

Supporting Information

Iridium-Catalyzed Asymmetric Hydrogenation of Cyclic Iminium Salts

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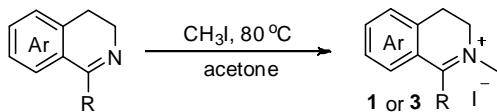
1. General

Commercially available reagents were used without further purification. ^1H NMR, ^{13}C NMR and ^{19}F NMR spectra were recorded at room temperature in CDCl_3 or DMSO on 400 MHz instrument with TMS (tetramethylsilane) as the internal standard. Enantiomeric excess was determined by HPLC using the chiral column described below in detail. Optical rotations were measured by polarimeter. Flash column chromatography was performed on silica gel (200-300 mesh).

2. General Procedure for Synthesis of Cyclic Iminium Salts

1-Substituted 3,4-dihydrosioquinolines and 1-substituted 4,9-dihydro-3- H - β -carbolines were synthesized according to the known literature procedures.¹ The cyclic iminium salts **1** and **3** were synthesized from the corresponding imines according to the known literature procedures.²

2.1 General Procedure for Synthesis of Cyclic Iminium Iodides



General Procedure: A mixture of 1-substituted 3,4-dihydrosioquinoline (or 1-substituted 4,9-dihydro-3- H - β -caroline) (1.0 mmol) and iodomethane (0.50 mL) in acetone (2.0 mL) was stirred at 80 °C overnight and the solid gradually formed. Afterwards the precipitate was filtered from the solution and washed with ethyl acetate to give the corresponding cyclic iminium iodides.

2-Methyl-1-phenyl-3,4-dihydroisoquinolin-2-ium iodide (1a): 5.0 mmol scale, 1.667 g, 95% yield, unknown compound, pale yellow solid, m.p. 217–218 °C; ^1H NMR (400 MHz, CDCl_3) δ 7.77 (dd, $J = 5.2, 3.3$ Hz, 2H), 7.64 (m, 4H), 7.42 (d, $J = 7.5$ Hz, 1H), 7.33 – 7.26 (m, 1H), 7.02 (d, $J = 7.9$ Hz, 1H), 4.56 (t, $J = 7.8$ Hz, 2H), 3.81 (s, 3H), 3.59 (t, $J = 7.8$ Hz, 2H); ^{13}C NMR (100 MHz, CDCl_3) δ 175.2, 137.4, 137.1, 134.0, 132.1, 129.4, 129.3, 129.0, 128.4, 128.0, 127.5, 53.5, 48.1, 26.0; HRMS calculated for $\text{C}_{16}\text{H}_{16}\text{N} [\text{M}-\text{I}]^+$ 222.1277, found 222.1280.

2-Methyl-1-(*m*-tolyl)-3,4-dihydroisoquinolin-2-ium iodide (1b): 1.0 mmol scale, 0.323 g, 89% yield, unknown compound, pale yellow solid, m.p. 201–202 °C; ^1H NMR (400 MHz, CDCl_3) δ 7.77 (t, $J = 7.5$ Hz, 1H), 7.61 – 7.54 (m, 3H), 7.42 – 7.39 (m, 3H), 6.97 (d, $J = 7.9$ Hz, 1H), 4.25 (t, $J = 7.7$ Hz, 2H), 3.50 (s, 3H), 3.34 (t, $J = 7.7$ Hz, 2H), 2.42 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 175.2, 139.4, 137.4, 137.0, 134.0, 132.8, 129.3, 129.2, 128.9, 128.4, 128.0, 127.5, 126.0, 53.4, 48.0, 26.0, 21.4; HRMS calculated for $\text{C}_{17}\text{H}_{18}\text{N} [\text{M}-\text{I}]^+$ 236.1434, found 236.1436.

2-Methyl-1-(*p*-tolyl)-3,4-dihydroisoquinolin-2-ium iodide (1c): 1.0 mmol scale, 0.330 g, 91% yield, unknown compound, pale yellow solid, m.p. 197–198 °C; ^1H NMR (400 MHz, CDCl_3) δ 7.67 – 7.61 (m, 3H), 7.42 – 7.40 (m, 3H), 7.30 (t, $J = 7.7$ Hz, 1H), 7.05 (d, $J = 7.8$ Hz, 1H), 4.55 (t, $J = 7.7$ Hz, 2H), 3.84 (s, 3H), 3.56 (t, $J = 7.8$ Hz, 2H), 2.48 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 175.4, 143.1, 137.5, 137.0, 134.2, 129.9, 129.2, 128.4, 128.0, 127.7, 126.4, 53.5, 48.0, 26.0, 21.7; HRMS calculated for $\text{C}_{17}\text{H}_{18}\text{N} [\text{M}-\text{I}]^+$ 236.1434, found 236.1439.

1-(4-Methoxyphenyl)-2-methyl-3,4-dihydroisoquinolin-2-ium iodide (1d): 1.0 mmol scale,

0.231 g, 61% yield, unknown compound, pale yellow solid, m.p. 215-216 °C; ¹H NMR (400 MHz, DMSO) δ 7.78 (t, *J* = 7.5 Hz, 1H), 7.59 (t, *J* = 7.8 Hz, 3H), 7.43 (t, *J* = 7.7 Hz, 1H), 7.27 (d, *J* = 8.7 Hz, 2H), 7.02 (d, *J* = 7.8 Hz, 1H), 4.26 (t, *J* = 7.6 Hz, 2H), 3.90 (s, 3H), 3.59 (s, 3H), 3.36 (t, *J* = 7.5 Hz, 2H); ¹³C NMR (100 MHz, DMSO) δ 178.7, 167.2, 143.4, 141.5, 138.5, 136.5, 133.4, 133.1, 133.0, 126.6, 119.7, 60.9, 57.5, 51.2, 29.9; HRMS calculated for C₁₇H₁₈NO [M-I]⁺ 252.1383, found 252.1388.

1-(4-Fluorophenyl)-2-methyl-3,4-dihydroisoquinolin-2-i um iodide (1e): 0.88 mmol scale, 0.283 g, 81% yield, unknown compound, pale yellow solid, m.p. 188-189 °C; ¹H NMR (400 MHz, CDCl₃) δ 7.91 – 7.87 (m, 2H), 7.67 (t, *J* = 7.5 Hz, 1H), 7.41 (d, *J* = 7.5 Hz, 1H), 7.33 – 7.28 (m, 3H), 7.00 (d, *J* = 7.9 Hz, 1H), 4.52 (t, *J* = 7.8 Hz, 2H), 3.80 (s, 3H), 3.58 (t, *J* = 7.8 Hz, 2H); ¹³C NMR (100 MHz, CDCl₃) δ 174.4, 164.6 (d, *J* = 255.2 Hz), 137.5, 137.2, 133.9, 132.1 (d, *J* = 8.8 Hz), 128.5, 128.1, 127.6, 125.3 (d, *J* = 3.5 Hz), 116.7 (d, *J* = 22.3 Hz), 53.6, 48.2, 25.9; ¹⁹F NMR (376 MHz, CDCl₃) δ -105.4; HRMS calculated for C₁₆H₁₅NF [M-I]⁺ 240.1183, found 240.1185.

1-(4-Chlorophenyl)-2-methyl-3,4-dihydroisoquinolin-2-i um iodide (1f): 1.0 mmol scale, 0.355 g, 92% yield, unknown compound, pale yellow solid, m.p. 234-235 °C; ¹H NMR (400 MHz, DMSO) δ 7.83 – 7.71 (m, 3H), 7.73 (d, *J* = 8.5 Hz, 2H), 7.59 (d, *J* = 7.5 Hz, 1H), 7.42 (t, *J* = 7.7 Hz, 1H), 6.99 (d, *J* = 7.9 Hz, 1H), 4.30 (t, *J* = 7.7 Hz, 2H), 3.54 (s, 3H), 3.40 (t, *J* = 7.7 Hz, 2H); ¹³C NMR (100 MHz, DMSO) δ 173.3, 138.3, 137.3, 137.1, 133.4, 131.0, 129.8, 128.9, 128.8, 128.4, 128.0, 52.9, 46.7, 25.0; HRMS calculated for C₁₆H₁₅NCl [M-I]⁺ 256.0888, found 256.0888.

1-(4-Bromophenyl)-2-methyl-3,4-dihydroisoquinolin-2-i um iodide (1g): 1.3 mmol scale, 0.523 g, 94% yield, unknown compound, pale yellow solid, m.p. 234-235 °C; ¹H NMR (400 MHz, DMSO) δ 7.96 (d, *J* = 8.4 Hz, 2H), 7.79 (t, *J* = 7.5 Hz, 1H), 7.65 (d, *J* = 8.4 Hz, 2H), 7.59 (d, *J* = 7.5 Hz, 1H), 7.42 (t, *J* = 7.6 Hz, 1H), 6.99 (d, *J* = 7.8 Hz, 1H), 4.30 (t, *J* = 7.7 Hz, 2H), 3.54 (s, 3H), 3.40 (t, *J* = 7.7 Hz, 2H); ¹³C NMR (100 MHz, DMSO) δ 173.3, 138.3, 137.1, 133.4, 132.7, 131.1, 129.3, 128.7, 128.4, 127.9, 126.2, 52.9, 46.7, 25.0; HRMS calculated for C₁₆H₁₅NBr [M-I]⁺ 300.0382, found 300.0386.

2,7-Dimethyl-1-phenyl-3,4-dihydroisoquinolin-2-i um iodide (1h): 1.0 mmol scale, 0.280 g, 77% yield, unknown compound, pale yellow solid, m.p. 202-204 °C; ¹H NMR (400 MHz, DMSO) δ 7.81 – 7.57 (m, 6H), 7.49 (d, *J* = 7.7 Hz, 1H), 6.74 (s, 1H), 4.29 (t, *J* = 7.7 Hz, 2H), 3.52 (s, 3H), 3.38 – 3.33 (m, 2H), 2.22 (s, 3H); ¹³C NMR (100 MHz, DMSO) δ 174.2, 137.7, 135.5, 133.3, 132.3, 130.2, 129.6, 128.8, 128.8, 128.7, 128.0, 52.9, 46.6, 24.7, 21.0; HRMS calculated for C₁₇H₁₈N [M-I]⁺ 236.1434, found 236.1435.

7-Methoxy-2-methyl-1-phenyl-3,4-dihydroisoquinolin-2-i um iodide (1i): 1.0 mmol scale, 0.385 g, 93% yield, unknown compound, pale yellow solid, m.p. 233-235 °C; ¹H NMR (400 MHz, DMSO) δ 7.83 – 7.69 (m, 5H), 7.55 (d, *J* = 8.4 Hz, 1H), 7.44 – 7.42 (m, 1H), 6.35 (d, *J* = 2.5 Hz, 1H), 4.31 (t, *J* = 7.7 Hz, 2H), 3.65 (s, 3H), 3.56 (s, 3H), 3.34 (t, *J* = 7.7 Hz, 2H); ¹³C NMR (100 MHz, DMSO) δ 173.9,

158.5, 132.4, 130.2, 130.0, 129.6, 128.9, 128.9, 128.8, 122.1, 118.7, 56.1, 53.3, 46.8, 24.3; HRMS calculated for C₁₇H₁₈NO [M-I]⁺ 252.1383, found 252.1388.

7-Chloro-2-methyl-1-phenyl-3,4-dihydroisoquinolin-2-ium iodide (1j): 0.8 mmol scale, 0.244 g, 79% yield, unknown compound, yellow solid, m.p. 237-238 °C; ¹H NMR (400 MHz, DMSO) δ 7.86 (dd, *J* = 8.2, 1.6 Hz, 1H), 7.80 – 7.67 (m, 5H), 7.64 (d, *J* = 8.2 Hz, 1H), 6.81 (d, *J* = 1.4 Hz, 1H), 4.32 (t, *J* = 7.7 Hz, 2H), 3.55 (s, 3H), 3.41 (t, *J* = 7.7 Hz, 2H); ¹³C NMR (100 MHz, DMSO) δ 175.1, 139.2, 138.3, 134.5, 134.3, 133.8, 132.7, 131.7, 131.6, 131.5, 130.9, 54.9, 48.9, 26.5; HRMS calculated for C₁₆H₁₅ClN [M-I]⁺ 256.0888, found 256.0887.

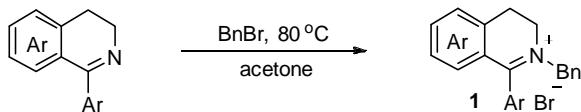
2-Methyl-1-phenethyl-3,4-dihydroisoquinolinium iodide (1k): 1.0 mmol scale, 0.309 g, 82% yield, unknown compound, pale yellow solid, m.p. 164-165 °C; ¹H NMR (400 MHz, DMSO) δ 8.13 (d, *J* = 7.9 Hz, 1H), 7.76 (t, *J* = 7.4 Hz, 1H), 7.52 (dd, *J* = 15.8, 7.7 Hz, 2H), 7.31 (d, *J* = 4.5 Hz, 4H), 7.27 – 7.20 (m, 1H), 4.05 (t, *J* = 7.5 Hz, 2H), 3.74 (s, 3H), 3.56 (t, *J* = 7.9 Hz, 2H), 3.12 (t, *J* = 7.5 Hz, 2H), 3.02 – 2.90 (m, 2H); ¹³C NMR (100 MHz, DMSO) δ 177.4, 139.2, 137.8, 136.5, 130.8, 129.1, 129.0, 128.7, 128.6, 127.3, 126.7, 52.9, 45.3, 33.0, 32.4, 25.4; HRMS calculated for C₁₈H₂₀N [M-I]⁺ 250.1590, found 250.1593.

2-Methyl-1-phenyl-4,9-dihydro-3*H*-pyrido[3,4-*b*]indol-2-ium iodide (3a): 2.0 mmol scale, 0.630 g, 81% yield, unknown compound, yellow solid, m.p. 230-231 °C; ¹H NMR (400 MHz, DMSO) δ 11.58 (brs, 1H), 7.87 – 7.68 (m, 6H), 7.50 – 7.36 (m, 2H), 7.23 – 7.19 (m, 1H), 4.34 (t, *J* = 8.6 Hz, 2H), 3.57 – 3.38 (m, 5H); ¹³C NMR (100 MHz, DMSO) δ 164.3, 142.1, 132.7, 129.8, 129.4, 129.0, 128.8, 127.6, 124.4, 124.3, 122.3, 122.0, 114.1, 53.8, 44.2, 19.5; HRMS calculated for C₁₇H₁₈N [M-I]⁺ 261.1386, found 261.1391.

2-Methyl-1-(*p*-tolyl)-4,9-dihydro-3*H*-pyrido[3,4-*b*]indol-2-ium iodide (3b): 5.0 mmol scale, 1.875 g, 93% yield, unknown compound, yellow solid, m.p. 237-239 °C; ¹H NMR (400 MHz, DMSO) δ 11.56 (brs, 1H), 7.81 (d, *J* = 8.2 Hz, 1H), 7.64 (d, *J* = 8.1 Hz, 2H), 7.57 (d, *J* = 8.1 Hz, 2H), 7.47 – 7.37 (m, 2H), 7.21 (ddd, *J* = 7.9, 5.9, 1.8 Hz, 1H), 4.33 (t, *J* = 8.6 Hz, 2H), 3.51 (s, 3H), 3.44 (t, *J* = 8.6 Hz, 2H); ¹³C NMR (100 MHz, DMSO) δ 164.4, 143.0, 142.1, 130.3, 129.6, 128.9, 127.6, 125.9, 124.4, 124.2, 122.3, 122.0, 114.1, 53.7, 44.2, 21.7, 19.4; HRMS calculated for C₁₇H₁₈N [M-I]⁺ 275.1543, found 275.1550.

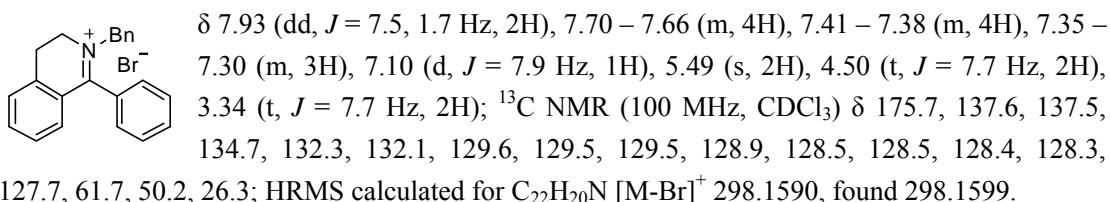
1-(4-Methoxyphenyl)-2-methyl-4,9-dihydro-3*H*-pyrido[3,4-*b*]indol-2-ium iodide (3c): 2.0 mmol scale, 0.795 g, 95% yield, unknown compound, yellow solid, m.p. 254-255 °C; ¹H NMR (400 MHz, DMSO) δ 11.57 (brs, 1H), 7.81 (d, *J* = 8.2 Hz, 1H), 7.71 (d, *J* = 8.8 Hz, 2H), 7.49 – 7.38 (m, 2H), 7.31 (d, *J* = 8.8 Hz, 2H), 7.25 – 7.17 (m, 1H), 4.31 (t, *J* = 8.5 Hz, 2H), 3.93 (s, 3H), 3.55 (s, 3H), 3.42 (t, *J* = 8.5 Hz, 2H); ¹³C NMR (100 MHz, DMSO) δ 164.0, 162.8, 141.9, 132.2, 128.8, 127.7, 124.5, 124.1, 122.2, 122.0, 120.5, 115.2, 114.0, 56.3, 53.8, 44.1, 19.4; HRMS calculated for C₁₇H₁₈N [M-I]⁺ 291.1492, found 291.1497.

2.2 General Procedure for Synthesis of Cyclic Iminium Bromides



General Procedure: A mixture of 1-substituted 3,4-dihydrosioquinoline (1.0 mmol) and (bromomethyl)benzene (1.2 mmol) in acetone (3.0 mL) was stirred at 80 °C overnight and the solid gradually formed. Afterwards the precipitate was filtered from the solution and washed with ethyl acetate. If no precipitate generated, the mixture should be concentrated in vacuum and further purification was performed by a silica gel column eluted with dichloromethane/methanol to give the cyclic iminium bromides **1**.

2-Benzyl-1-phenyl-3,4-dihydroisoquinolin-2-i um bromide (1l): 2.7 mmol scale, 0.840 g, 83% yield, unknown compound, pale yellow solid, m.p. 168–170 °C; ^1H NMR (400 MHz, CDCl_3)

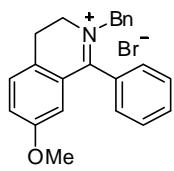


2-Benzyl-1-(*p*-tolyl)-3,4-dihydroisoquinolin-2-i um bromide (1m): 1.0 mmol scale, 0.274 g, 70% yield, unknown compound, pale yellow oil; ^1H NMR (400 MHz, CDCl_3) δ 7.79 (d, $J = 8.0$ Hz, 2H), 7.66 (t, $J = 7.5$ Hz, 1H), 7.46 (d, $J = 8.0$ Hz, 2H), 7.39 – 7.30 (m, 7H), 7.13 (d, $J = 7.9$ Hz, 1H), 5.54 (s, 2H), 4.50 (t, $J = 7.7$ Hz, 2H), 3.28 (t, $J = 7.7$ Hz, 2H), 2.49 (s, 3H); ^{13}C NMR (100 MHz, DMSO) δ 176.6, 142.6, 138.7, 137.5, 134.3, 132.9, 130.2, 129.6, 129.3, 128.8, 128.7, 128.6, 128.5, 127.3, 60.6, 50.3, 25.6, 21.6; HRMS calculated for $\text{C}_{23}\text{H}_{22}\text{N} [\text{M-Br}]^+$ 312.1747, found 312.1752.

2-Benzyl-1-(4-methoxyphenyl)-3,4-dihydroisoquinolin-2-i um bromide (1n): 1.0 mmol scale, 0.224 g, 55% yield, unknown compound, pale yellow oil; ^1H NMR (400 MHz, CDCl_3) δ 7.87 (d, $J = 8.6$ Hz, 2H), 7.65 (t, $J = 7.5$ Hz, 1H), 7.38 – 7.27 (m, 7H), 7.15 (t, $J = 7.4$ Hz, 3H), 5.57 (s, 2H), 4.46 (t, $J = 7.6$ Hz, 2H), 3.91 (s, 3H), 3.22 (t, $J = 7.5$ Hz, 2H); ^{13}C NMR (100 MHz, CDCl_3) δ 175.6, 162.8, 138.0, 137.2, 134.9, 132.6, 131.7, 129.5, 129.5, 128.4, 128.3, 128.2, 128.0, 121.1, 114.9, 61.3, 55.8, 50.1, 26.4; HRMS calculated for $\text{C}_{23}\text{H}_{22}\text{NO} [\text{M-Br}]^+$ 328.1696, found 328.1698.

2-Benzyl-1-(4-chlorophenyl)-3,4-dihydroisoquinolin-2-i um bromide (1o): 1.0 mmol scale, 0.217 g, 53% yield, unknown compound, pale yellow oil; ^1H NMR (400 MHz, CDCl_3) δ 7.99 (d, $J = 8.4$ Hz, 2H), 7.71 – 7.63 (m, 3H), 7.42 – 7.39 (m, 4H), 7.36 – 7.28 (m, 3H), 7.08 (d, $J = 7.9$ Hz, 1H), 5.43 (s, 2H), 4.47 (t, $J = 7.7$ Hz, 2H), 3.32 (t, $J = 7.6$ Hz, 2H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.9, 138.8, 137.7, 134.5, 131.9, 130.7, 129.9, 129.6, 129.6, 128.5, 128.4, 128.4, 128.3, 127.8, 127.6, 61.8, 50.4, 26.2; HRMS calculated for $\text{C}_{22}\text{H}_{19}\text{NCl} [\text{M-Br}]^+$ 332.1201, found 332.1207.

2-Benzyl-7-methoxy-1-phenyl-3,4-dihydroisoquinolin-2-i um bromide (1p): 1.0 mmol scale, 0.286 g, 64% yield, unknown compound, pale yellow oil; ^1H NMR (400 MHz, DMSO) δ 7.74 – 7.70 (m, 5H), 7.53 (d, $J = 8.4$ Hz, 1H), 7.46 – 7.40 (m, 6H), 6.40 (d, $J = 2.4$ Hz, 1H), 5.06 (s, 2H),

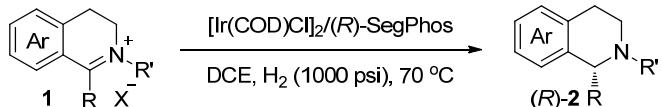


4.12 (t, J = 7.5 Hz, 2H), 3.65 (s, 3H), 3.19 (t, J = 7.7 Hz, 2H); ^{13}C NMR (100 MHz, DMSO) δ 177.0, 159.4, 133.5, 133.3, 131.4, 130.8, 130.8, 130.5, 130.4, 130.2, 130.0, 129.5, 129.5, 123.6, 120.1, 61.7, 57.0, 51.7, 25.5; HRMS calculated for $\text{C}_{23}\text{H}_{22}\text{NO} [\text{M}-\text{Br}]^+$ 328.1696, found 328.1698.

