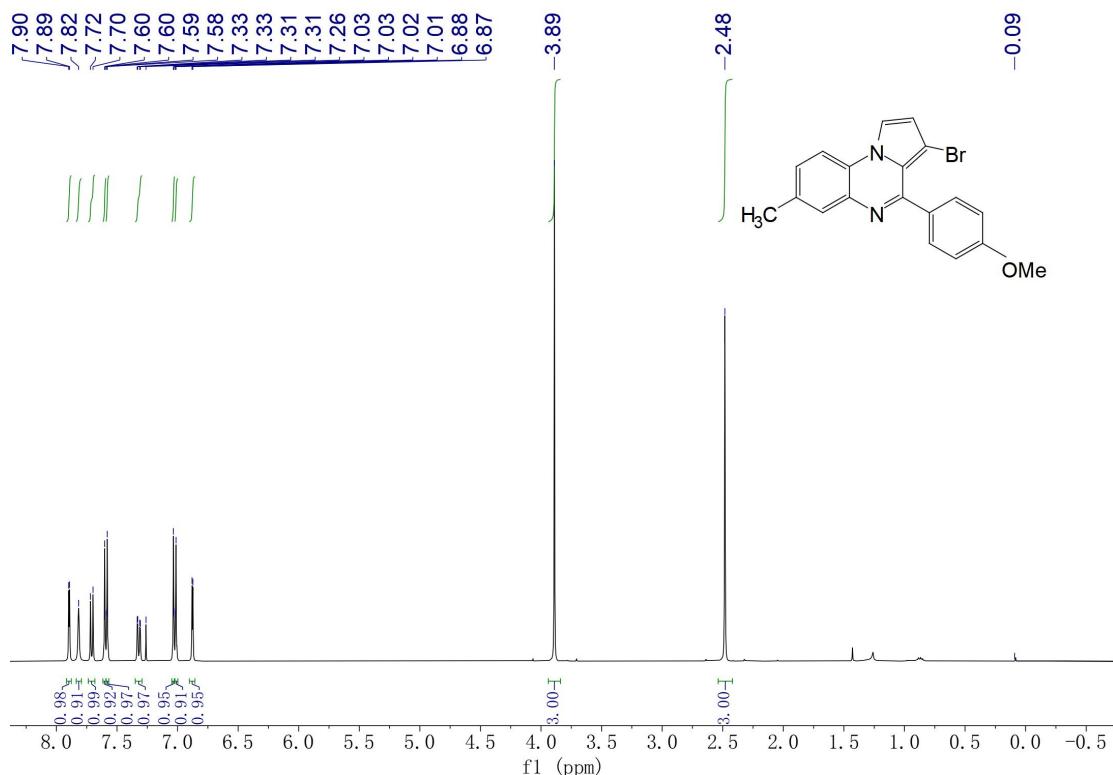


1,3-dibromo-8-fluoropyrrolo[1,2-*a*]quinoxaline [4n]