2-Benzyl-7-chloro-1-phenyl-3,4-dihydroisoquinolin-2-ium bromide (1q): 0.8 mmol scale, 0.275 g, 83% yield, unknown compound, pale yellow solid, m.p. 214–216 °C; ^1H NMR (400 MHz, DMSO) δ 7.93 (dd, J = 8.2, 2.1 Hz, 1H), 7.87 (dd, J = 7.6, 1.5 Hz, 2H), 7.81 – 7.72 (m, 3H), 7.68 (d, J = 8.2 Hz, 1H), 7.53 – 7.39 (m, 5H), 6.90 (d, J = 2.1 Hz, 1H), 5.13 (s, 2H), 4.21 (t, J = 7.6 Hz, 2H), 3.35 (t, J = 7.6 Hz, 2H); ^{13}C NMR (100 MHz, DMSO) δ 175.5, 137.6, 136.8, 132.7, 132.5, 132.5, 132.4, 130.8, 130.4, 129.9, 129.6, 129.5, 129.4, 128.8, 128.8, 61.1, 50.6, 25.0; HRMS calculated for $\text{C}_{22}\text{H}_{19}\text{ClN} [\text{M}-\text{Br}]^+$ 332.1201, found 332.1206.

3. General Procedure for Iridium-Catalyzed Asymmetric Hydrogenation

3.1. Hydrogenation of Cyclic Iminium Salts with a 3,4-Dihydroisoquinoline Core



General Procedure: A mixture of $[\text{Ir}(\text{COD})\text{Cl}]_2$ (1.3 mg, 0.002 mmol) and (R) -SegPhos (2.7 mg, 0.0044 mmol) in dichloroethane (1.0 mL) was stirred at room temperature for 10 min in glove box. Subsequently, the catalyst was transferred by a syringe to the mixture of 1-substituted 3,4-dihydroisoquinolin-2-ium salts **1** (0.2 mmol) using 4.0 mL dichloroethane. The hydrogenation was performed at 70 °C at a hydrogen pressure of 1000 psi for 24 h. After carefully releasing the hydrogen gas, the mixture was concentrated in vacuum and further purification was performed by a silica gel column eluted with hexanes/ethyl acetate to give the corresponding chiral hydrogenation products (R) -**2**.

(R)-(−)-2-Methyl-1-phenyl-1,2,3,4-tetrahydroisoquinoline (2a): 44 mg, 99% yield, 91% ee, known compound,³ pale yellow oil; $[\alpha]^{20}_D = -117.26$ (*c* 0.88, CHCl_3) [lit³: $[\alpha]^{20}_D = -103.4$ (*c* 0.50, CHCl_3) for 95% ee], $R_f = 0.32$ (hexanes/ethyl acetate 5:1); ^1H NMR (400 MHz, CDCl_3) δ 7.33 – 7.24 (m, 5H), 7.10 (q, *J* = 7.2 Hz, 2H), 6.97 (t, *J* = 7.3 Hz, 1H), 6.63 (d, *J* = 7.8 Hz, 1H), 4.24 (s, 1H), 3.31 – 3.23 (ddd, *J* = 16.3, 11.2, 5.4 Hz, 1H), 3.12 (ddd, *J* = 11.4, 5.4, 2.7 Hz, 1H), 2.82 (dt, *J* = 16.1, 2.8 Hz, 1H), 2.69 – 2.55 (m, 1H), 2.23 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 144.0, 138.6, 134.3, 129.7, 128.6, 128.3, 128.3, 127.3, 126.0, 125.7, 71.6, 52.4, 44.4, 29.6; Enantiomeric excess was determined by HPLC (OD-H column, hexane/*i*PrOH 98/2, 0.70 mL/min, 230 nm): $t_1 = 6.0$ min, $t_2 = 6.3$ min (major).

(R)-(−)-2-Methyl-1-*m*-tolyl-1,2,3,4-tetrahydroisoquinoline (2b): 47 mg, 99% yield, 90% ee, known compound,³ pale yellow oil; $[\alpha]^{20}_D = -115.77$ (*c* 0.90, CHCl_3) [lit³: $[\alpha]^{20}_D = -118.2$ (*c* 0.50, CHCl_3) for 91% ee], $R_f = 0.23$ (hexanes/ethyl acetate 5:1); ^1H NMR (400 MHz, CDCl_3) δ 7.24 – 7.20 (m, 1H), 7.13 – 7.07 (m, 5H), 6.98 (t, *J* = 7.3 Hz, 1H), 6.64 (d, *J* = 7.8 Hz, 1H), 4.19 (s, 1H), 3.32 – 3.24 (m, 1H), 3.15 – 3.11 (m, 1H), 2.83 – 2.79 (m, 1H), 2.66 – 2.59 (m, 1H), 2.31 (s, 3H), 2.23 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 144.0, 138.7, 138.0, 134.3, 130.1, 128.6, 128.3, 128.1, 128.1, 126.9, 125.9, 125.7, 71.7, 52.6, 44.5, 29.6, 21.5; Enantiomeric excess was determined by HPLC (OJ-H column, hexane/*i*PrOH 95/5, 0.70 mL/min, 230 nm): $t_1 = 6.7$ min (major), $t_2 = 7.7$ min.

(R)-(−)-2-Methyl-1-*p*-tolyl-1,2,3,4-tetrahydroisoquinoline (2c): 46 mg, 97% yield, 90% ee, known compound,³ pale yellow oil; $[\alpha]^{20}_D = -114.66$ (*c* 0.90, CHCl_3) [lit³: $[\alpha]^{20}_D = -111.0$ (*c* 0.50, CHCl_3) for 94% ee], $R_f = 0.25$ (hexanes/ethyl acetate 5:1); ^1H NMR (400 MHz, CDCl_3) δ 7.16 – 7.06 (m, 6H), 6.97 (t, *J* = 7.3 Hz, 1H), 6.64 (d, *J* = 7.8 Hz, 1H), 4.20 (s, 1H), 3.30 – 3.22 (M, 1H), 3.11 (ddd, *J* = 11.5, 5.5, 2.8 Hz, 1H), 2.81 (dt, *J* = 16.1, 2.8 Hz, 1H), 2.62 (td, *J* = 11.3, 3.8 Hz, 1H), 2.33 (s, 3H), 2.23 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 140.9, 138.8, 136.9, 134.3, 129.5, 129.0, 128.6, 128.3, 125.9, 125.7, 71.3, 52.4, 44.4, 29.6, 21.2; Enantiomeric excess was determined by HPLC (OJ-H column, hexane/*i*PrOH 95/5, 0.70 mL/min, 230 nm): $t_1 = 6.8$ min (major), $t_2 = 9.7$ min.

(R)-(−)-1-(4-Methoxyphenyl)-2-methyl-1,2,3,4-tetrahydroisoquinoline (2d): 46 mg, 91% yield, 89% ee, known compound,³ pale yellow solid; $[\alpha]^{20}_D = -110.96$ (*c* 0.62, CHCl_3) [lit³: $[\alpha]^{20}_D$

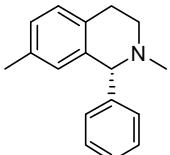
$\delta = -111.2$ (c 0.50, CHCl_3) for 91% ee], $R_f = 0.25$ (hexanes/ethyl acetate 2:1); ^1H NMR (400 MHz, CDCl_3) δ 7.17 (d, $J = 8.5$ Hz, 2H), 7.09 (q, $J = 7.6$ Hz, 2H), 6.98 (t, $J = 7.2$ Hz, 1H), 6.85 (d, $J = 8.6$ Hz, 2H), 6.64 (d, $J = 7.8$ Hz, 1H), 4.20 (s, 1H), 3.80 (s, 3H), 3.29 – 3.21 (m, 1H), 3.11 (ddd, $J = 11.4, 5.4, 2.8$ Hz, 1H), 2.86 – 2.77 (m, 1H), 2.62 (td, $J = 11.3, 3.8$ Hz, 1H), 2.23 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 158.8, 138.9, 136.0, 134.3, 130.6, 128.6, 128.3, 125.9, 125.6, 125.6, 113.6, 70.9, 55.2, 52.4, 44.3, 29.5. Enantiomeric excess was determined by HPLC (OD-H column, hexane/ $i\text{PrOH}$ 90/10, 1.0 mL/min, 230 nm): $t_1 = 4.1$ min, $t_2 = 4.4$ min (major).

(R)-(-)-1-(4-Fluorophenyl)-2-methyl-1,2,3,4-tetrahydroisoquinoline (2e): 45 mg, 94% yield, 90% ee, unknown compound, colorless oil; $[\alpha]^{20}_D = -114.38$ (c 0.82, CHCl_3), $R_f = 0.30$ (hexanes/ethyl acetate 5:1); ^1H NMR (400 MHz, CDCl_3) δ 7.25 – 7.21 (m, 2H), 7.13 – 7.07 (m, 2H), 7.02 – 6.96 (m, 3H), 6.60 (d, $J = 7.8$ Hz, 1H), 4.22 (s, 1H), 3.29 – 3.21 (m, 1H), 3.11 (ddd, $J = 11.5, 5.5, 2.7$ Hz, 1H), 2.81 (dt, $J = 16.1, 3.0$ Hz, 1H), 2.63 (td, $J = 11.3, 3.8$ Hz, 1H), 2.22 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 162.1 (d, $J = 245.2$ Hz), 139.9 (d, $J = 3.2$ Hz), 138.4, 134.4, 131.1 (d, $J = 7.9$ Hz), 128.5, 128.4, 126.1, 125.7, 115.1 (d, $J = 21.3$ Hz), 70.8, 52.4, 44.3, 29.5; ^{19}F NMR (376 MHz, CDCl_3) δ -115.41; Enantiomeric excess was determined by HPLC (OD-H column, hexane/ $i\text{PrOH}$ 95/5, 1.0 mL/min, 230 nm): $t_1 = 3.8$ min, $t_2 = 4.2$ min (major); HRMS calculated for $\text{C}_{16}\text{H}_{17}\text{NF}$ $[\text{M}+\text{H}]^+$ 242.1340, found 242.1350.

(R)-(-)-1-(4-Chlorophenyl)-2-methyl-1,2,3,4-tetrahydroisoquinoline (2f): 50 mg, 97% yield, 87% ee, known compound,³ white solid; $[\alpha]^{20}_D = -118.55$ (c 0.90, CHCl_3) [lit³: $[\alpha]^{20}_D = -130.0$ (c 0.50, CHCl_3) for 91% ee], $R_f = 0.35$ (hexanes/ethyl acetate 5:1); ^1H NMR (400 MHz, CDCl_3) δ 7.28 (d, $J = 8.4$ Hz, 2H), 7.22 (t, $J = 7.9$ Hz, 2H), 7.13 – 7.07 (m, 2H), 6.98 (t, $J = 7.2$ Hz, 1H), 6.59 (d, $J = 7.8$ Hz, 1H), 4.21 (s, 1H), 3.29 – 3.20 (m, 1H), 3.10 (ddd, $J = 11.4, 5.4, 2.6$ Hz, 1H), 2.88 – 2.74 (m, 1H), 2.62 (td, $J = 11.3, 3.7$ Hz, 1H), 2.22 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 142.7, 138.1, 134.3, 133.0, 130.9, 128.5, 128.5, 128.4, 126.1, 125.8, 70.8, 52.3, 44.3, 29.5; Enantiomeric excess was determined by HPLC (OD-H column, hexane/ $i\text{PrOH}$ 90/10, 1.0 mL/min, 230 nm): $t_1 = 3.7$ min, $t_2 = 4.1$ min (major).

(R)-(-)-1-(4-Bromophenyl)-2-methyl-1,2,3,4-tetrahydroisoquinoline (2g): 59 mg, 98% yield, 87% ee, unknown compound, white solid, m.p. 97–99 °C; $[\alpha]^{20}_D = -106.86$ (c 1.18, CHCl_3), $R_f = 0.30$ (hexanes/ethyl acetate 5:1); ^1H NMR (400 MHz, CDCl_3) δ 7.45 (d, $J = 8.3$ Hz, 2H), 7.18 – 7.09 (m, 4H), 7.00 (t, $J = 7.1$ Hz, 1H), 6.60 (d, $J = 7.8$ Hz, 1H), 4.22 (s, 1H), 3.30 – 7.21 (m, 1H), 3.12 (ddd, $J = 11.4, 5.4, 2.6$ Hz, 1H), 2.88 – 2.76 (m, 1H), 2.64 (td, $J = 11.4, 3.8$ Hz, 1H), 2.23 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 143.3, 138.0, 134.3, 131.5, 131.3, 128.5, 128.4, 126.2, 125.8, 121.2, 70.9, 52.3, 44.4, 29.5; Enantiomeric excess was determined by HPLC (OD-H column, hexane/ $i\text{PrOH}$ 98/2, 0.7 mL/min, 230 nm): $t_1 = 6.0$ min, $t_2 = 7.2$ min (major); HRMS calculated for $\text{C}_{16}\text{H}_{17}\text{NBr}$ $[\text{M}+\text{H}]^+$ 302.0539, found 302.0541.

(R)-(-)-2,7-Dimethyl-1-phenyl-1,2,3,4-tetrahydroisoquinoline (2h): 46 mg, 97% yield, 93% ee, known compound,³ pale yellow oil; $[\alpha]^{20}_D = -72.07$ (c 0.82, CHCl_3) [lit³: $[\alpha]^{20}_D = -73.8$ (c 0.50, CHCl_3) for 93% ee], $R_f = 0.50$ (hexanes/ethyl acetate 2:1); ^1H NMR (400 MHz, CDCl_3) δ 7.33 – 7.23 (m, 5H), 7.01 (d, $J = 7.7$ Hz, 1H), 6.91 (d, $J = 7.7$ Hz, 1H), 6.44 (s, 1H), 4.19 (s, 1H), 3.25 –


 3.17 (m, 1H), 3.10 (ddd, J = 11.3, 5.3, 2.8 Hz, 1H), 2.83 – 2.73 (m, 1H), 2.60 (td, J = 11.2, 3.8 Hz, 1H), 2.22 (s, 3H), 2.12 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 144.1, 138.4, 135.0, 131.4, 129.7, 129.0, 128.3, 128.2, 127.3, 126.9, 71.5, 52.4, 44.4, 29.2, 21.1; Enantiomeric excess was determined by HPLC (OG column, hexane/ $i\text{PrOH}$ 99.6/0.4, 0.7 mL/min, 230 nm): t_1 = 5.9 min, t_2 = 6.5 min (major).

(R)-(-)-7-Methoxy-2-methyl-1-phenyl-1,2,3,4-tetrahydroisoquinoline (2i): 46 mg, 91% yield, 86% ee, known compound,³ pale yellow oil; $[\alpha]^{20}_{\text{D}} = -52.50$ (c 0.80, CHCl_3) [lit³: $[\alpha]^{20}_{\text{D}} = -58.4$ (c 0.50, CHCl_3) for 86% ee], $R_f = 0.35$ (hexanes/ethyl acetate 5:1); ^1H NMR (400 MHz, CDCl_3) δ 7.32 – 7.23 (m, 5H), 7.04 (d, J = 8.4 Hz, 1H), 6.68 (dd, J = 8.4, 2.6 Hz, 1H), 6.17 (d, J = 2.4 Hz, 1H), 4.19 (s, 1H), 3.59 (s, 3H), 3.22 – 3.08 (m, 2H), 2.77 (dd, J = 15.7, 3.2 Hz, 1H), 2.60 (td, J = 10.9, 3.8 Hz, 1H), 2.23 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 157.5, 143.8, 139.7, 129.6, 129.2, 128.3, 127.3, 126.7, 113.8, 112.0, 71.6, 55.1, 52.5, 44.4, 28.7; Enantiomeric excess was determined by HPLC (OJ-H column, hexane/ $i\text{PrOH}$ 90/10, 0.7 mL/min, 230 nm): t_1 = 12.3 min (major), t_2 = 17.4 min.

(R)-(-)-7-Chloro-2-methyl-1-phenyl-1,2,3,4-tetrahydroisoquinoline (2j): 48 mg, 93% yield, 84% ee, unknown compound, colourless oil; $[\alpha]^{20}_{\text{D}} = -34.79$ (c 0.96, CHCl_3), $R_f = 0.45$ (hexanes/ethyl acetate 2:1); ^1H NMR (400 MHz, CDCl_3) δ 7.36 – 7.22 (m, 5H), 7.13 – 6.99 (m, 2H), 6.61 (s, 1H), 4.18 (s, 1H), 3.25 – 3.06 (m, 2H), 2.84 – 2.73 (m, 1H), 2.63 – 2.56 (m, 1H), 2.21 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 143.1, 140.5, 132.9, 131.3, 129.7, 129.6, 128.5, 128.4, 127.6, 126.2, 71.21, 52.1, 44.3, 29.0; Enantiomeric excess was determined by HPLC (AD-H column, hexane/ $i\text{PrOH}$ 95/5, 0.8 mL/min, 230 nm): t_1 = 4.7 min, t_2 = 5.1 min (major); HRMS calculated for $\text{C}_{16}\text{H}_{17}\text{ClN}$ [$\text{M}+\text{H}]^+$ 258.1044, found 258.1047.

(R)-(+)-2-Methyl-1-phenethyl-1,2,3,4-tetrahydroisoquinoline (2k): 50 mg, 99% yield, 26% ee, known compound,⁴ pale yellow oil; $[\alpha]^{20}_{\text{D}} = +1.20$ (c 1.00, CHCl_3), $R_f = 0.30$ (hexanes/ethyl acetate 2:1); ^1H NMR (400 MHz, CDCl_3) δ 7.28 – 7.24 (m, 2H), 7.22 – 7.04 (m, 7H), 3.50 (t, J = 5.3 Hz, 1H), 3.21 – 3.10 (m, 1H), 2.89 – 2.66 (m, 4H), 2.59 – 2.50 (m, 1H), 2.48 (s, 3H), 2.16 – 2.01 (m, 2H); ^{13}C NMR (100 MHz, CDCl_3) δ 143.0, 138.1, 134.9, 128.8, 128.5, 128.3, 127.1, 125.9, 125.8, 125.6, 63.1, 48.5, 42.9, 36.7, 31.4, 26.3; Enantiomeric excess was determined by HPLC (OD-H column, hexane/ $i\text{PrOH}$ 98/2, 0.7 mL/min, 230 nm): t_1 = 6.6 min, t_2 = 7.2 min (major).

(R)-(-)-2-Benzyl-1-phenyl-1,2,3,4-tetrahydroisoquinoline (2l): The hydrogenation was performed in dichloroethane (3 mL) at 50 °C; 57 mg, 95% yield, 96% ee, known compound,³ white solid; $[\alpha]^{20}_{\text{D}} = -100.72$ (c 1.10, CHCl_3) [lit³: $[\alpha]^{20}_{\text{D}} = -94.6$ (c 0.50, CHCl_3) for 95% ee], $R_f = 0.41$ (hexanes/ethyl acetate 20:1); ^1H NMR (400 MHz, CDCl_3) δ 7.38 (d, J = 7.4 Hz, 2H), 7.34 – 7.20 (m, 8H), 7.12 – 7.07 (m, 2H), 7.02 – 6.98 (m, 1H), 6.72 (d, J = 7.8 Hz, 1H), 4.60 (s, 1H), 3.81 (d, J = 13.6 Hz, 1H), 3.24 (d, J = 13.6 Hz, 1H), 3.12 – 3.02 (m, 2H), 2.78 – 2.74 (m, 1H), 2.54 – 2.48 (m, 1H); ^{13}C NMR (100 MHz, CDCl_3) δ 144.5, 139.6, 138.6, 134.9, 129.7, 128.9, 128.8, 128.5, 128.3, 128.2, 127.3, 126.9, 125.9, 125.7, 68.9, 58.9, 47.4, 29.3; Enantiomeric excess was determined by HPLC (AD-H column, hexane/ $i\text{PrOH}$ 90/10, 1.0 mL/min, 230 nm): t_1 = 3.5 min (major), t_2 = 3.9 min.

(R)-(-)-2-Benzyl-1-(*p*-tolyl)-1,2,3,4-tetrahydroisoquinoline (2m): The hydrogenation was performed in dichloroethane (3 mL) at 50 °C; 61 mg, 97% yield, 95% ee, unknown compound, white

solid, m.p. 97-99 °C; $[\alpha]^{20}_D = -92.45$ (*c* 1.22, CHCl₃), R_f = 0.58 (hexanes/ethyl acetate 10:1); ¹H NMR (400 MHz, CDCl₃) δ 7.31 – 7.24 (m, 6H), 7.21 – 7.18 (m, 1H), 7.12 – 7.05 (m, 4H), 7.00 – 6.96 (m, 1H), 6.73 (d, *J* = 7.8 Hz, 1H), 4.56 (s, 1H), 3.81 (d, *J* = 13.6 Hz, 1H), 3.22 (d, *J* = 13.6 Hz, 1H), 3.11 – 3.01 (m, 2H), 2.78 – 2.72 (m, 1H), 2.52 – 2.45 (m, 1H), 2.31 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 141.4, 139.7, 138.9, 136.8, 134.9, 129.6, 129.1, 128.9, 128.8, 128.5, 128.2, 126.8, 125.9, 125.7, 68.6, 58.9, 47.4, 29.4, 21.2; Enantiomeric excess was determined by HPLC (AD-H column, hexane/ⁱPrOH 98/2, 0.7 mL/min, 230 nm): t₁ = 5.2 min (major), t₂ = 5.9 min; HRMS calculated for C₂₃H₂₄N [M+H]⁺ 314.1903, found 314.1908.

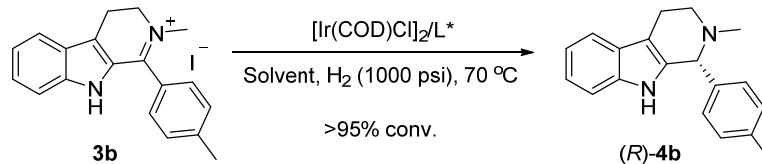
(R)-(-)-2-Benzyl-1-(4-methoxyphenyl)-1,2,3,4-tetrahydroisoquinoline (2n): The hydrogenation was performed in dichloroethane (3 mL) at 50 °C; 63 mg, 96% yield, 96% ee, known compound,² colorless oil; $[\alpha]^{20}_D = -98.39$ (*c* 1.00, CHCl₃), [lit²: $[\alpha]^{29}_D = -94.6$ (*c* 2.05, CHCl₃) for 94% ee], R_f = 0.45 (hexanes/ethyl acetate 10:1); ¹H NMR (400 MHz, CDCl₃) δ 7.31 – 7.25 (m, 6H), 7.22 – 7.18 (m, 1H), 7.10 – 7.06 (m, 2H), 7.01 – 6.96 (m, 1H), 6.87 – 6.83 (m, 2H), 6.73 (d, *J* = 7.8 Hz, 1H), 4.55 (s, 1H), 3.82 (d, *J* = 13.6 Hz, 1H), 3.77 (s, 3H), 3.22 (d, *J* = 13.6 Hz, 1H), 3.11 – 3.00 (m, 2H), 2.75 (dt, *J* = 7.2, 4.5 Hz, 1H), 2.53 – 2.46 (m, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 158.8, 139.7, 138.9, 136.5, 134.9, 130.7, 128.9, 128.8, 128.5, 128.2, 126.8, 125.9, 125.6, 113.7, 68.2, 58.8, 55.3, 47.4, 29.3; Enantiomeric excess was determined by HPLC (AD-H column, hexane/ⁱPrOH 99.2/0.8, 0.7 mL/min, 230 nm): t₁ = 8.9 min (major), t₂ = 10.1 min.

(R)-(-)-2-Benzyl-1-(4-chlorophenyl)-1,2,3,4-tetrahydroisoquinoline (2o): The hydrogenation was performed in dichloroethane (3 mL) at 50 °C; 62 mg, 93% yield, 95% ee, unknown compound, white solid, mp 108-109 °C; $[\alpha]^{20}_D = -92.54$ (*c* 1.14, CHCl₃), R_f = 0.43 (hexanes/ethyl acetate 20:1); ¹H NMR (400 MHz, CDCl₃) δ 7.37 – 7.24 (m, 9H), 7.15 – 7.12 (m, 2H), 7.07 – 7.03 (m, 1H), 6.73 (d, *J* = 7.7 Hz, 1H), 4.63 (s, 1H), 3.82 (d, *J* = 13.5 Hz, 1H), 3.28 (d, *J* = 13.5 Hz, 1H), 3.15 – 3.04 (m, 2H), 2.80 (dt, *J* = 6.8, 4.2 Hz, 1H), 2.58 – 2.52 (m, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 143.9, 140.1, 138.8, 135.7, 133.8, 131.8, 129.6, 129.5, 129.5, 129.3, 129.1, 127.8, 127.0, 126.6, 68.9, 59.7, 48.1, 30.0; Enantiomeric excess was determined by HPLC (AD-H column, hexane/ⁱPrOH 98/2, 0.7 mL/min, 230 nm): t₁ = 5.9 min (major), t₂ = 6.3 min; HRMS calculated for C₂₂H₂₁NCl [M+H]⁺ 334.1357, found 334.1360.

(R)-(-)-2-Benzyl-7-methoxy-1-phenyl-1,2,3,4-tetrahydroisoquinoline (2p): The hydrogenation was performed in dichloroethane (3 mL) at 50 °C; 60 mg, 91% yield, 93% ee, unknown compound, colorless oil; $[\alpha]^{20}_D = -41.87$ (*c* 1.12, CHCl₃), R_f = 0.35 (hexanes/ethyl acetate 20:1); ¹H NMR (400 MHz, CDCl₃) δ 7.39 – 7.22 (m, 10H), 7.05 (d, *J* = 8.4 Hz, 1H), 6.70 (dd, *J* = 8.4, 2.5 Hz, 1H), 6.29 (d, *J* = 2.6 Hz, 1H), 4.58 (s, 1H), 3.82 (d, *J* = 13.6 Hz, 1H), 3.63 (s, 3H), 3.27 (d, *J* = 13.5 Hz, 1H), 3.10 (dt, *J* = 11.3, 4.7 Hz, 1H), 3.05 – 2.94 (m, 1H), 2.73 (dt, *J* = 15.8, 3.9 Hz, 1H), 2.51 (ddd, *J* = 11.6, 9.9, 4.0 Hz, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 157.5, 144.2, 139.6, 139.5, 129.6, 129.4, 128.8, 128.3, 128.2, 127.3, 127.1, 126.8, 114.0, 112.1, 68.8, 58.9, 55.1, 47.4, 28.2; Enantiomeric excess was determined by HPLC (AD-H column, hexane/ⁱPrOH 98/2, 0.7 mL/min, 230 nm): t₁ = 7.0 min (major), t₂ = 9.1 min; HRMS calculated for C₂₃H₂₄NO [M+H]⁺ 330.1852, found 330.1854.

(R)-(-)-2-Benzyl-7-chloro-1-phenyl-1,2,3,4-tetrahydroisoquinoline (2q): The hydrogenation was performed in dichloroethane (3 mL) at 50 °C; 64 mg, 96% yield, 95% ee, unknown compound, colorless oil; $[\alpha]^{20}_D = -27.89$ (c 1.28, CHCl₃), $R_f = 0.58$ (hexanes/ethyl acetate 10:1); ¹H NMR (400 MHz, CDCl₃) δ 7.37 – 7.18 (m, 10H), 7.10 – 6.97 (m, 2H), 6.71 (d, $J = 1.0$ Hz, 1H), 4.54 (s, 1H), 3.78 (d, $J = 13.5$ Hz, 1H), 3.23 (d, $J = 13.5$ Hz, 1H), 3.13 – 3.03 (m, 1H), 3.02 – 2.95 (m, 1H), 2.75 – 2.69 (m, 1H), 2.50 – 2.44 (m, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 143.5, 140.4, 139.3, 133.4, 131.2, 129.9, 129.6, 128.7, 128.5, 128.2, 127.6, 127.0, 126.2, 68.5, 58.7, 47.1, 28.7; Enantiomeric excess was determined by HPLC (AD-H column, hexane/ⁱPrOH 95/5, 0.8 mL/min, 230 nm): $t_1 = 4.5$ min (major), $t_2 = 5.2$ min; HRMS calculated for C₂₂H₂₁ClN [M+H]⁺ 334.1357, found 330. 334.1355.

Optimization of Iridium-Catalyzed Asymmetric Hydrogenation of Cyclic Iminium Salts with a 4,9-Dihydro-3-H-β-carboline Core



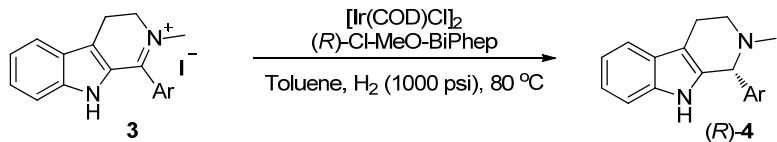
Entry ^a	Solvent	L*	Ee (%) ^c
1	DCE	L3	62
2	Toluene	L3	79
3	Dioxane	L3	76
4	THF	L3	73
5	ⁱ PrOH	L3	47
6 ^b	Toluene	L1	76
7	Toluene	L4	77
8	Toluene	L5	85
9 ^d	Toluene	L5	84

L1: (R)-SynPhos **L2:** (R)-BINAP **L3:** (R)-SegPhos **L4:** (R)-DifluorPhos **L5:** (R)-Cl-MeO-BiPhep

^a Reaction condition: **3b** (0.1 mmol), [Ir(COD)Cl]₂ (1 mol%), **L*** (2.2 mol%), H₂ (1000 psi), solvent (3.0 mL), 24 h, 70 °C; The reaction was conducted with full conversion (determined by ¹H NMR analysis with 1,3,5-trimethoxy-benzene as the internal standard). ^b 66% conversion ^c Determined by HPLC analysis. ^d 80 °C.

With the optimization of reaction conditions, (R)-Cl-MeO-BiPhep and toluene turned out to be the best ligand and solvent for hydrogenation of 2-methyl-1-(*p*-tolyl)-4,9-dihydro-3*H*-pyrido-[3,4-*b*]indol-2-ium iodide, providing the desired product in full conversion and 84% ee.

3.2. Hydrogenation of Cyclic Iminium Salts with 4,9-Dihydro-3-H- β -carboline Core



General Procedure: A mixture of $[\text{Ir}(\text{COD})\text{Cl}]_2$ (1.4 mg, 0.002 mmol) and $(R)\text{-Cl-MeO-BiPhep}$ (3.0 mg, 0.0044 mmol) in toluene (1.0 mL) was stirred at room temperature for 10 min in glove box. Subsequently, the catalyst was transferred by a syringe to the mixture of 1-substituted 2-methyl-2,3,4,9-tetrahydro-1*H*-pyrido[3,4-*b*]indol-2-ium iodide **3** (0.2 mmol) using 4.0 mL toluene. The hydrogenation was performed at 80°C and at a hydrogen pressure of 1000 psi for 24 h. After carefully releasing the hydrogen, the mixture was concentrated in vacuum and further purification was performed by a silica gel column eluted with hexanes/ethyl acetate to give the chiral hydrogenation product **(R)-4**.

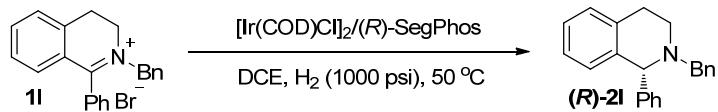
(R)-(-)-2-Methyl-1-phenyl-2,3,4,9-tetrahydro-1*H*-pyrido[3,4-*b*]indole (4a): 41 mg, 78% yield, 88% ee, known compound,⁵ pale yellow solid; $[\alpha]^{20}_D = -89.20$ (*c* 0.76, CHCl_3), $R_f = 0.45$ (hexanes/ethyl acetate 2:1); ^1H NMR (400 MHz, CDCl_3) δ 7.53 – 7.50 (m, 1H), 7.43 – 7.28 (m, 5H), 7.20 (brs, 1H), 7.16 – 7.11 (m, 1H), 7.11 – 7.04 (m, 2H), 4.28 (s, 1H), 3.34 – 3.15 (m, 1H), 3.15 – 3.06 (m, 1H), 2.93 – 2.79 (m, 1H), 2.74 (td, *J* = 11.0, 4.2 Hz, 1H), 2.32 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 140.8, 136.3, 134.9, 129.2, 128.8, 128.2, 127.1, 121.5, 119.4, 118.3, 110.8, 108.9, 67.4, 53.3, 43.6, 21.6; Enantiomeric excess was determined by HPLC (IC column, hexane/ $^i\text{PrOH}$ 70/30, 0.8 mL/min, 230 nm): $t_1 = 5.1$ min, $t_2 = 9.9$ min (major).

(R)-(-)-2-Methyl-1-p-tolyl-2,3,4,9-tetrahydro-1*H*-pyrido[3,4-*b*]indole (4b): 54 mg, 98% yield, 84% ee, unknown compound, pale yellow solid, m.p. 158–160 $^\circ\text{C}$; $[\alpha]^{20}_D = -85.19$ (*c* 1.04, CHCl_3), $R_f = 0.45$ (hexanes/ethyl acetate 2:1); ^1H NMR (400 MHz, CDCl_3) δ 7.53 – 7.51 (m, 1H), 7.24 – 7.20 (m, 3H), 7.18 – 7.13 (m, 3H), 7.13 – 7.05 (m, 2H), 4.26 (s, 1H), 3.31 – 3.19 (m, 1H), 3.14 – 3.06 (m, 1H), 2.88 – 2.67 (m, 2H), 2.34 (s, 3H+3H); ^{13}C NMR (100 MHz, CDCl_3) δ 138.0, 137.7, 136.3, 135.1, 129.4, 129.1, 127.2, 121.5, 119.4, 118.3, 110.8, 108.8, 67.1, 53.3, 43.6, 21.7, 21.2; Enantiomeric excess was determined by HPLC (IC column, hexane/ $^i\text{PrOH}$ 70/30, 0.7 mL/min, 230 nm): $t_1 = 6.3$ min, $t_2 = 11.3$ min (major); HRMS calculated for $\text{C}_{23}\text{H}_{24}\text{NO}$ [$\text{M}+\text{H}]^+$ 277.1705 found 277.1714.

(R)-(-)-1-(4-Methoxyphenyl)-2-methyl-2,3,4,9-tetrahydro-1*H*-pyrido[3,4-*b*]indole (4c): 48 mg, 83% yield, 88% ee, unknown compound, colorless oil; $[\alpha]^{20}_D = -69.88$ (*c* 0.90, CHCl_3), $R_f = 0.40$ (hexanes/ethyl acetate 2:1); ^1H NMR (400 MHz, CDCl_3) δ 7.57 – 7.49 (m, 1H), 7.26 – 7.20 (m, 3H), 7.19 – 7.13 (m, 1H), 7.13 – 7.04 (m, 2H), 6.93 – 6.83 (m, 2H), 4.24 (s, 1H), 3.81 (s, 3H), 2.86 – 2.81 (m, 1H), 3.13 – 3.06 (m, 1H), 2.83 (ddd, *J* = 13.3, 4.0, 2.0 Hz, 1H), 2.73 (td, *J* = 11.0, 4.2 Hz, 1H), 2.31 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 159.5, 136.3, 135.3, 132.8, 130.3, 127.2, 121.5, 119.4, 118.3, 114.1, 110.8, 108.8, 66.7, 55.3, 53.3, 43.5, 21.7; Enantiomeric excess was determined by HPLC (IC column, hexane/ $^i\text{PrOH}$ 70/30, 0.8 mL/min, 230 nm): $t_1 = 6.7$ min, $t_2 = 12.2$ (major) min; HRMS calculated for $\text{C}_{23}\text{H}_{24}\text{NO}$ [$\text{M}+\text{H}]^+$

293.1648, found 293.1646.

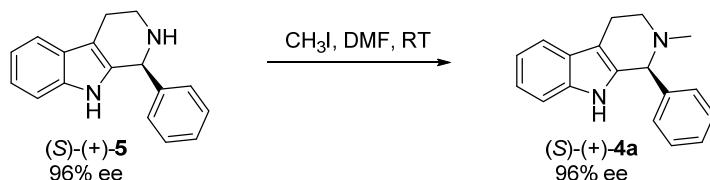
4. Asymmetric Hydrogenation at Gram Scale



A mixture of $[\text{Ir}(\text{COD})\text{Cl}]_2$ (26.8 mg, 0.04 mmol) and (*R*)-SegPhos (53.7 mg, 0.088 mmol) in dichloroethane (5.0 mL) was stirred at room temperature for 10 min in glove box. Subsequently, the catalyst was transferred by a syringe to the mixture of **1l** (1.513 g, 4.0 mmol) using 35 mL of dichloroethane. The hydrogenation was performed at 50°C at a hydrogen pressure of 1000 psi for 24 h. After carefully releasing the hydrogen, the mixture was concentrated in vacuum and further purification was performed by a silica gel column eluted with hexanes/ethyl acetate to give the product (*R*)-**2l** 1.141 g in 95% yield and 96% ee. Enantiomeric excess was determined by HPLC (AD-H column, hexane/ $^1\text{PrOH}$ 90/10, 1.0 mL/min, 230 nm): $t_1 = 3.6$ min (major), $t_2 = 3.9$ min.

5. The Determination of Absolute Configuration of (-)-4a

The absolute configuration of (-)-**4a** is unknown. Considering that the absolute configuration of compound (*S*)-(+)**5** is known,⁶ *N*-methylation of (*S*)-(+)**5** was performed to deliver the product (*S*)-(+)-**4a**, which is the same structure with our hydrogenation product **4a**. By comparison with the optical rotation datum, the absolute configuration of hydrogenation product **4a** could be unambiguously assigned.



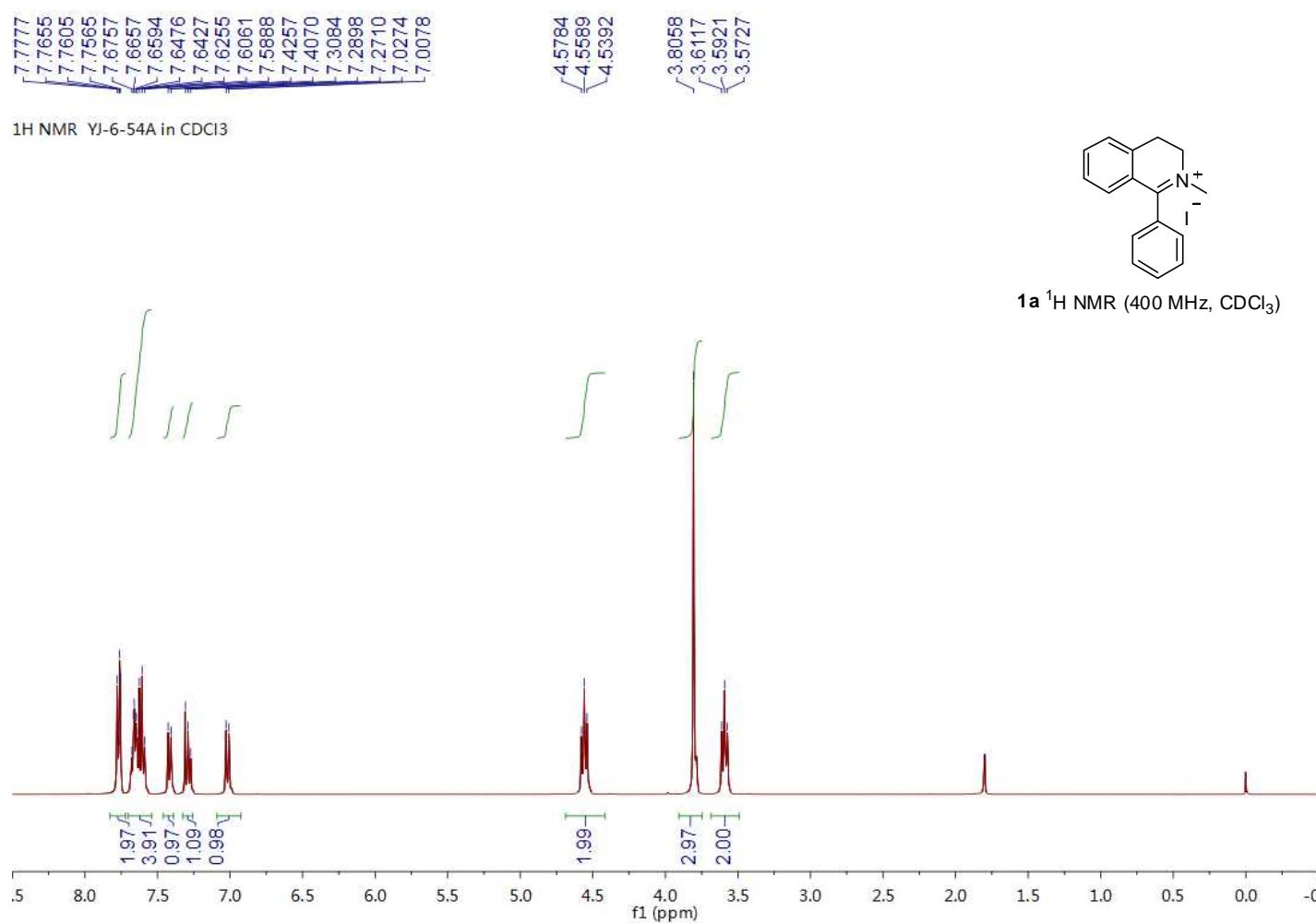
Iodomethane (22 mg, 0.30 mmol) was slowly added to the solution of (*S*)-(+)-**5** (37 mg, 0.15 mmol, 96% ee) and DMF (2 mL). After stirring 2 h, saturated aqueous sodium bicarbonate was added to the mixture until pH reaching to 8-9. The mixture was extracted with dichloromethane twice and the combined organic extracts dried over sodium sulfate. The resulting mixture was concentrated in vacuum and further purification was performed by a silica gel column eluted with hexanes/ethyl acetate to give product (*S*)-(+)-2-methyl-1-phenyl-2,3,4,9-tetrahydro-1*H*-pyrido-[3,4-*b*]indole **4a**. The optical rotation datum of the corresponding product is opposite to our hydrogenation product of iminium salt **3a**. By comparison with the optical rotation datum, the absolute configuration of hydrogenation product (-)-**4a** was assigned as (*R*)-(-)-**4a**.

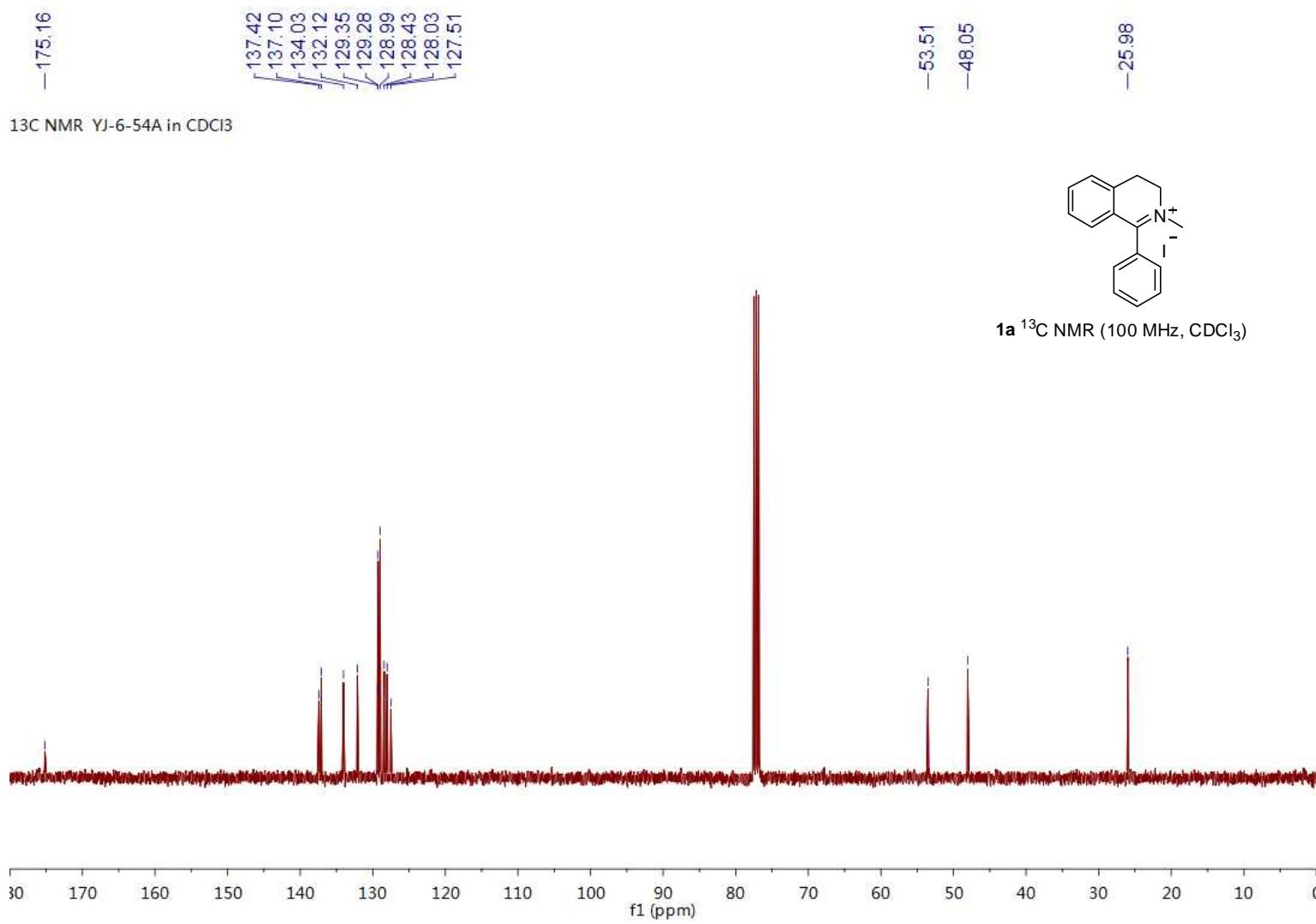
(*S*)-(+)-2-Methyl-1-phenyl-2,3,4,9-tetrahydro-1*H*-pyrido[3,4-*b*]indole ((*S*)-(+)-4a**):** 20 mg, yield: 51%, ee: 96%, the known compound,⁵ pale yellow solid; $[\alpha]^{20}_D = +102.49$ (*c* 0.16, CHCl₃), R_f = 0.32 (hexanes/ethyl acetate 2:1); ¹H NMR (400 MHz, CDCl₃) δ 7.53 – 7.50 (m, 1H), 7.39 – 7.27 (m, 5H), 7.19 (brs, 1H), 7.15 – 7.05 (m, 3H), 4.27 (s, 1H), 3.22 (ddd, *J* = 11.4, 5.4, 2.3 Hz, 1H), 3.13 – 3.05 (m, 1H), 2.86 – 2.81 (m, 1H), 2.74 (td, *J* = 11.0, 4.2 Hz, 1H), 2.31 (s, 3H); Enantiomeric excess was determined by HPLC (IC column, hexane/ⁱPrOH 70/30, 0.8 mL/min, 230 nm): t₁ = 5.2 min (major), t₂ = 10.1 min.