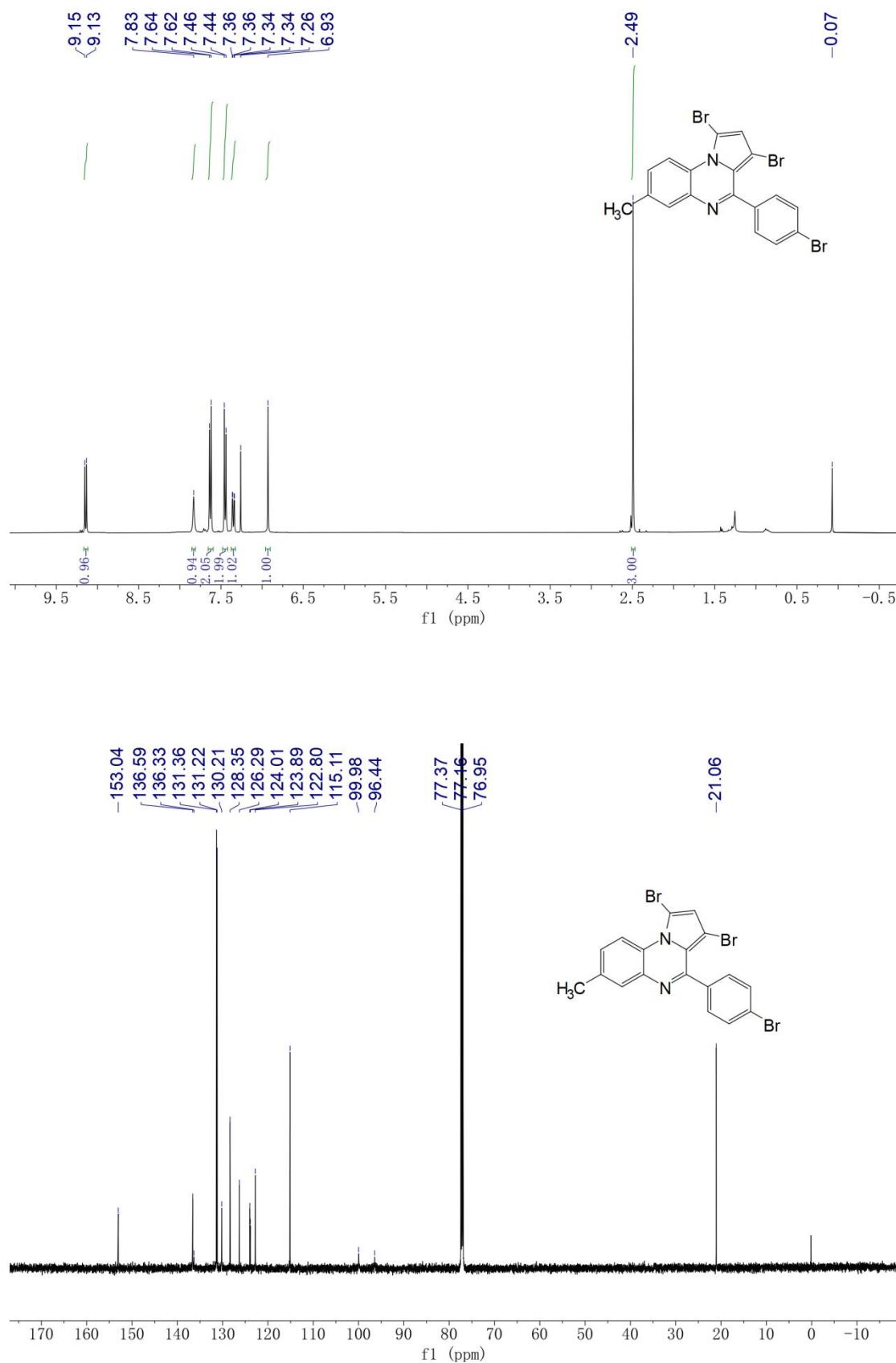




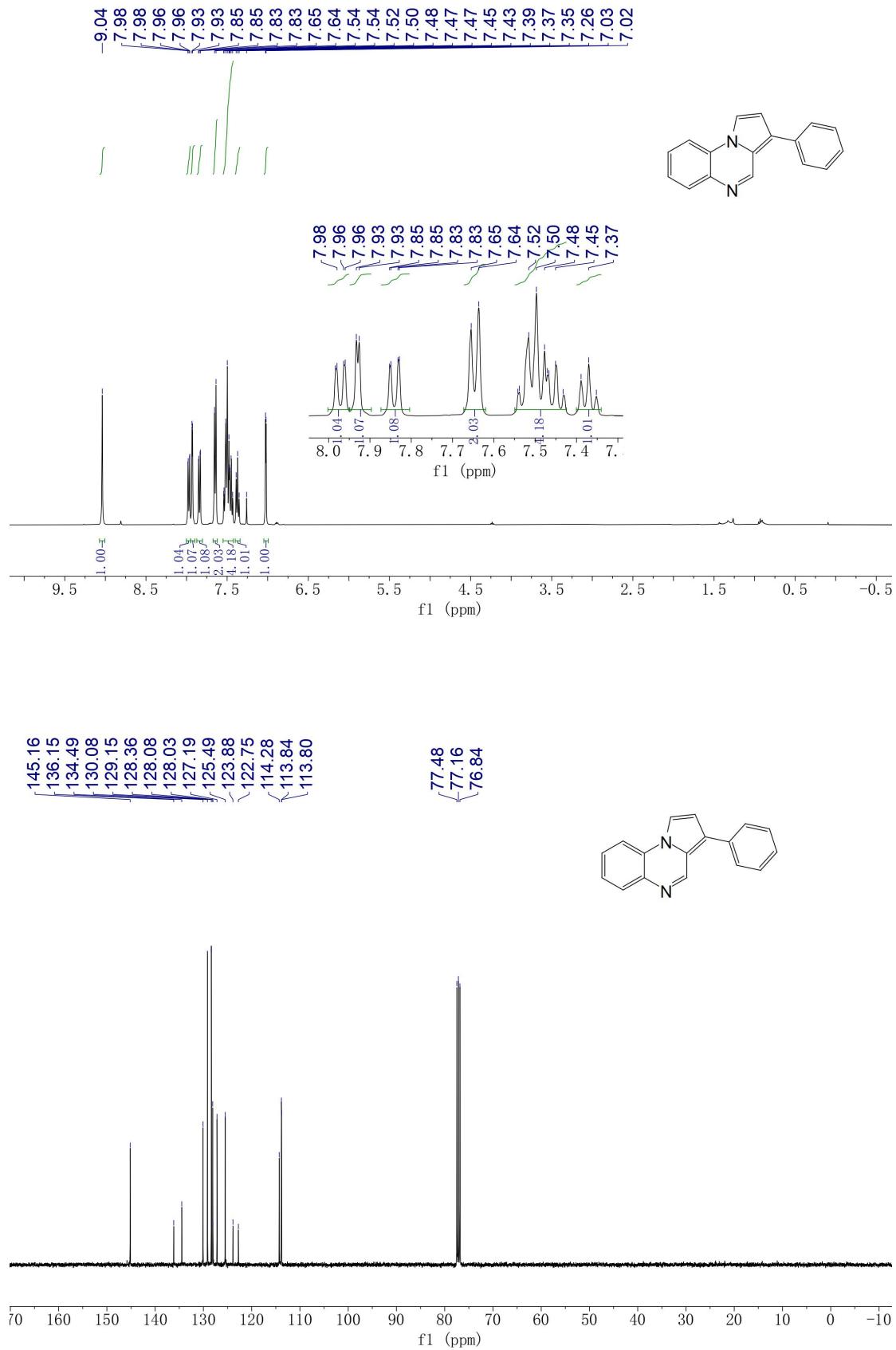
1,3-dibromo-4-(4-methoxyphenyl)-7-methylpyrrolo[1,2-*a*]quinoxaline [4o]