6. References

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6. Copy of NMR and HPLC for Compounds



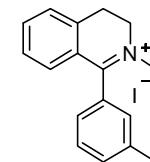


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7.5545
7.4178
7.4061
7.3965
7.3976
6.9567

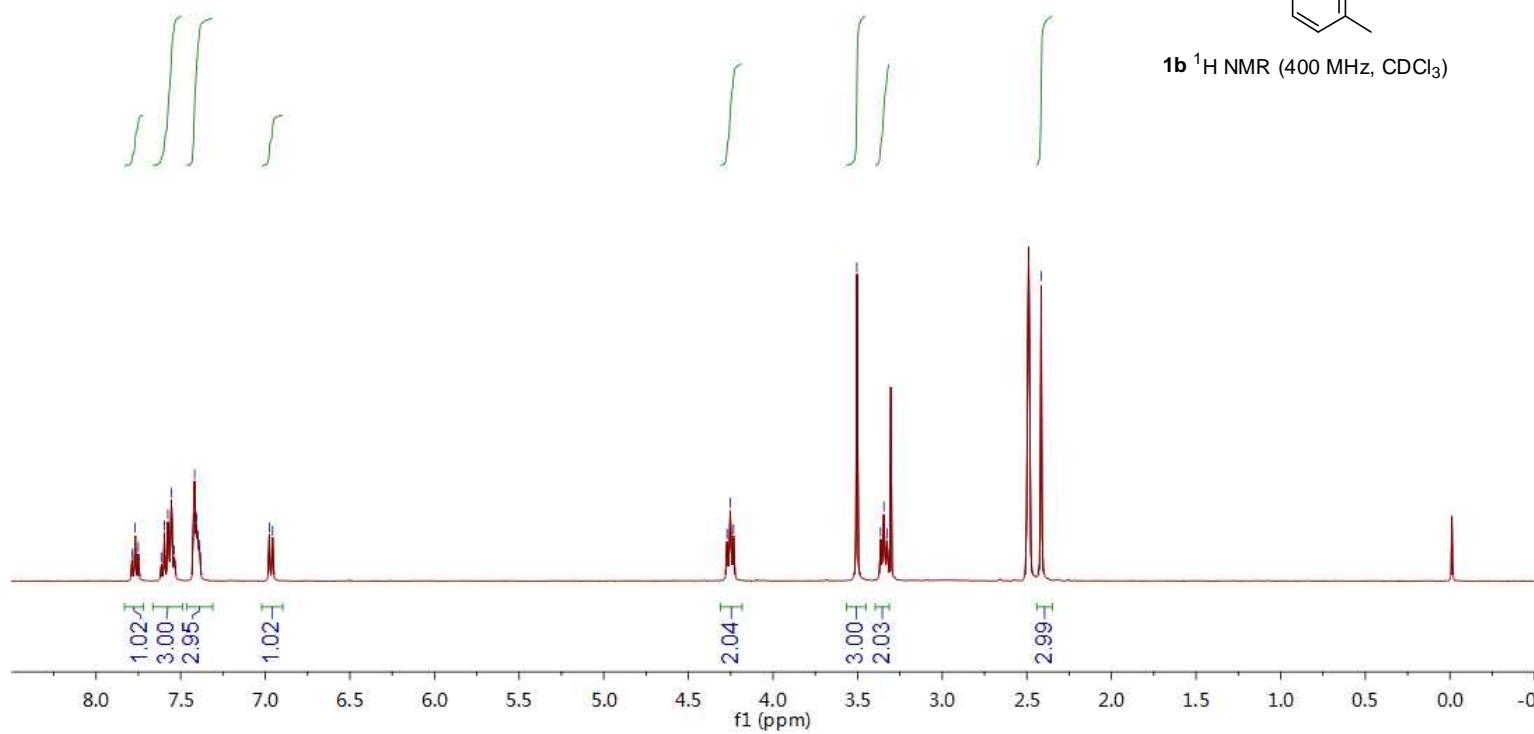
¹H NMR YJ-6-85C in DMSO

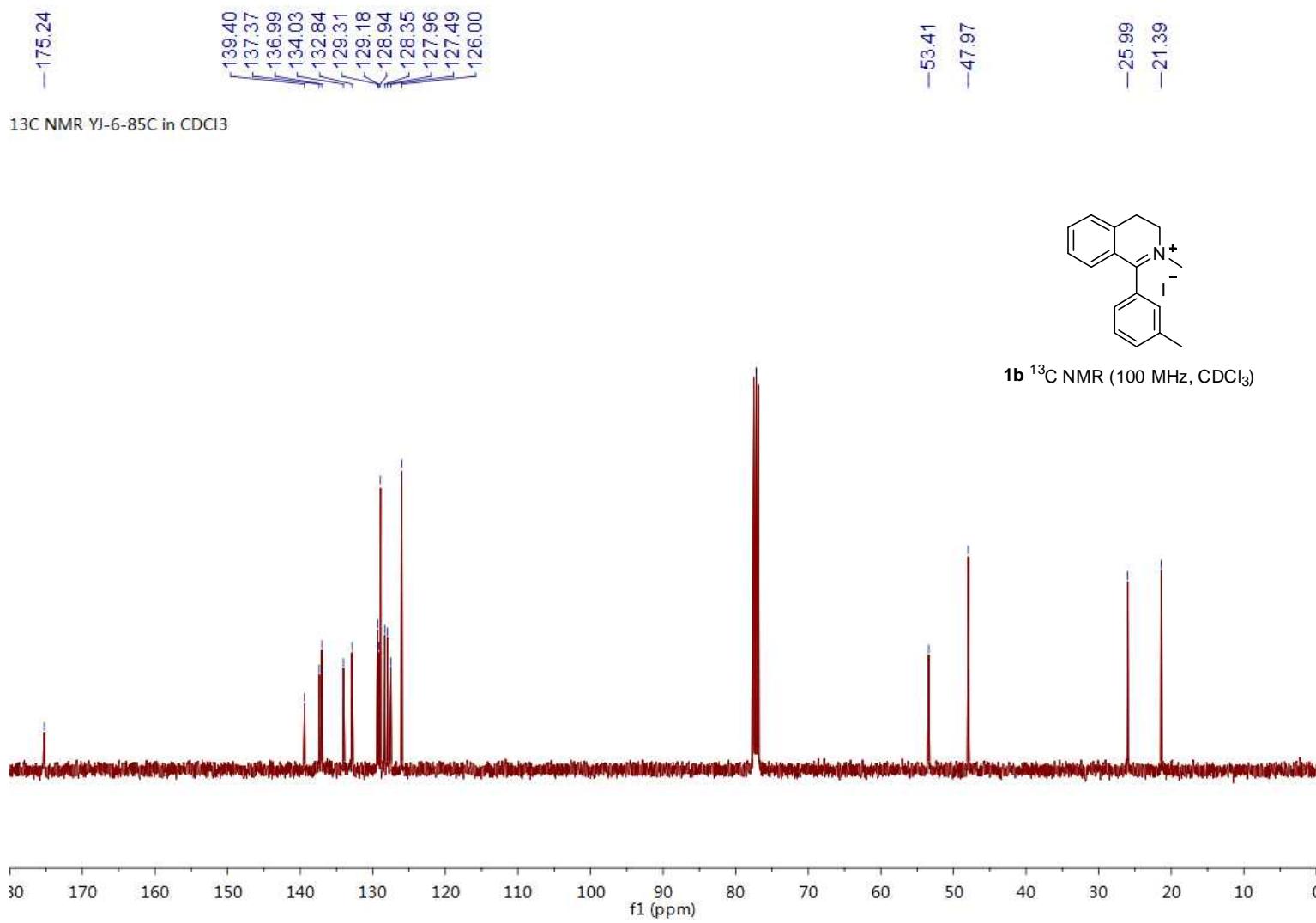
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3.3253

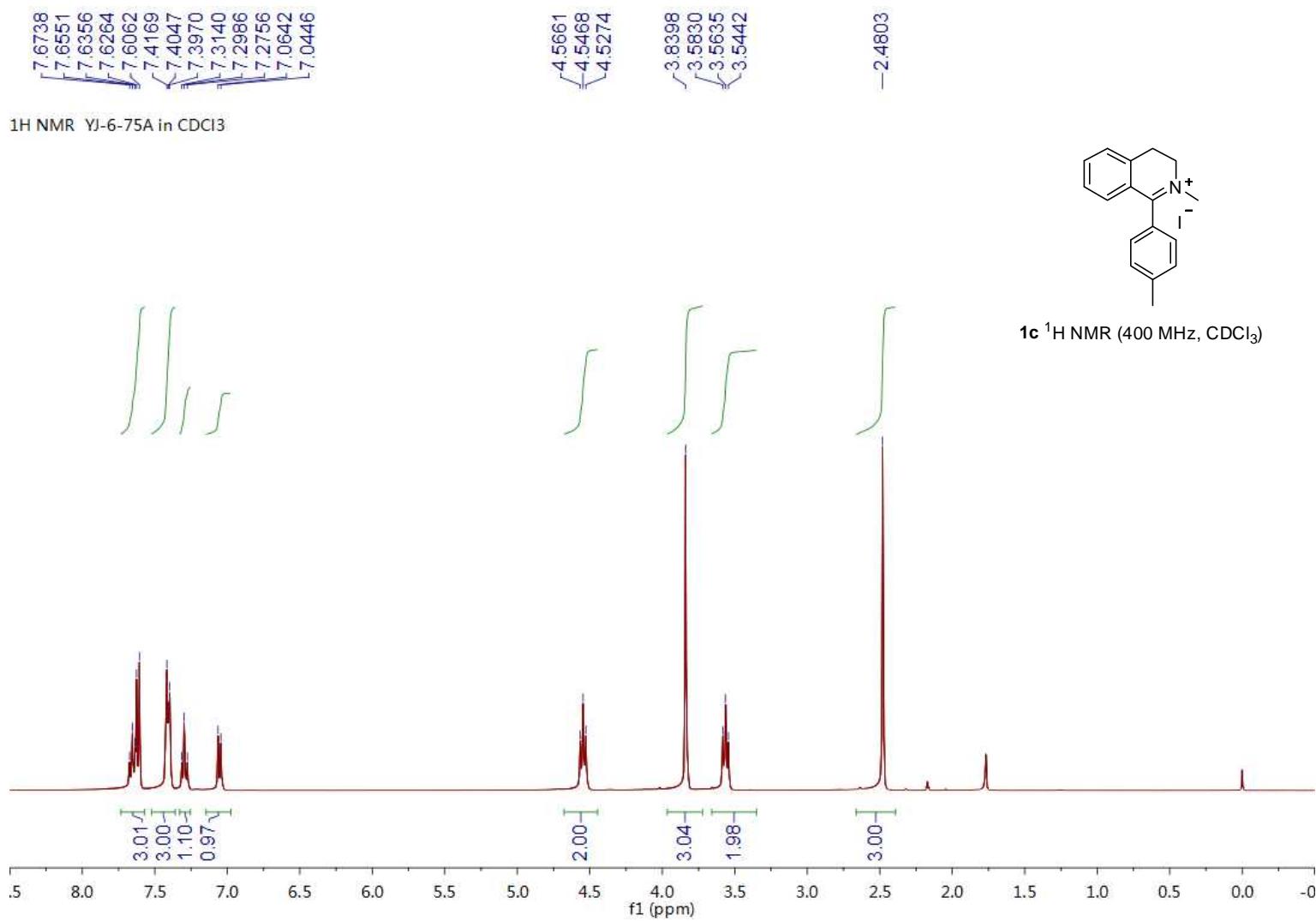
-2.4155



1b ¹H NMR (400 MHz, CDCl₃)







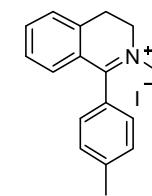
-175.41

143.05
137.54
136.97
134.16
129.87
129.23
128.38
127.98
127.65
126.36

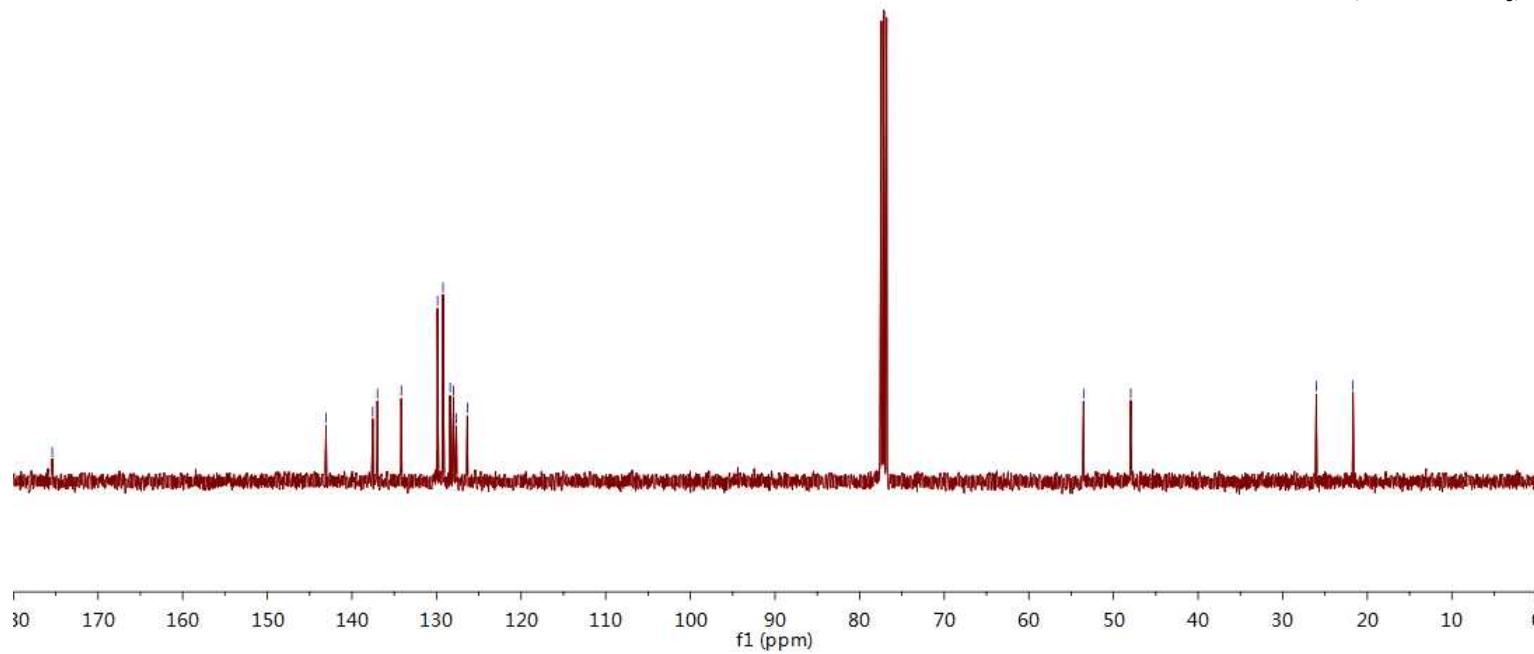
¹³C NMR YJ-6-75A in CDCl₃

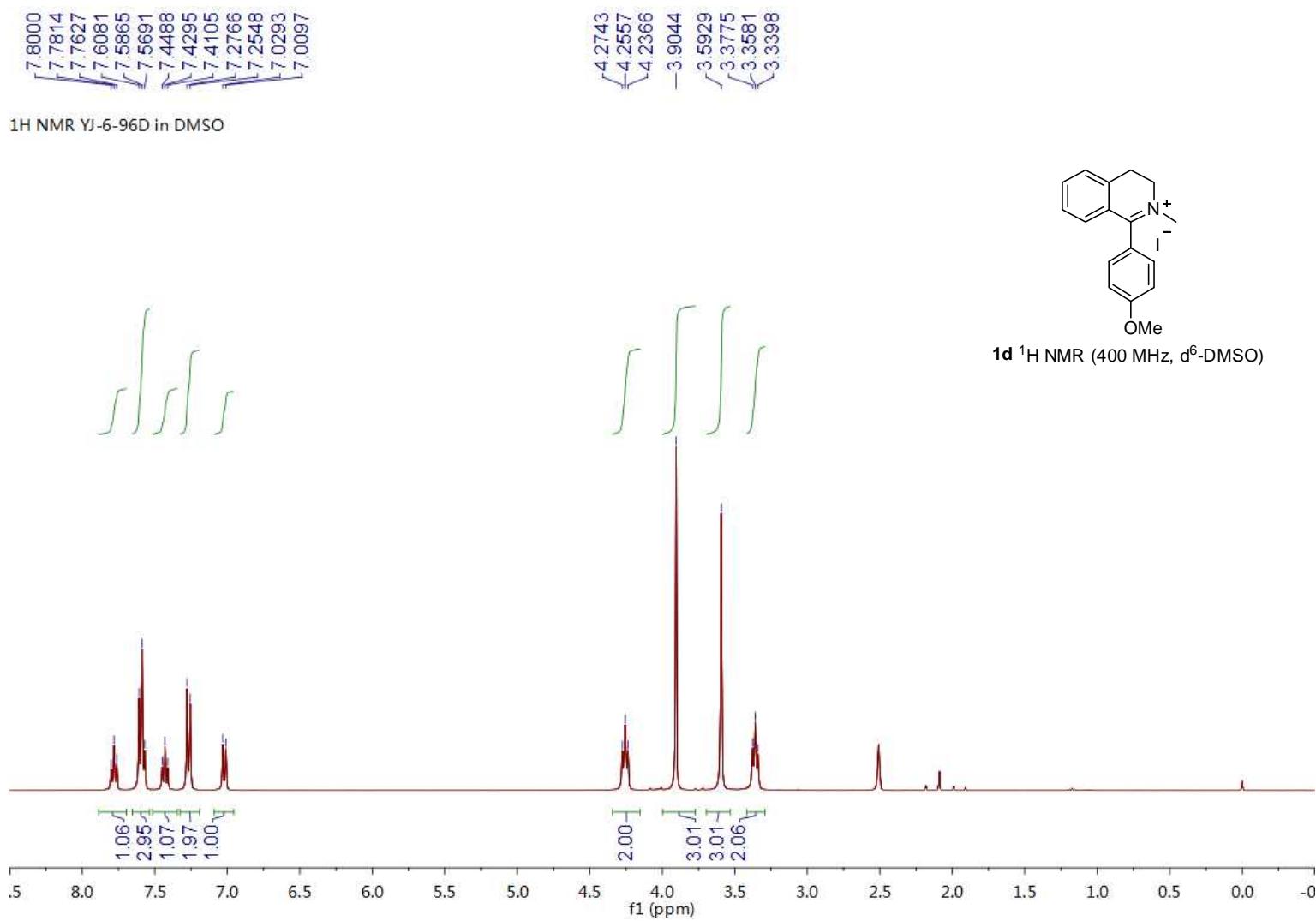
-53.52
-47.96

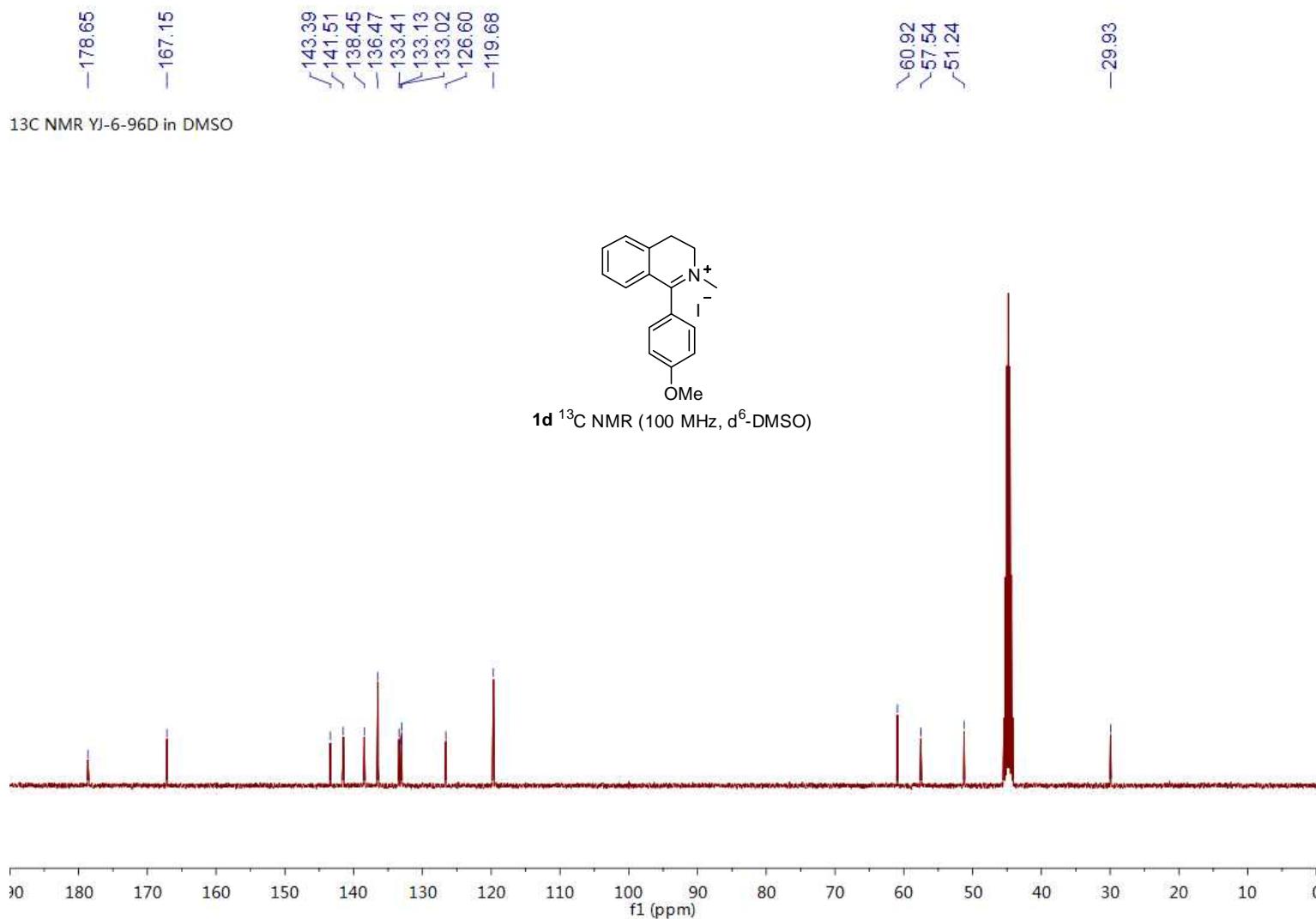
-26.03
-21.69



1c ¹³C NMR (400 MHz, CDCl₃)

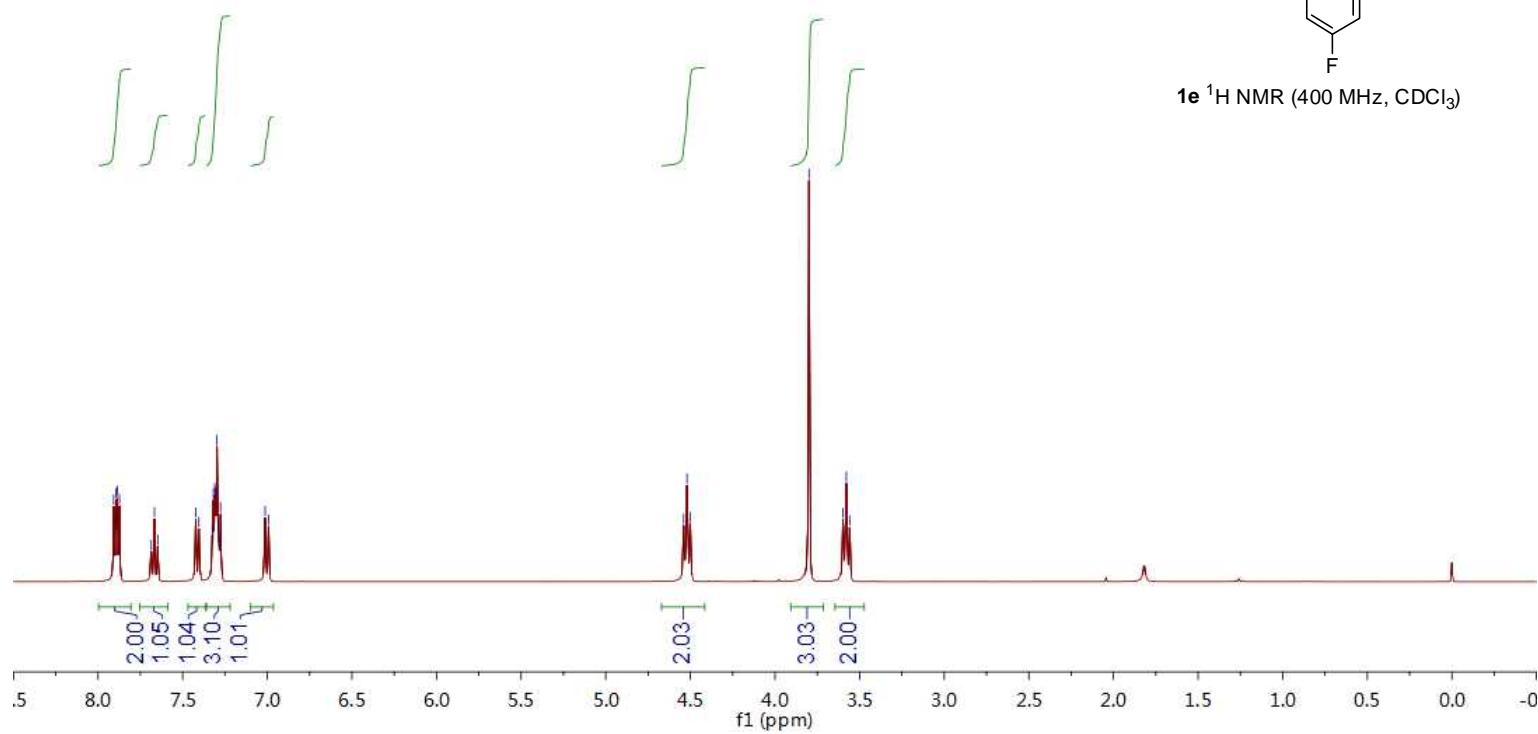






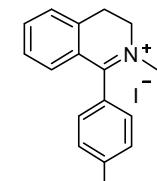
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7.3184
7.3103
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7.2758
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6.9930

1H NMR YJ-6-89E in CDCl₃

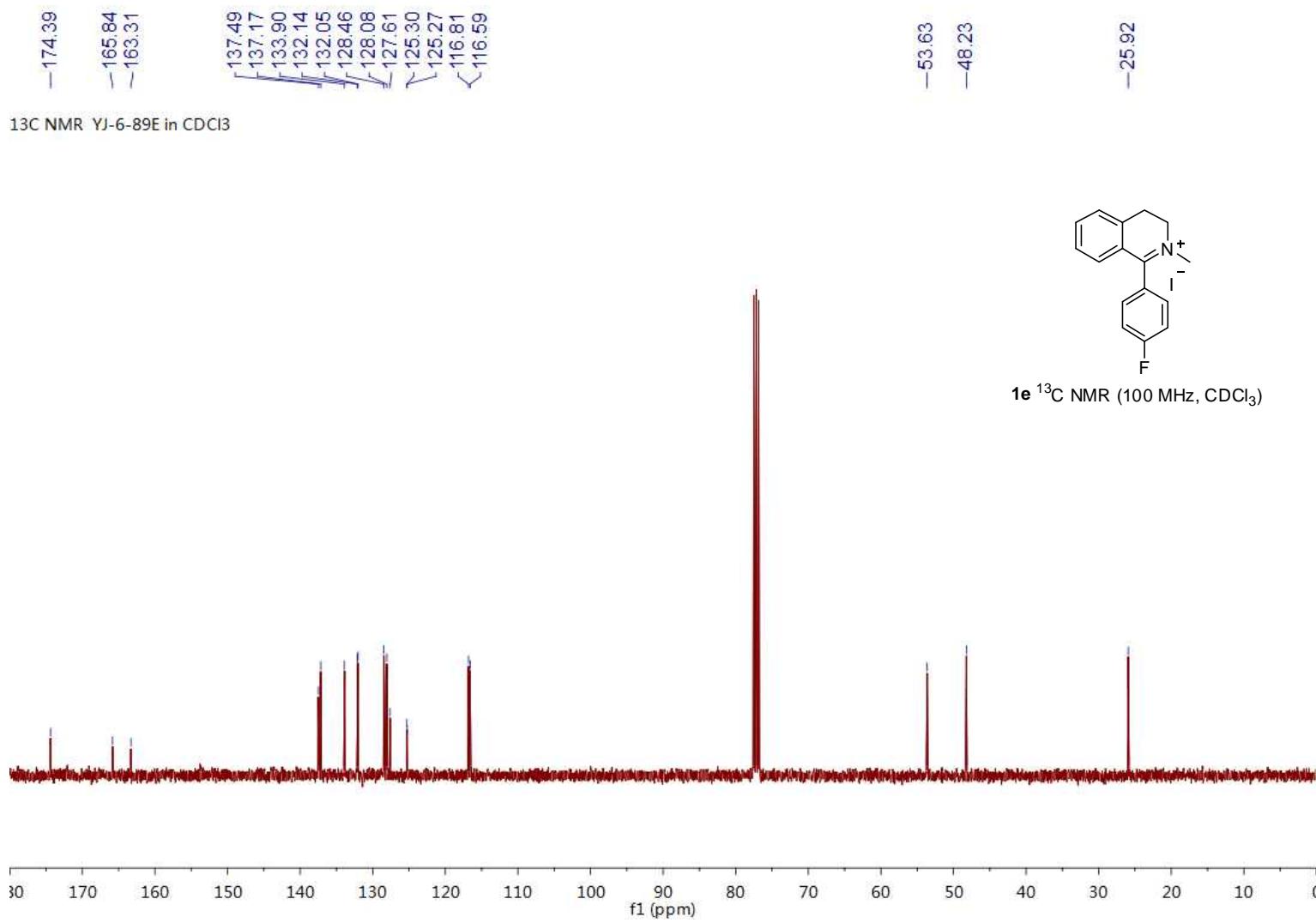


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4.5202
4.5006

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3.5588

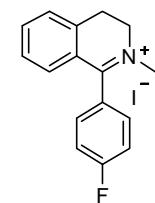


1e ¹H NMR (400 MHz, CDCl₃)

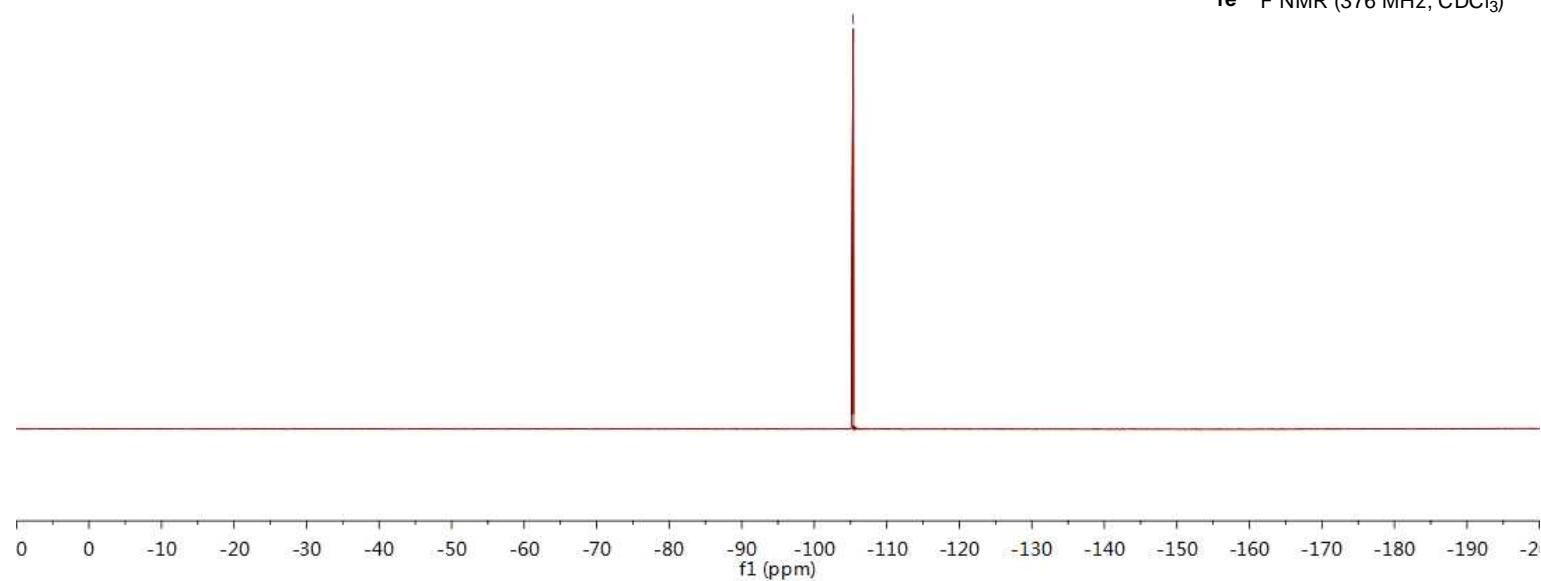


¹⁹F NMR YJ-6-89E in CDCl₃

--105.37



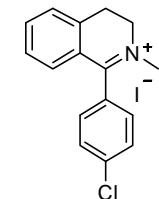
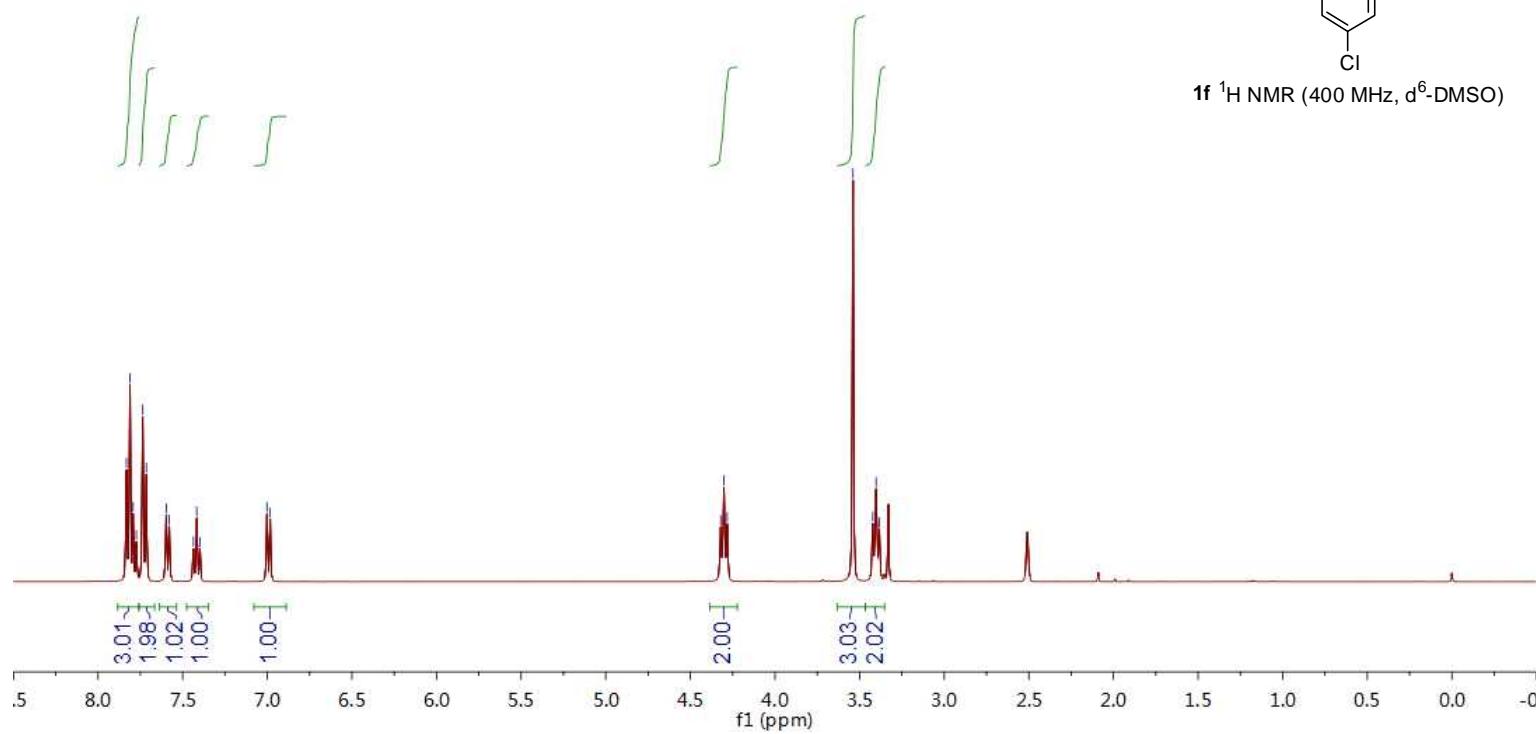
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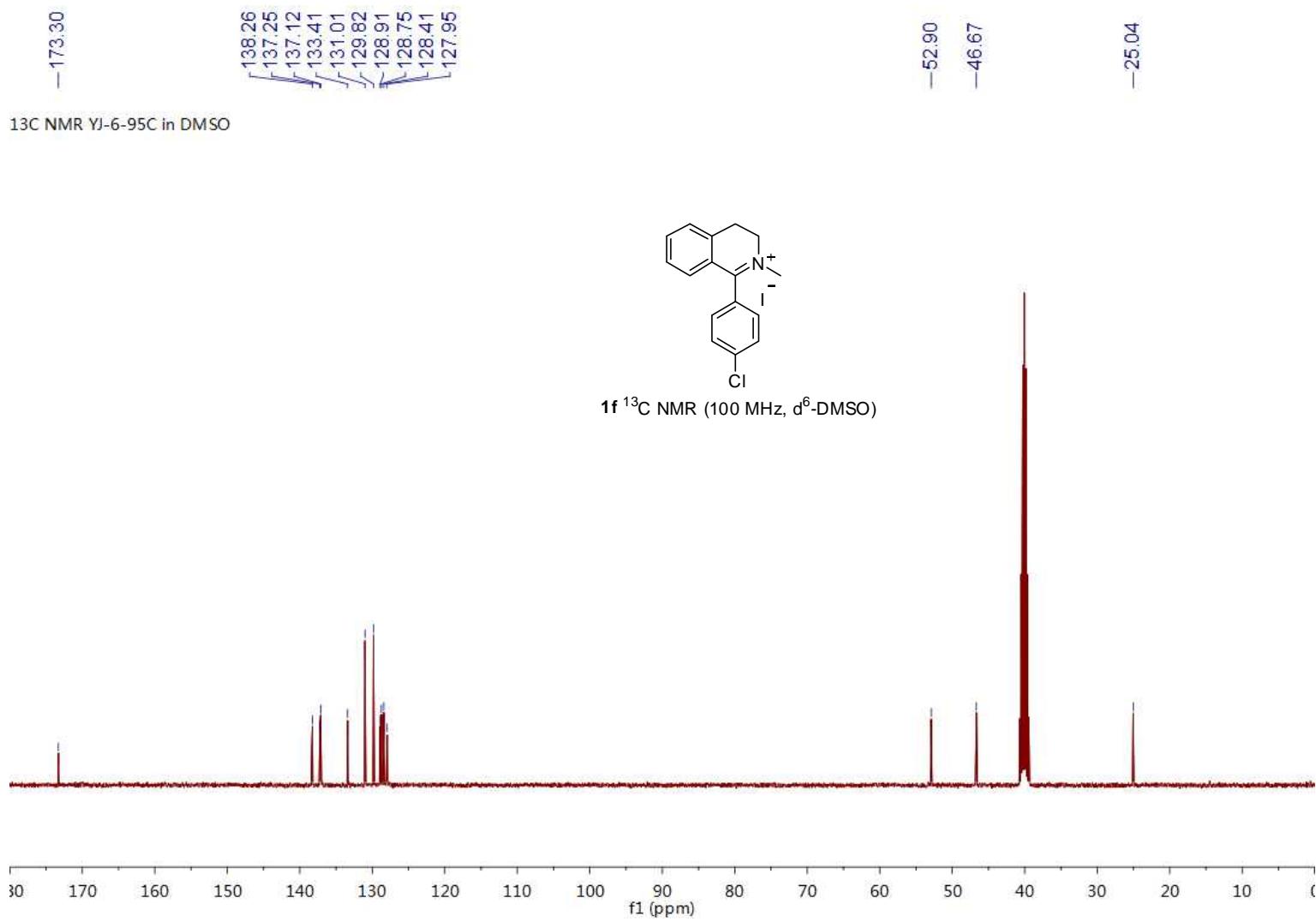
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¹H NMR YJ-6-95C in DMSO

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4.3017
4.2821
3.5387
3.4219
3.4025
3.3833



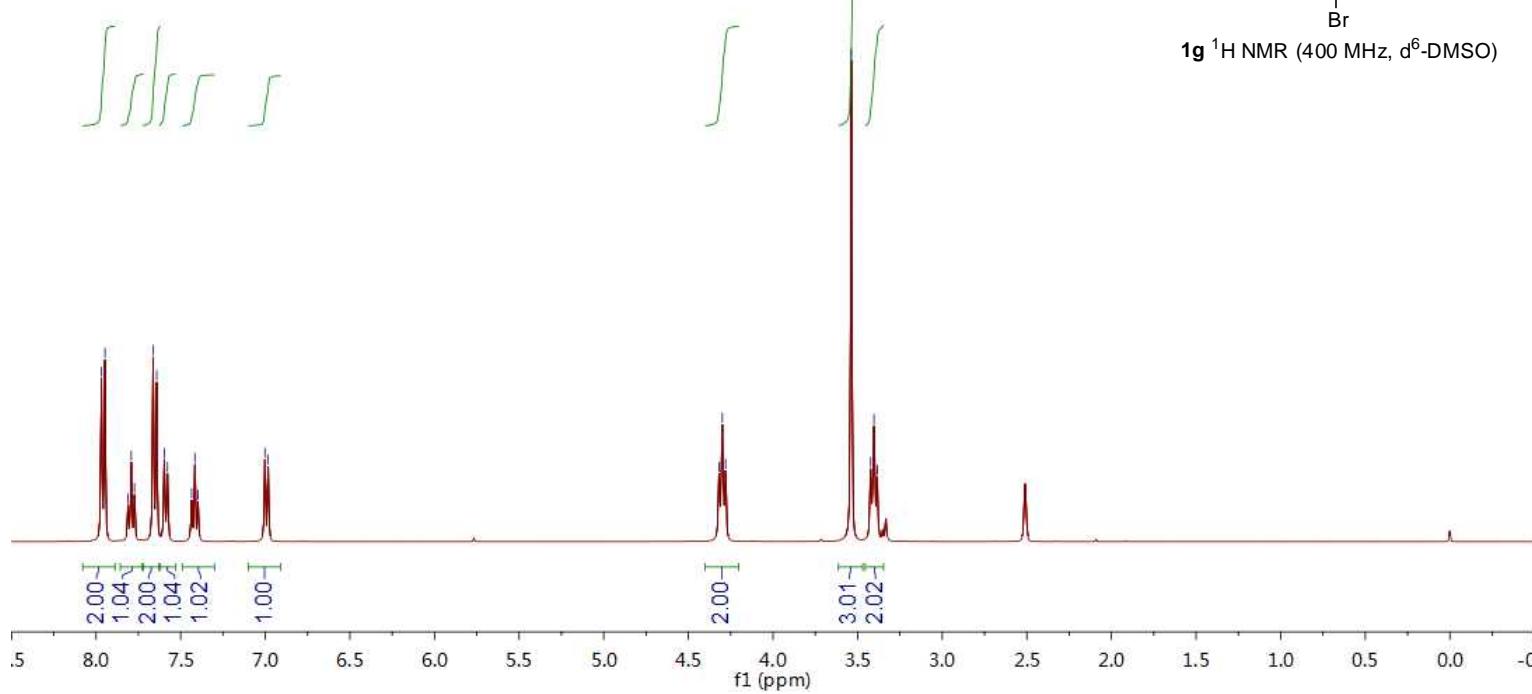
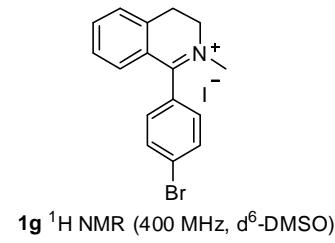
1f ¹H NMR (400 MHz, d⁶-DMSO)