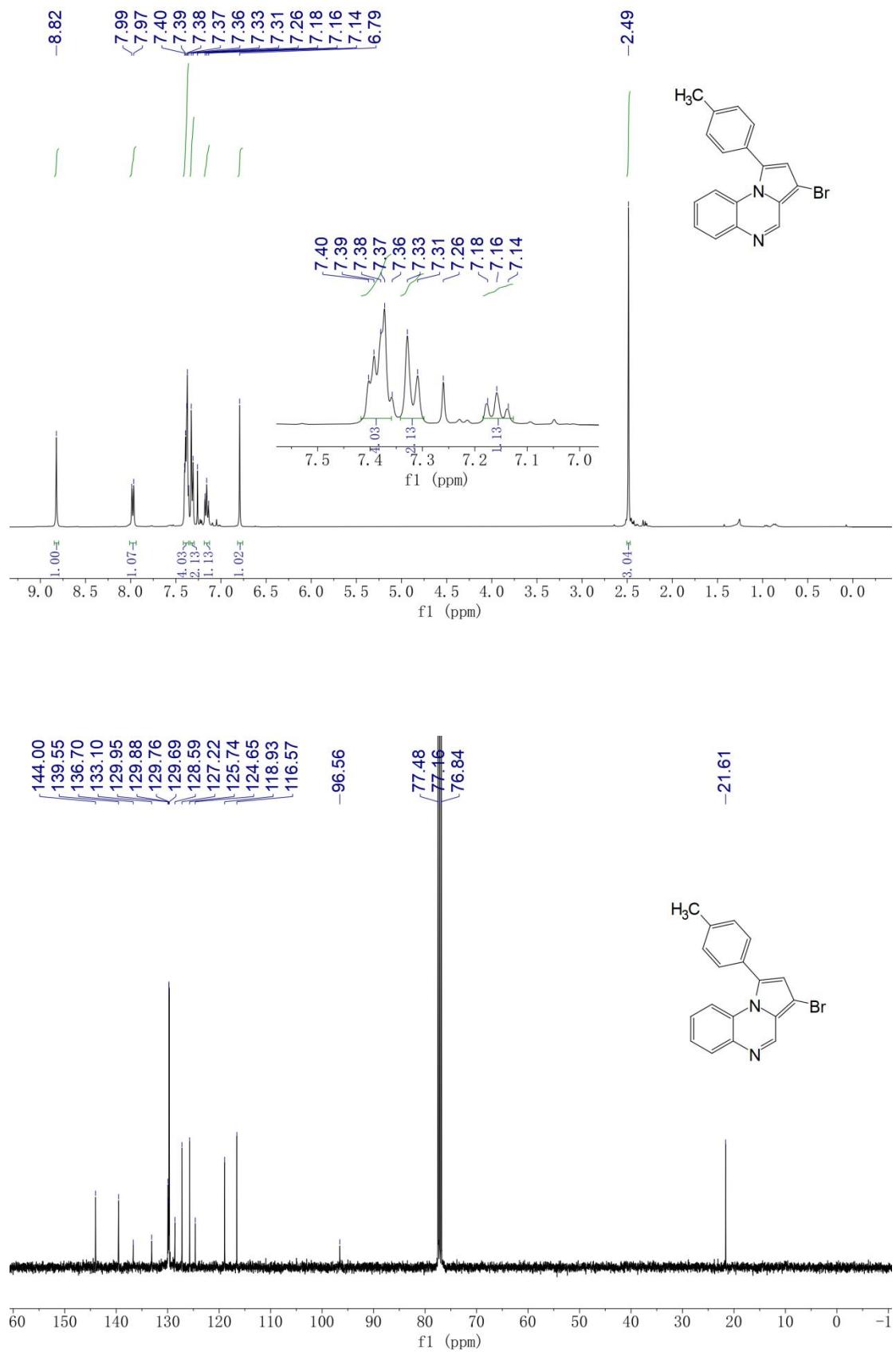
1,3-dibromo-4-(4-bromophenyl)-7-methylpyrrolo[1,2-*a*]quinoxaline [4q]



3-phenylpyrrolo[1,2-*a*]quinoxaline [3aa]



3-bromo-1-(p-tolyl)pyrrolo[1,2-*a*]quinoxaline [3ab]



3-bromo-1-chloropyrrolo[1,2-*a*]quinoxaline [3ac]