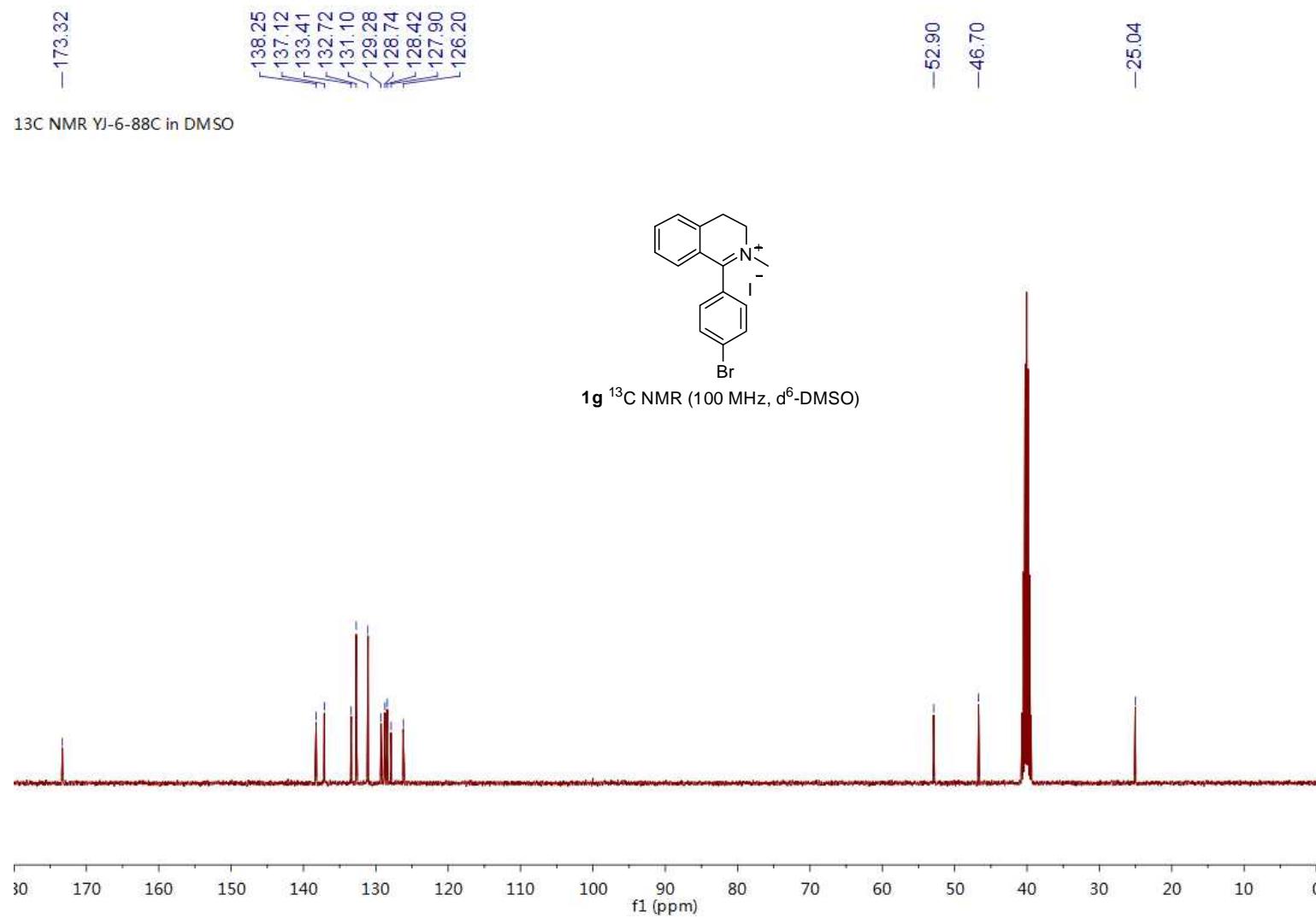


7.9681
7.9472
7.7919
7.6626
7.6416
7.5964
7.5777
7.4169
6.9827

1H NMR YJ-6-88C in DMSO

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3.5372
3.4226
3.4033
3.3841



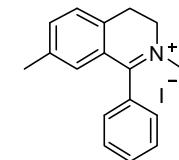


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6.4378

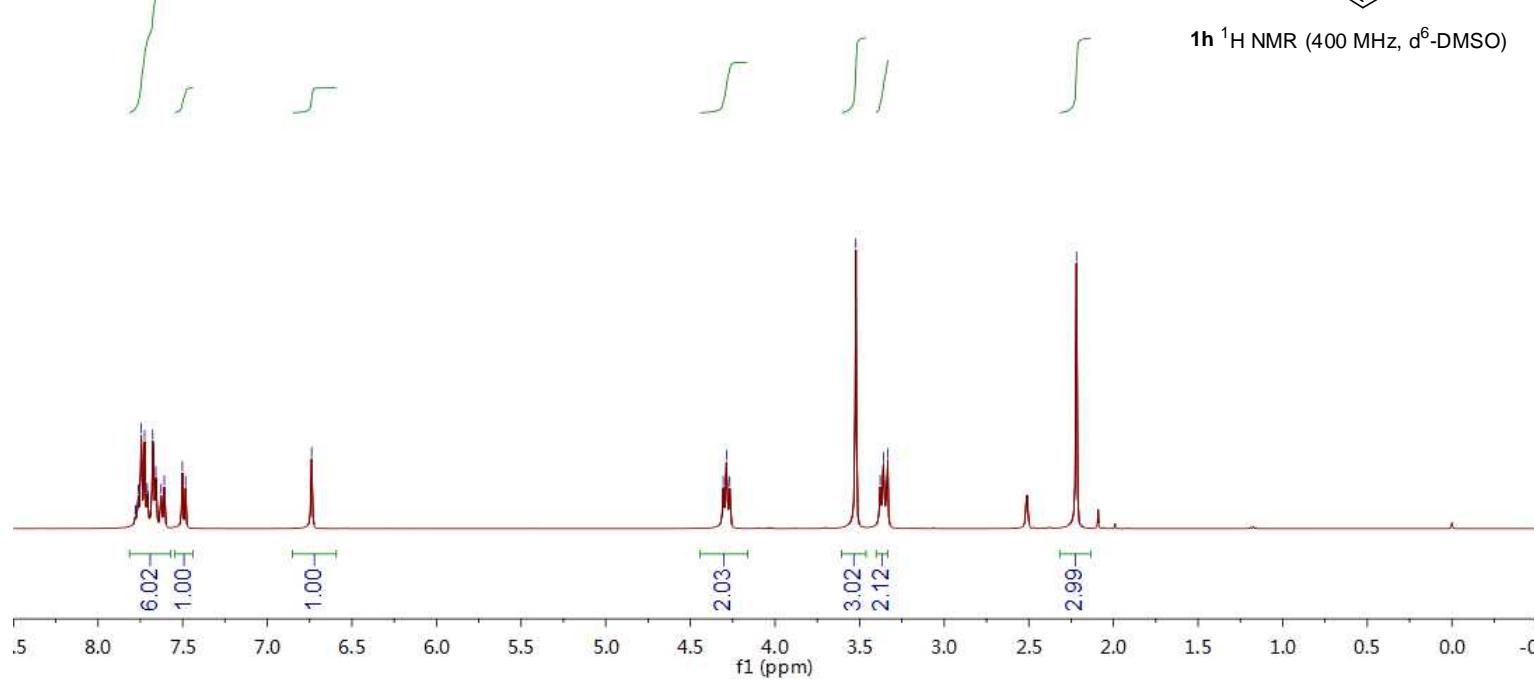
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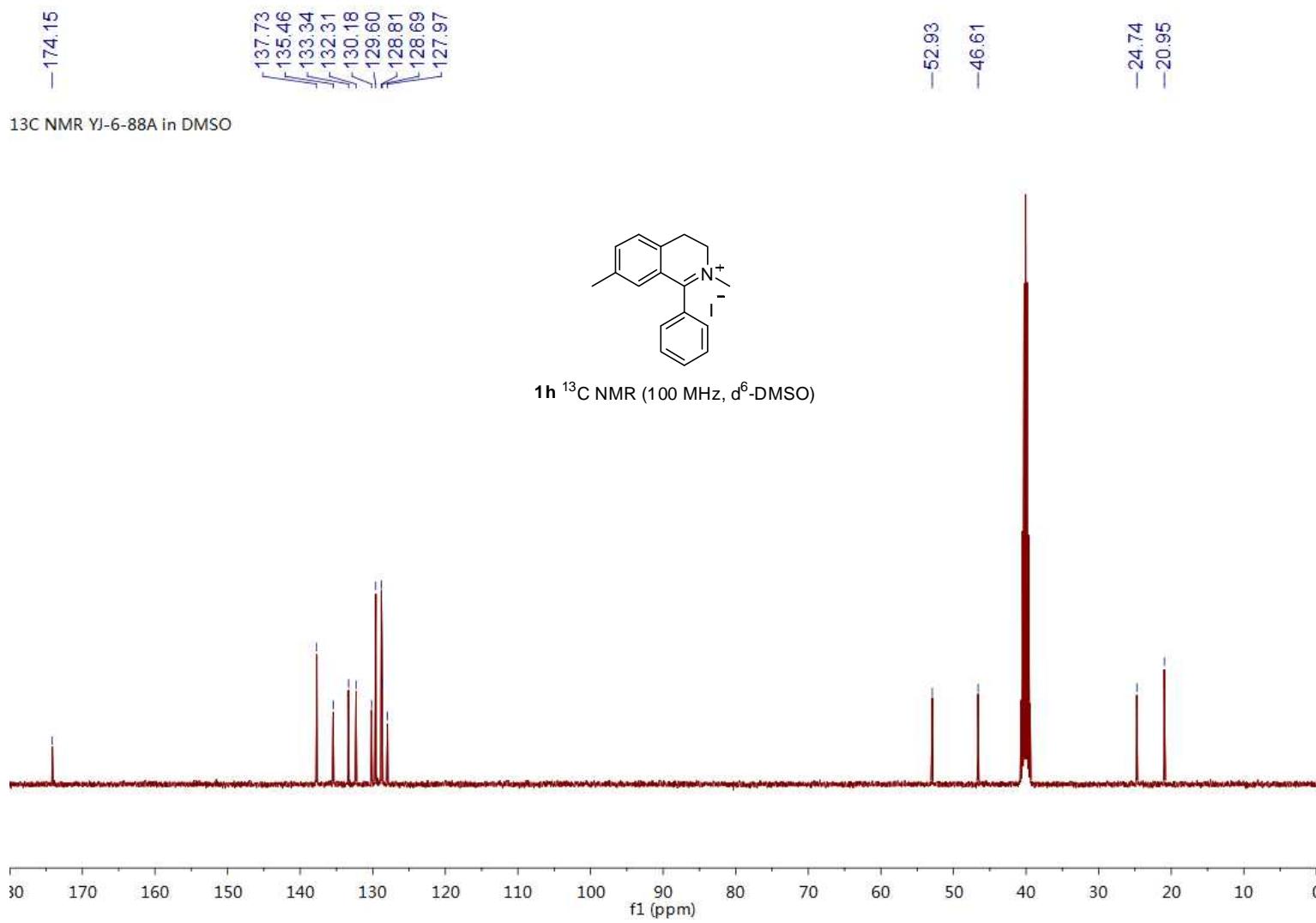
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-2.2181



1h ^1H NMR (400 MHz, $\text{d}^6\text{-DMSO}$)

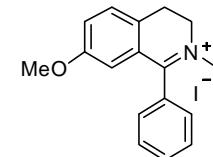




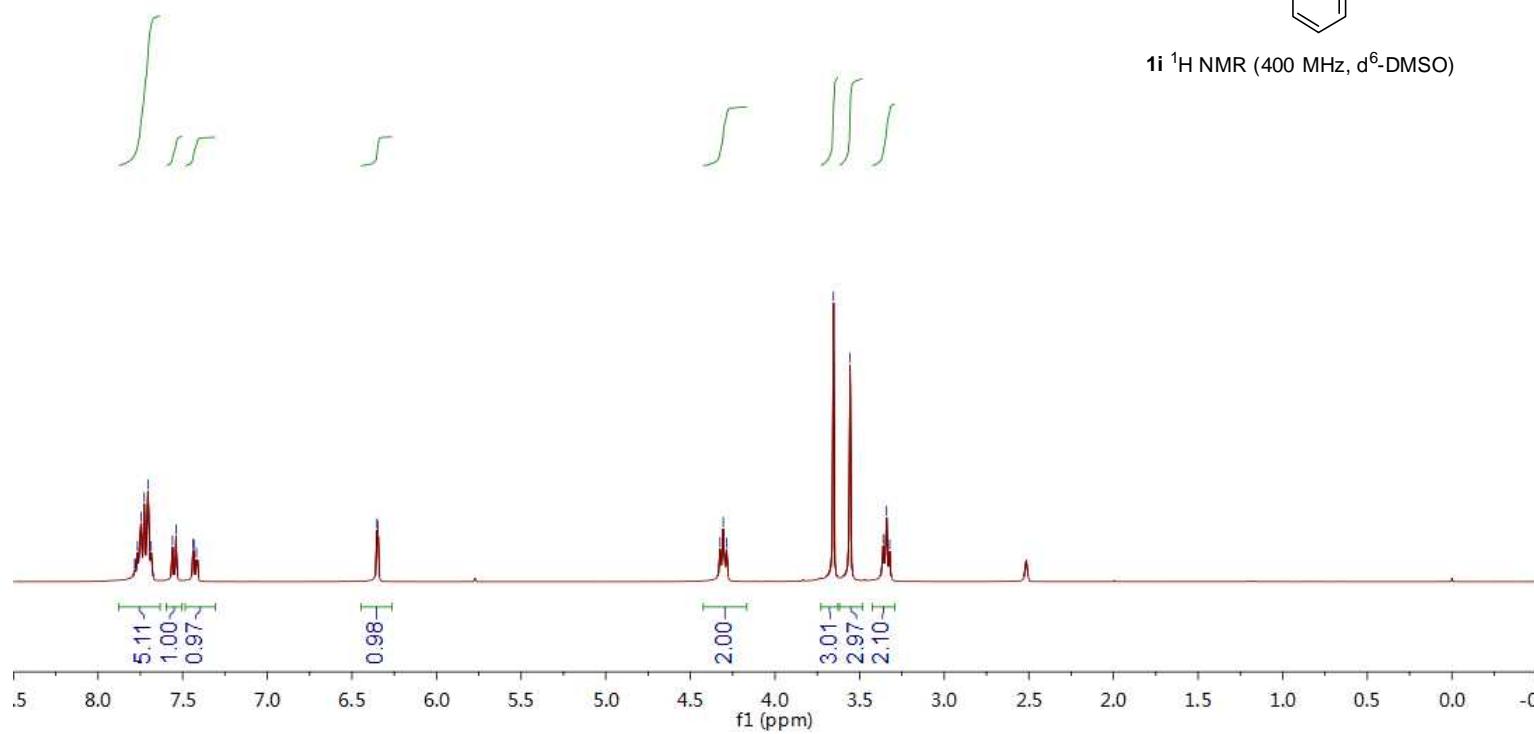
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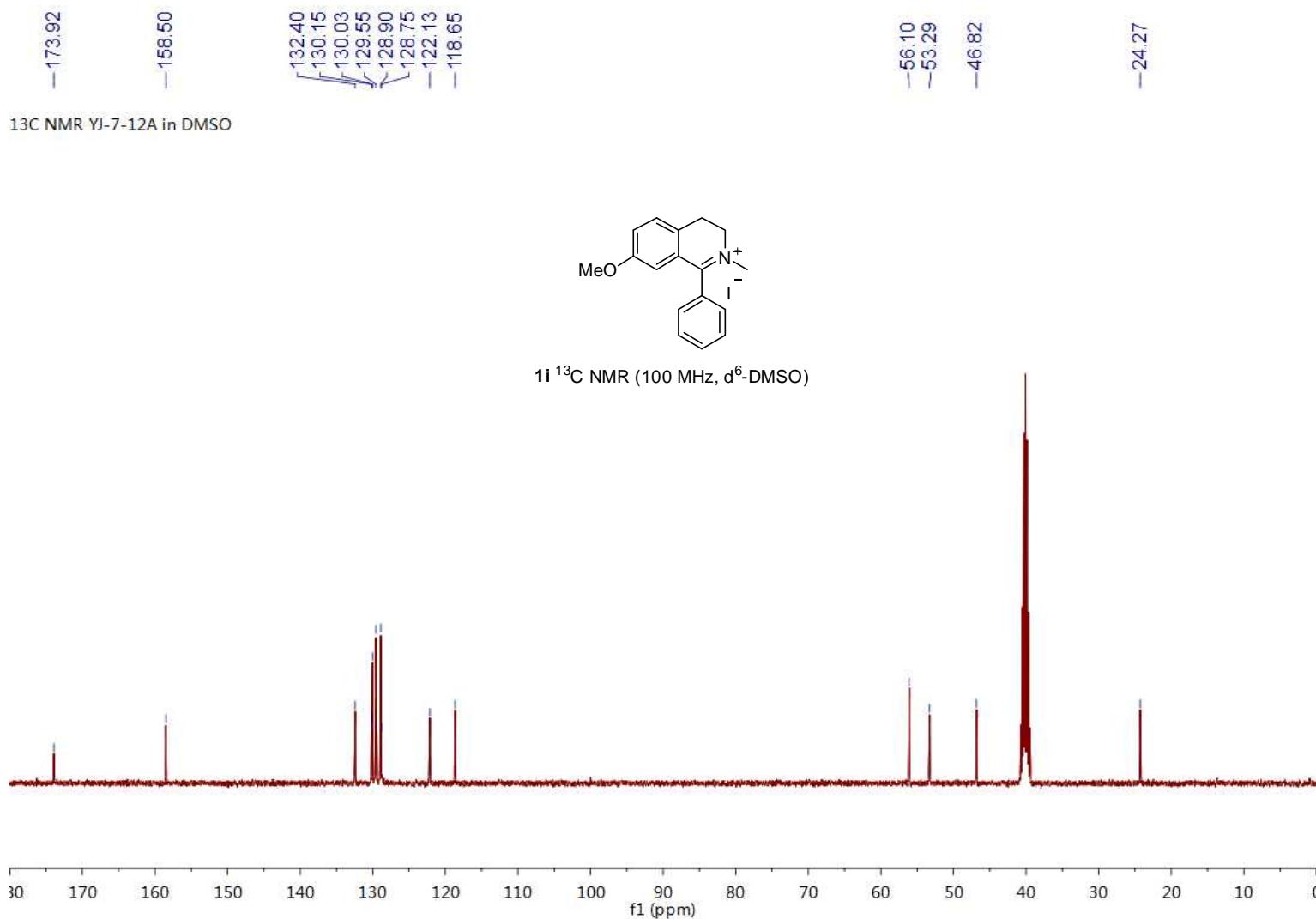
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 3.3403
 3.3220



1i ¹H NMR (400 MHz, d⁶-DMSO)

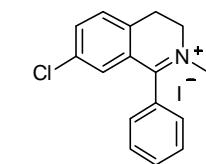




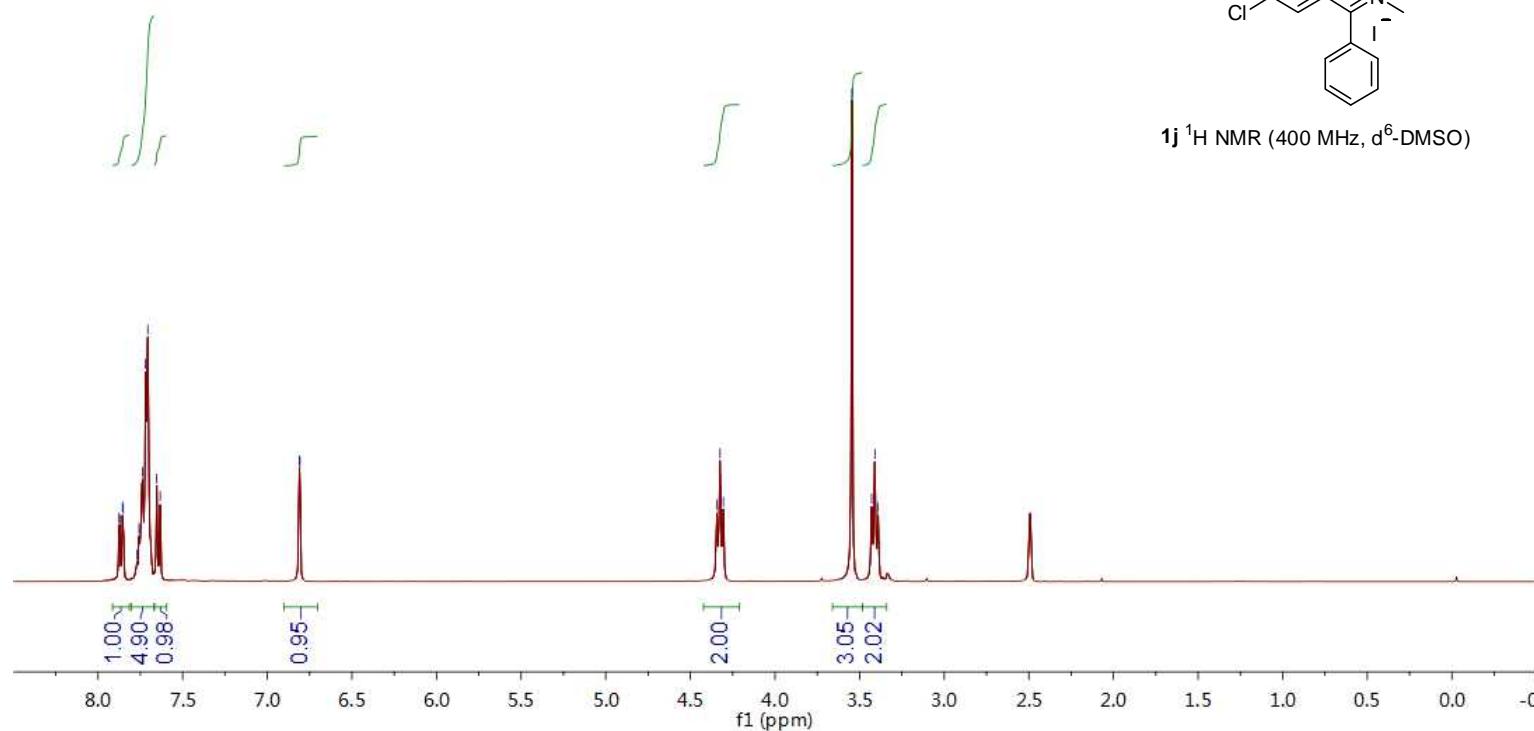
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6.8071

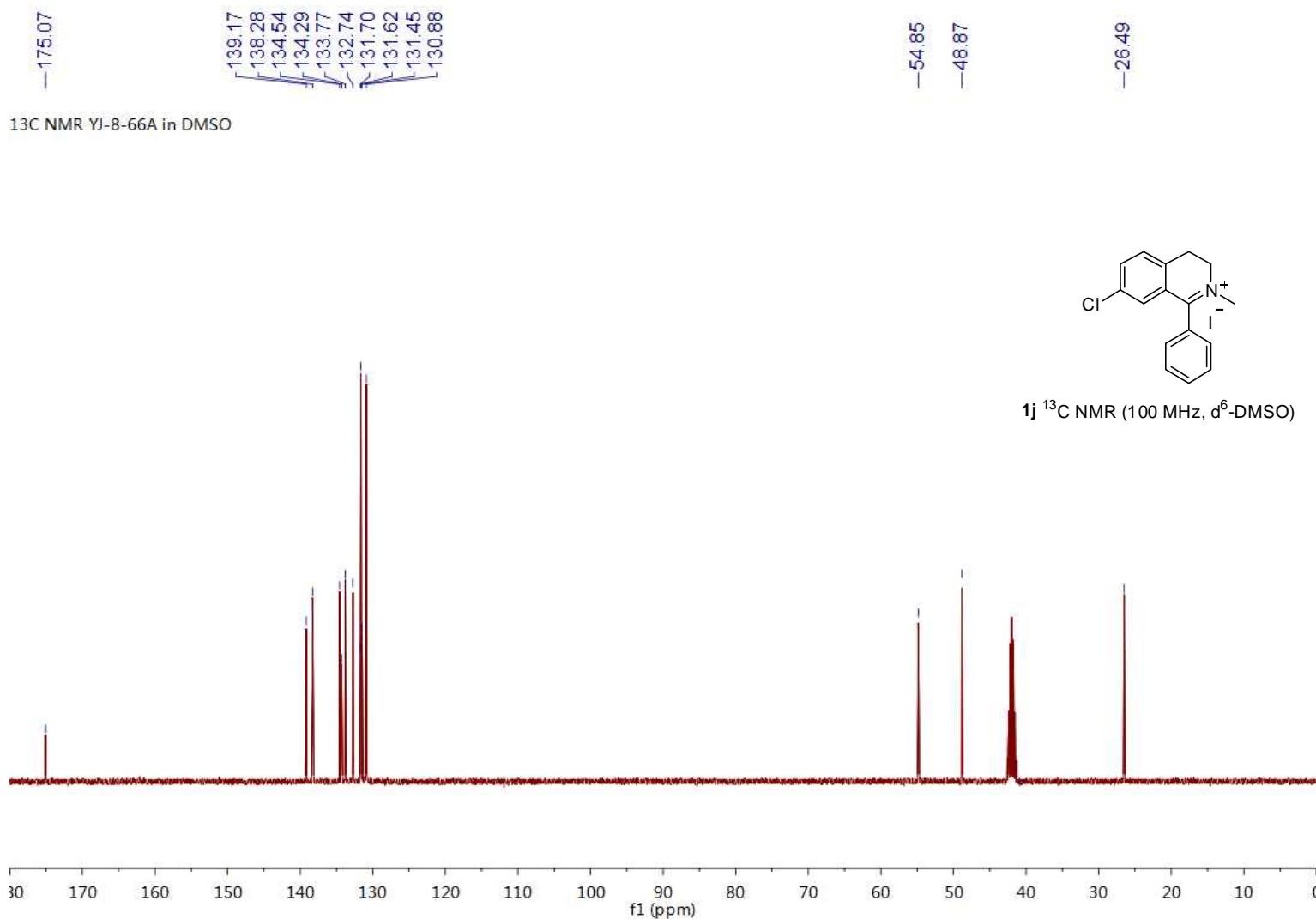
^1H NMR YJ-8-66A in DMSO

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3.3904



1j ^1H NMR (400 MHz, $\text{d}^6\text{-DMSO}$)

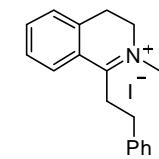




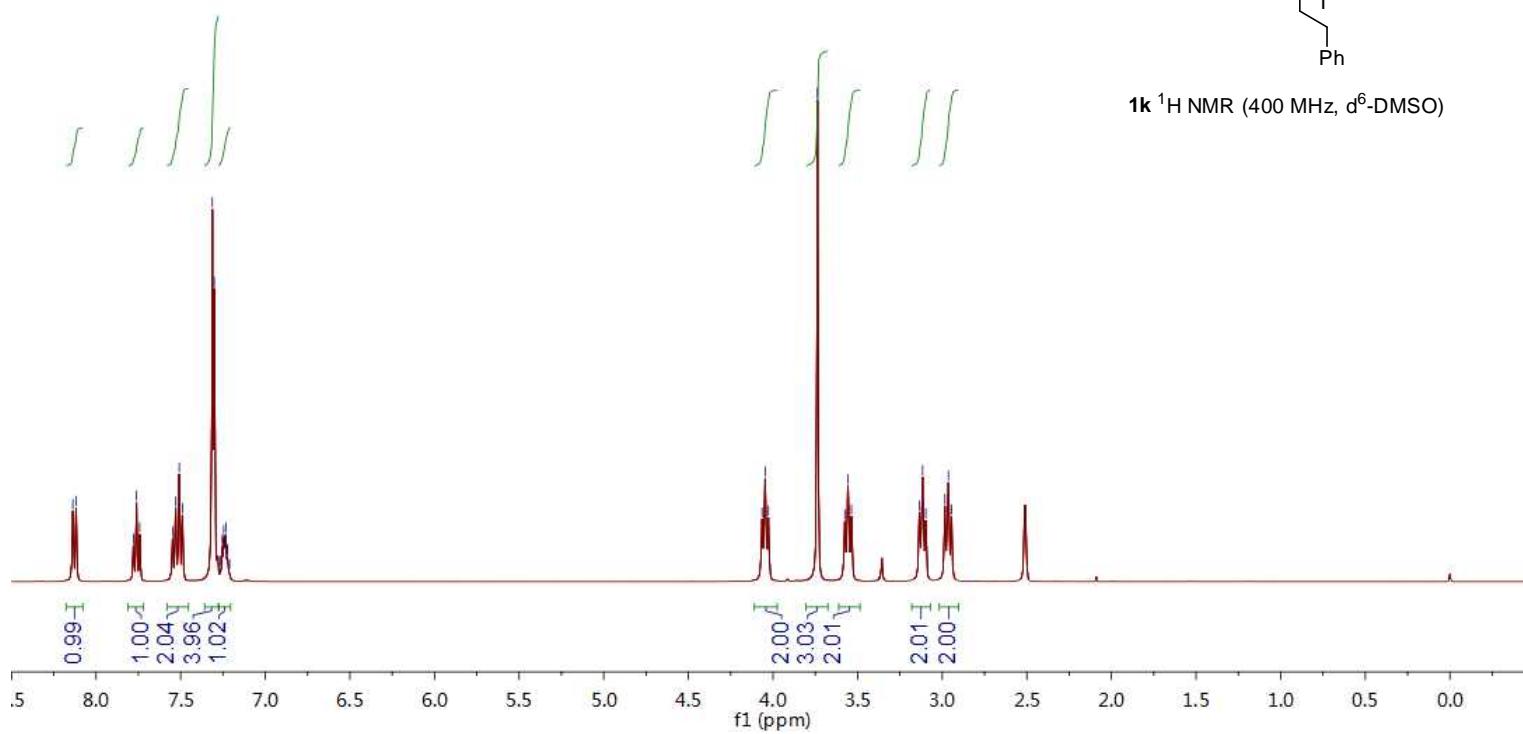
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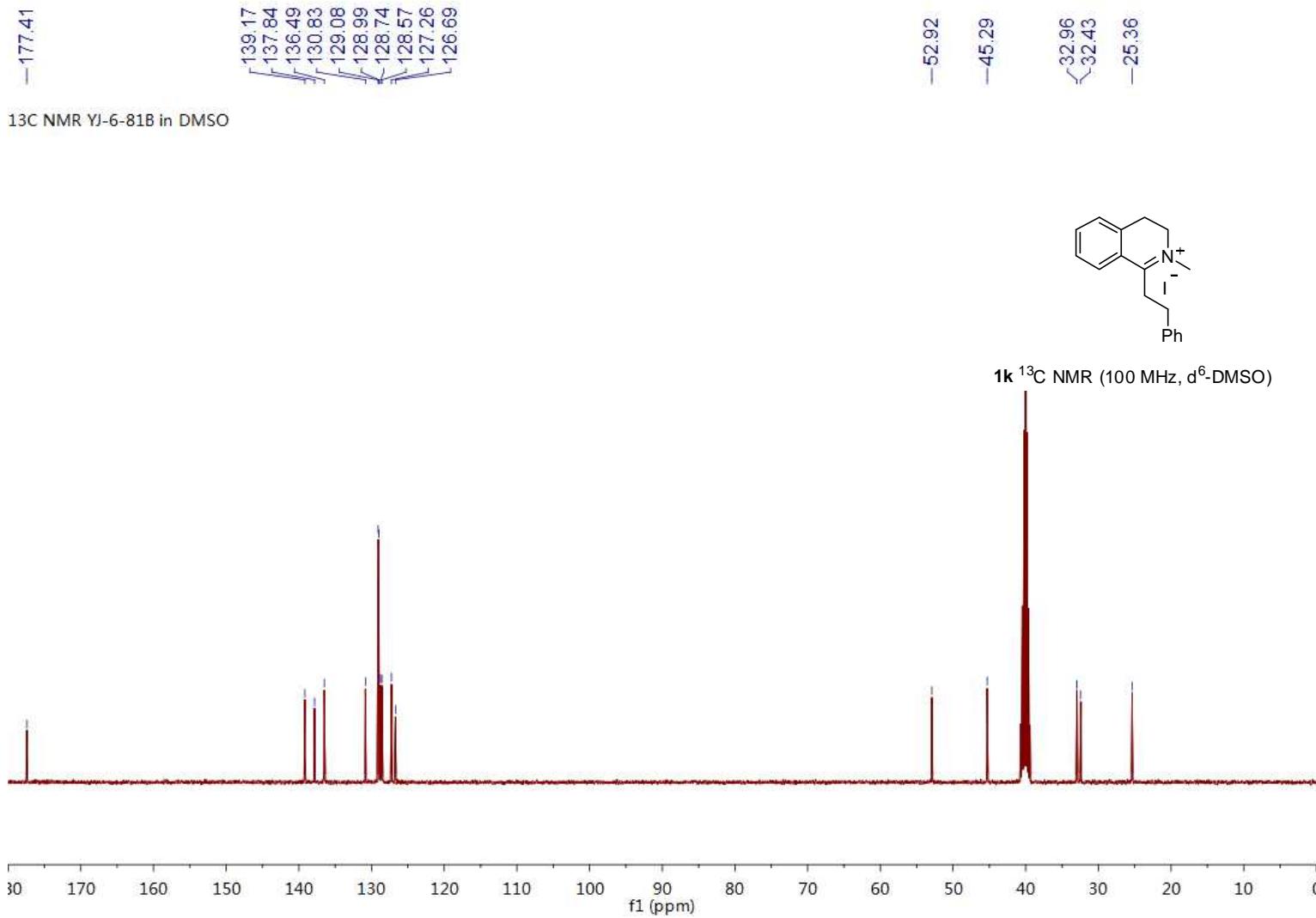
¹H NMR YJ-6-81B in DMSO

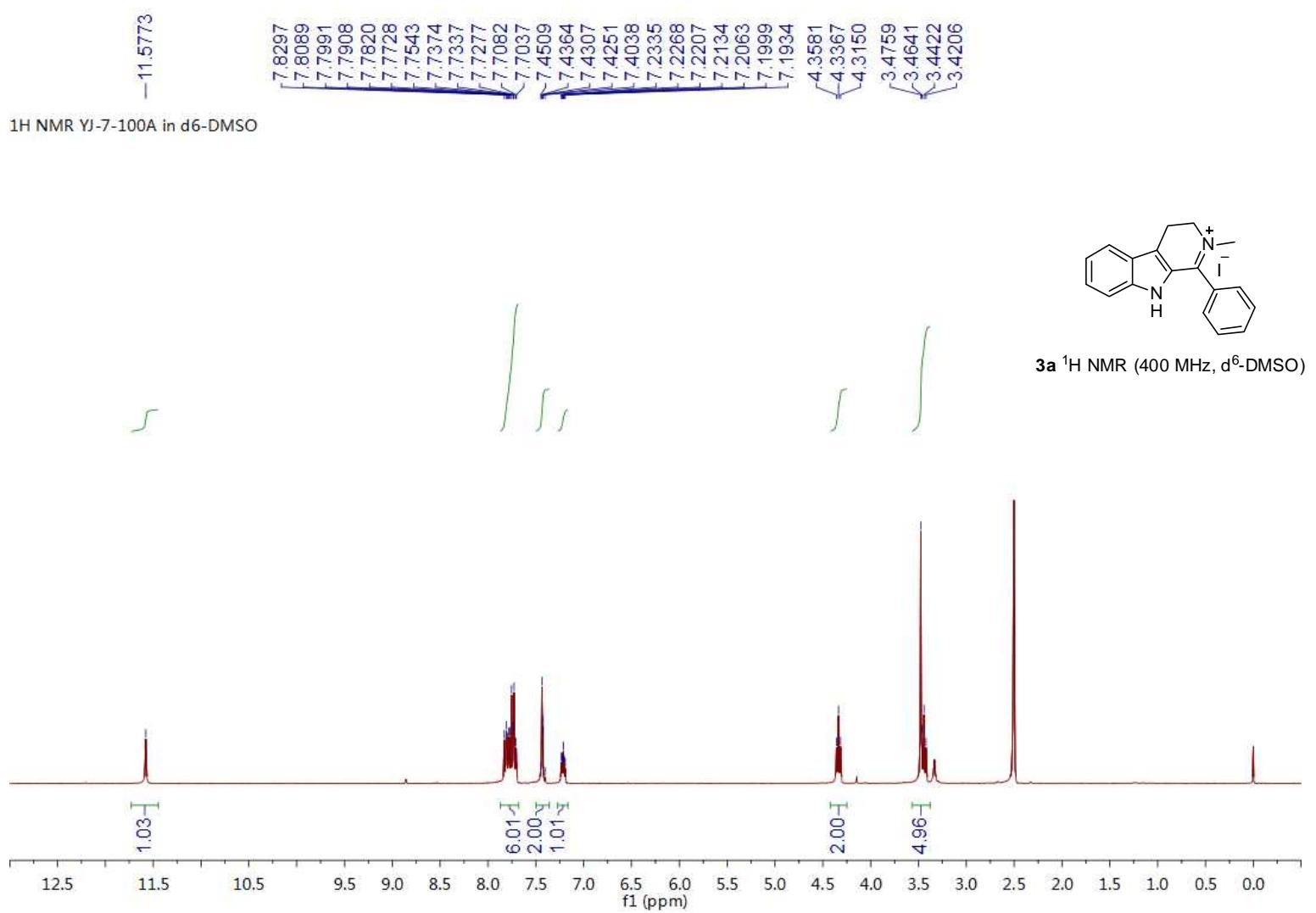
4.0646
 4.0461
 4.0271
 3.7368
 3.5756
 3.5567
 3.5362
 3.1346
 3.1157
 3.0971
 2.9839
 2.9635
 2.9445

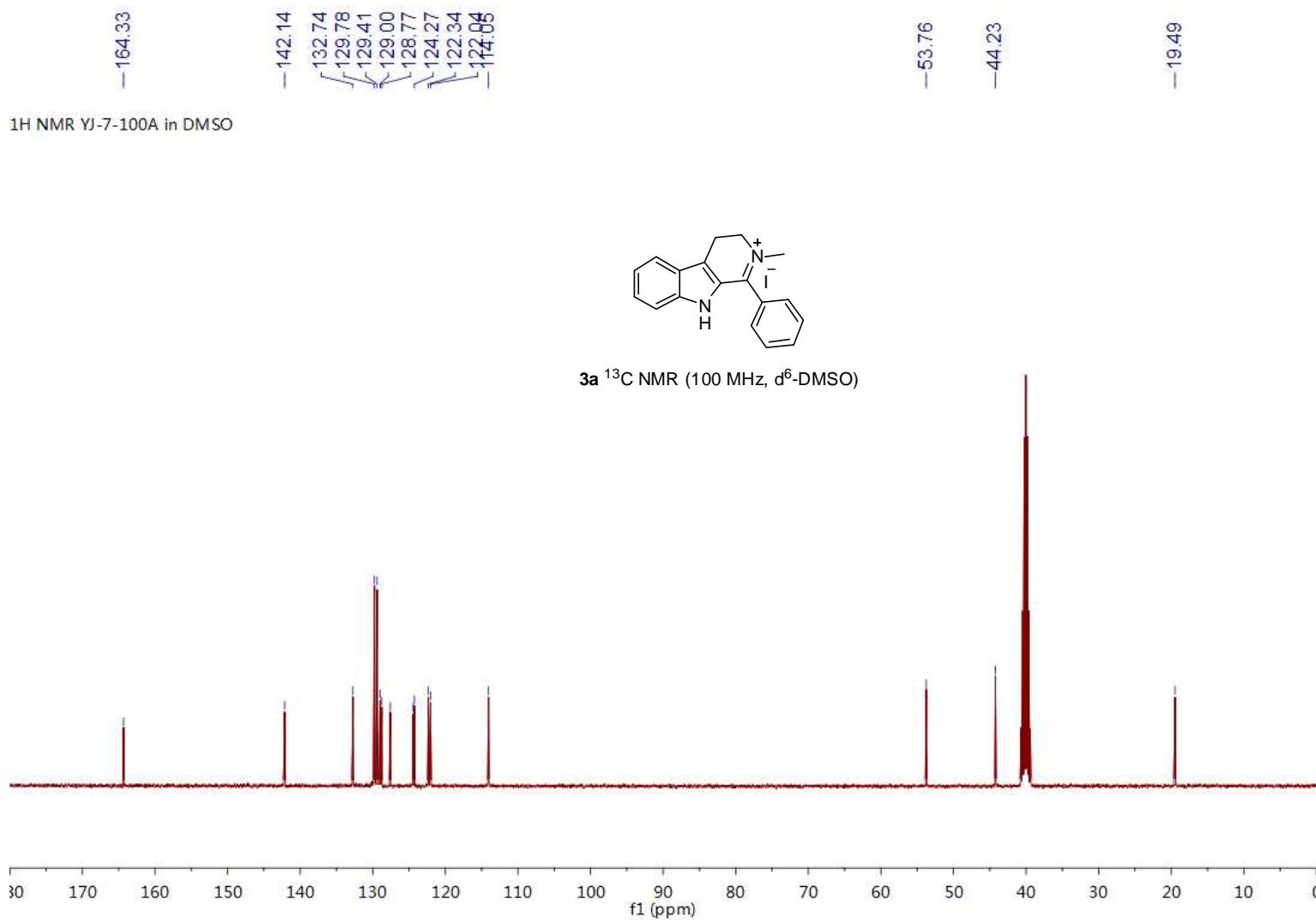


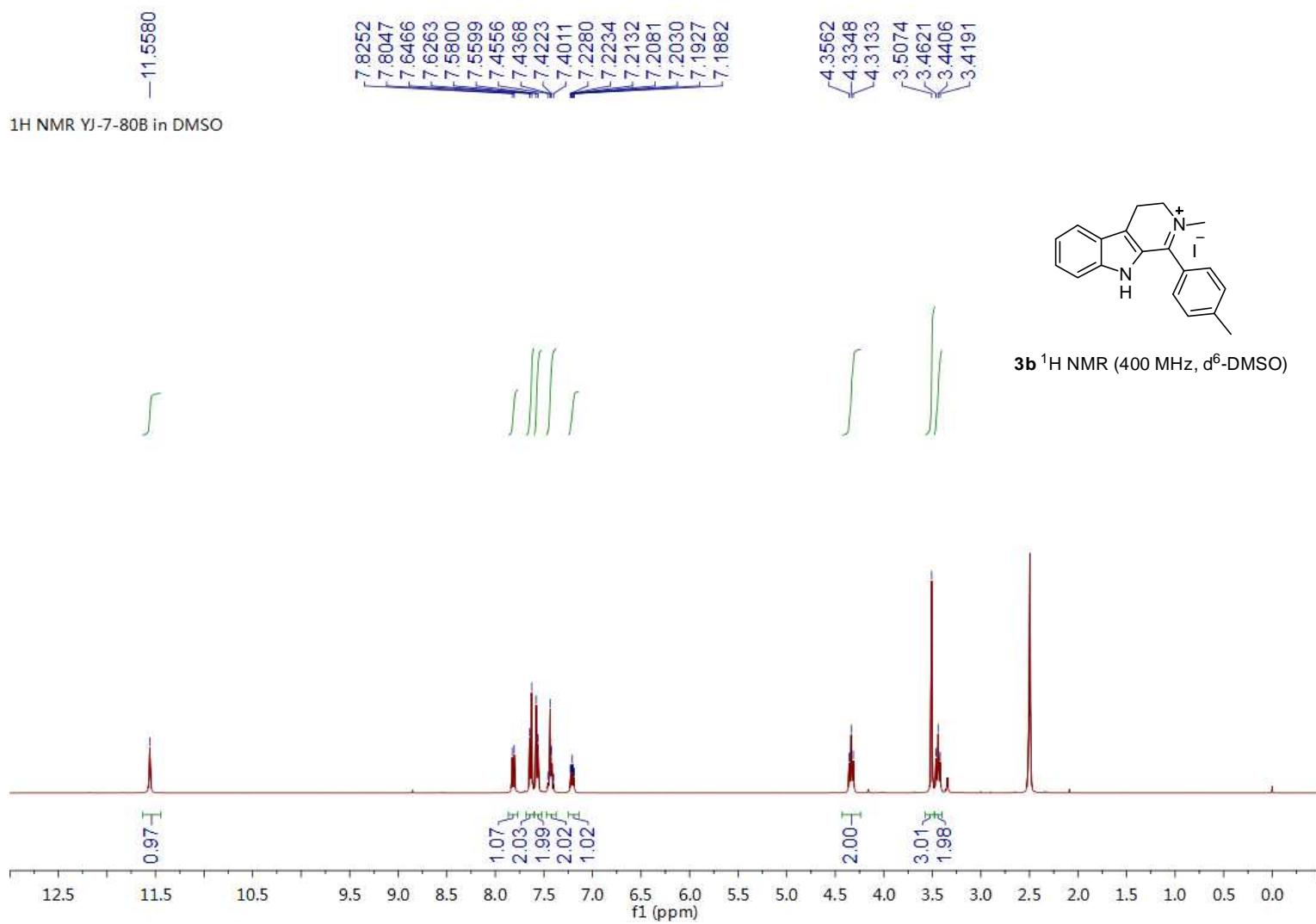
1k ¹H NMR (400 MHz, d⁶-DMSO)

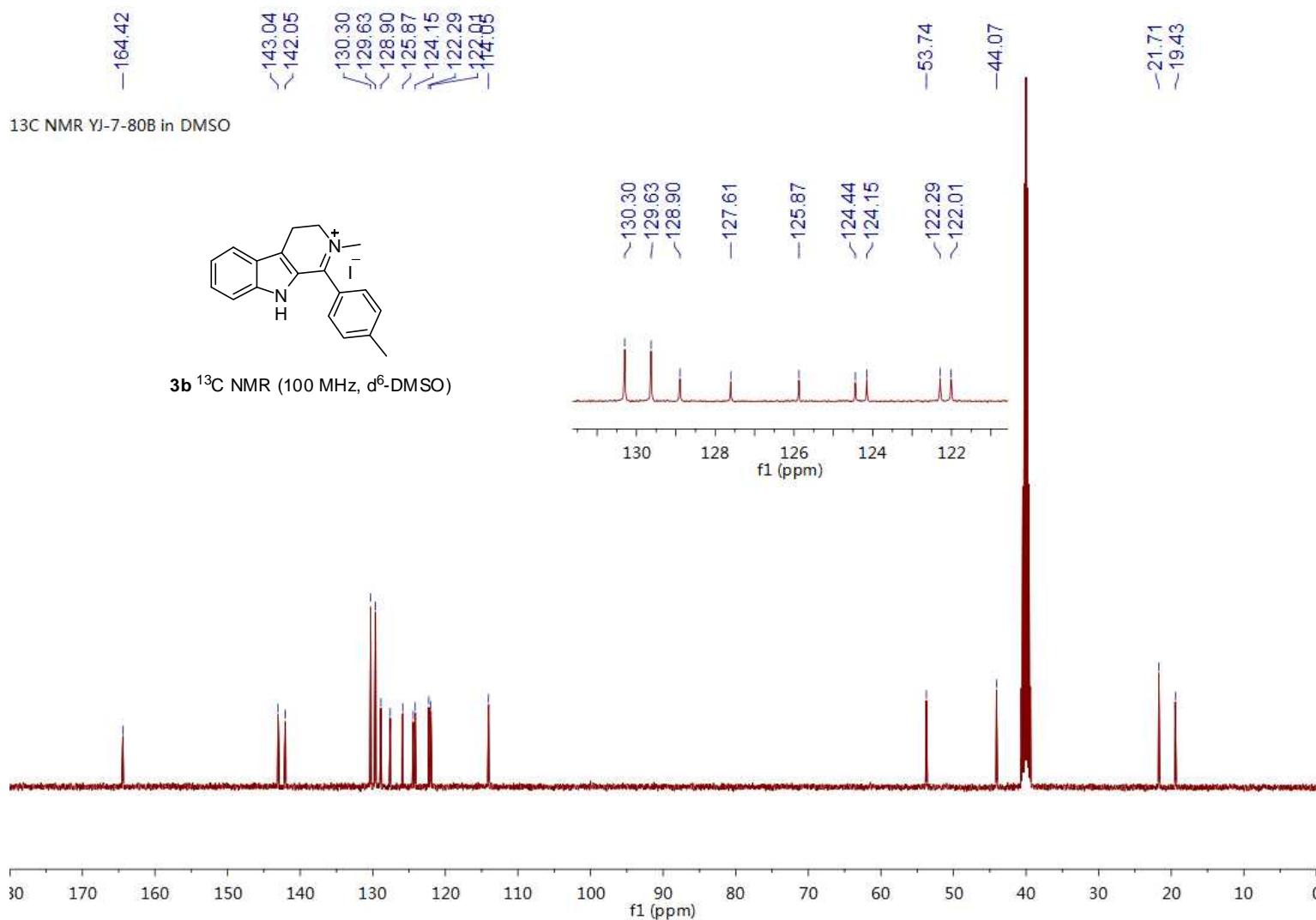


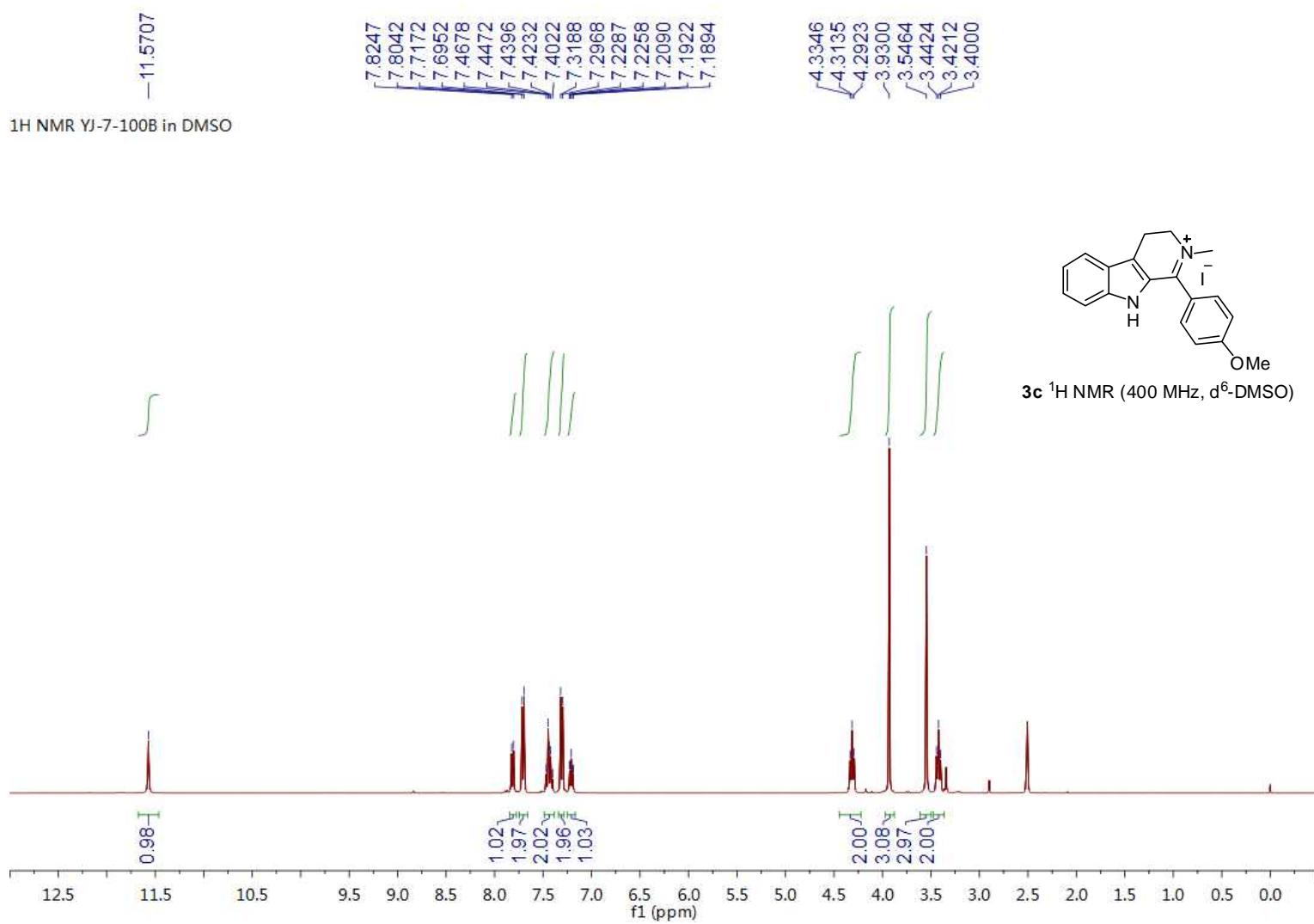


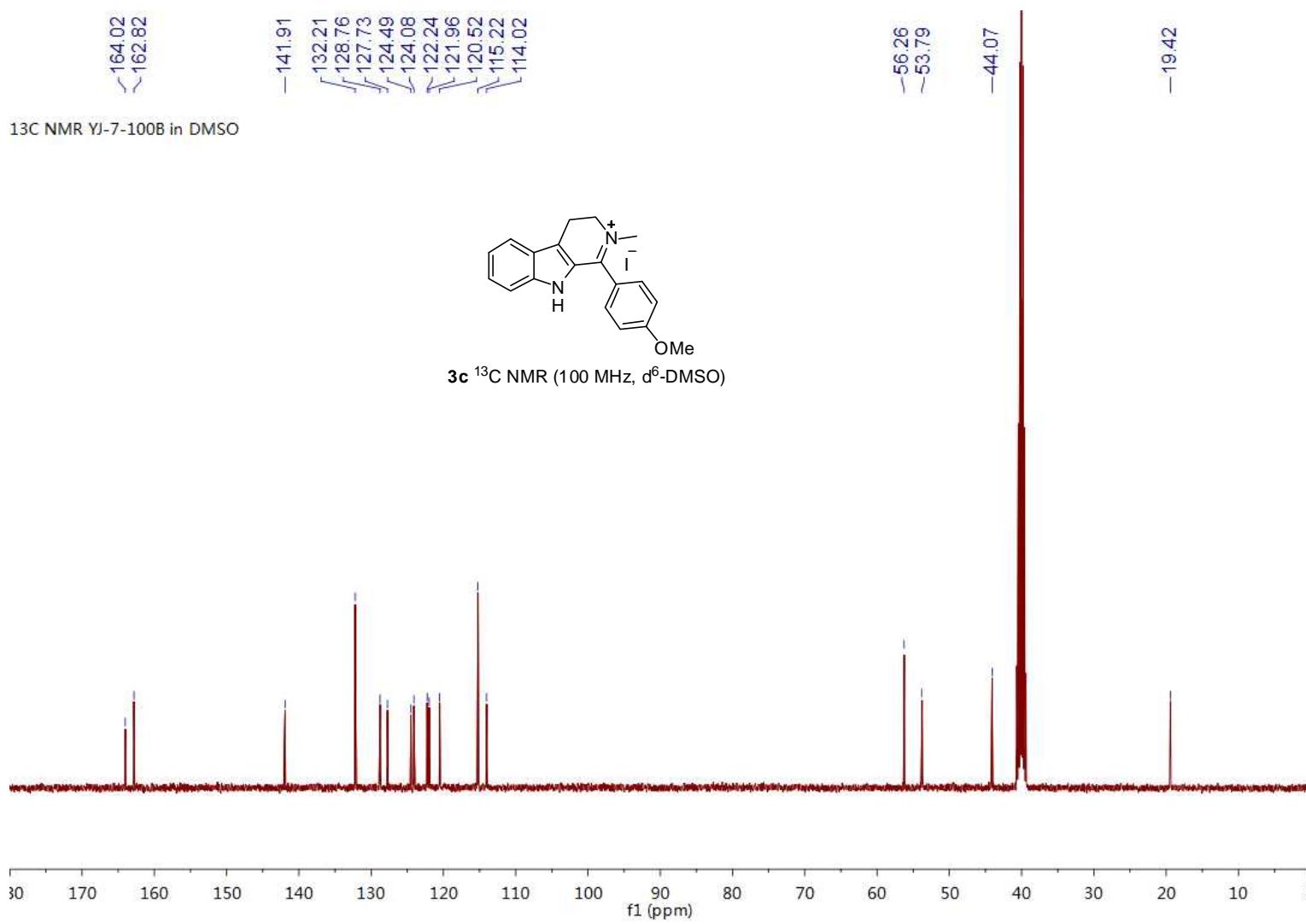


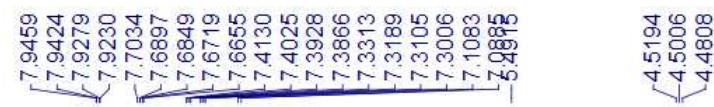








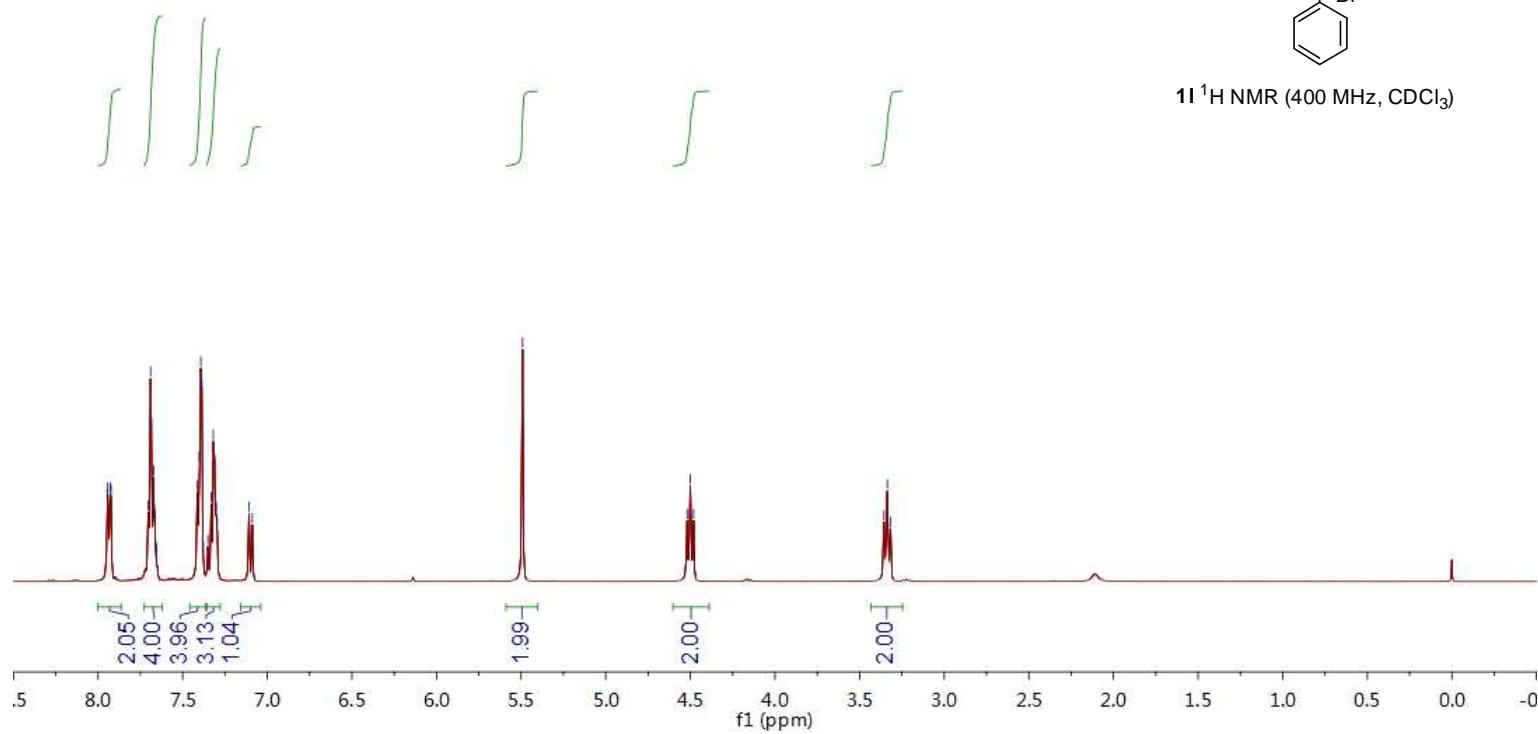


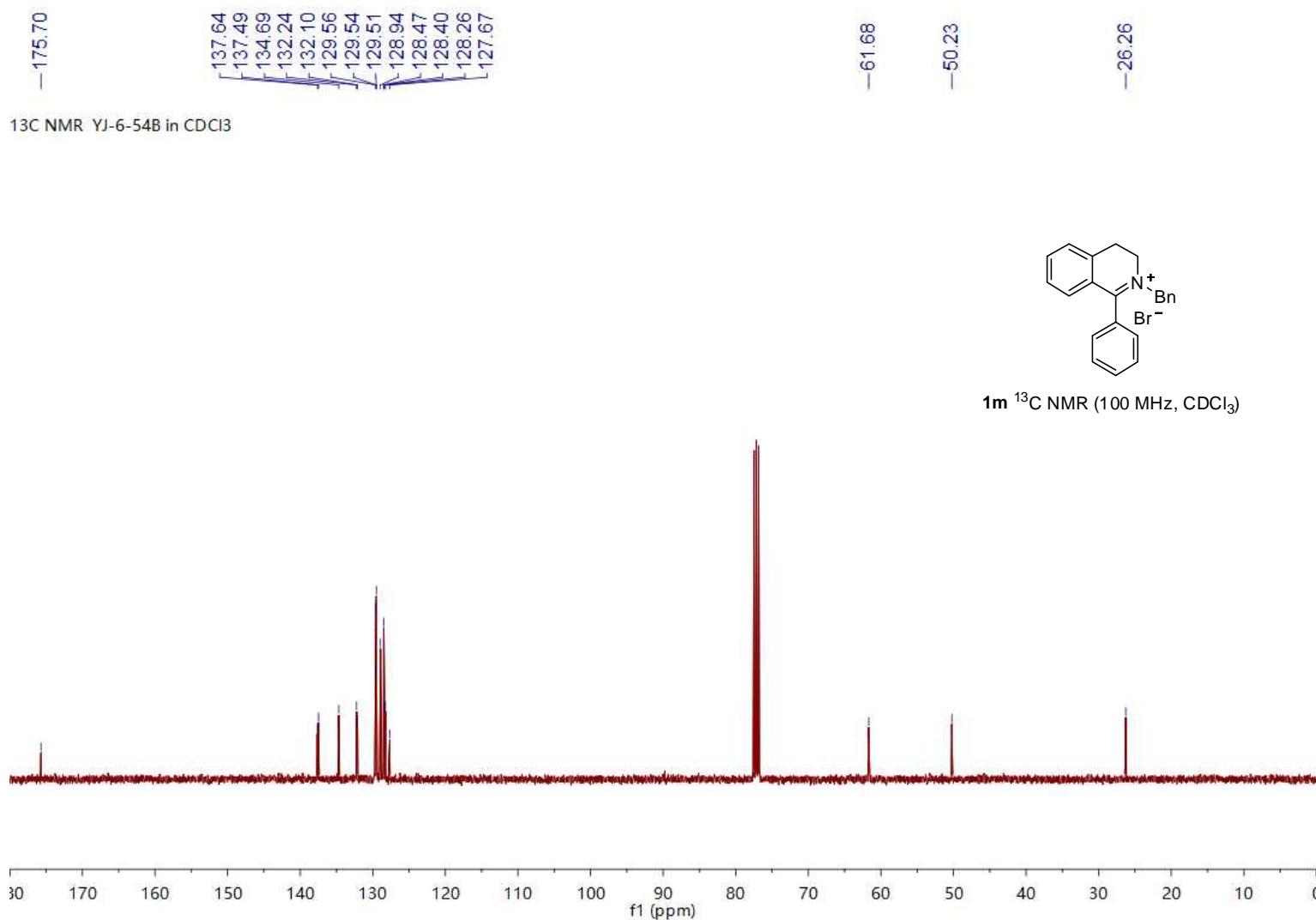


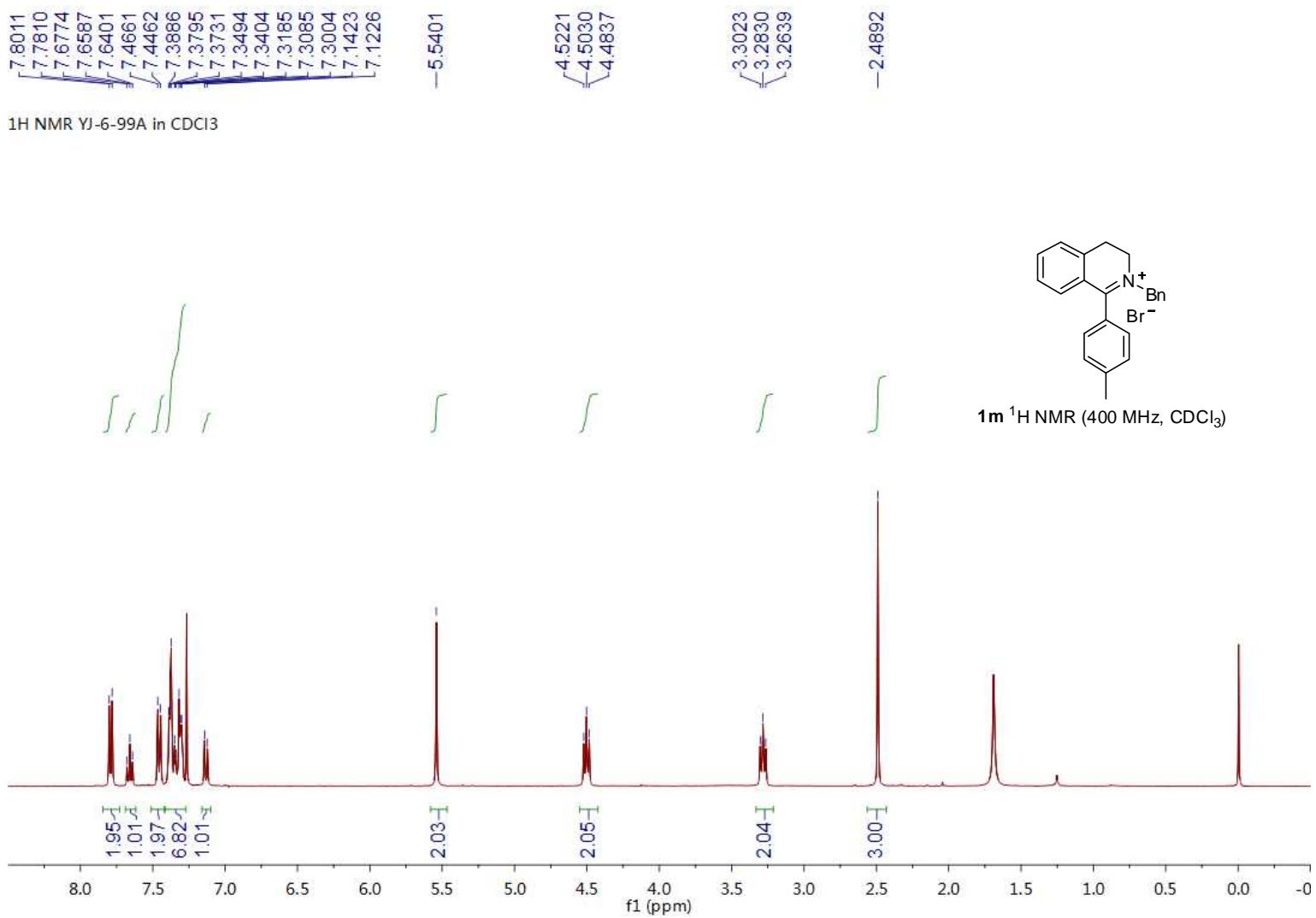
1H NMR YJ-6-54B in CDCl_3



11 ^1H NMR (400 MHz, CDCl_3)







-176.58

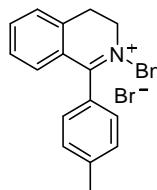
142.62
138.67
137.49
134.30
132.85
130.19
129.55
129.32
128.84
128.69
128.62
128.46
127.26

^{13}C NMR YJ-6-99A in DMSO

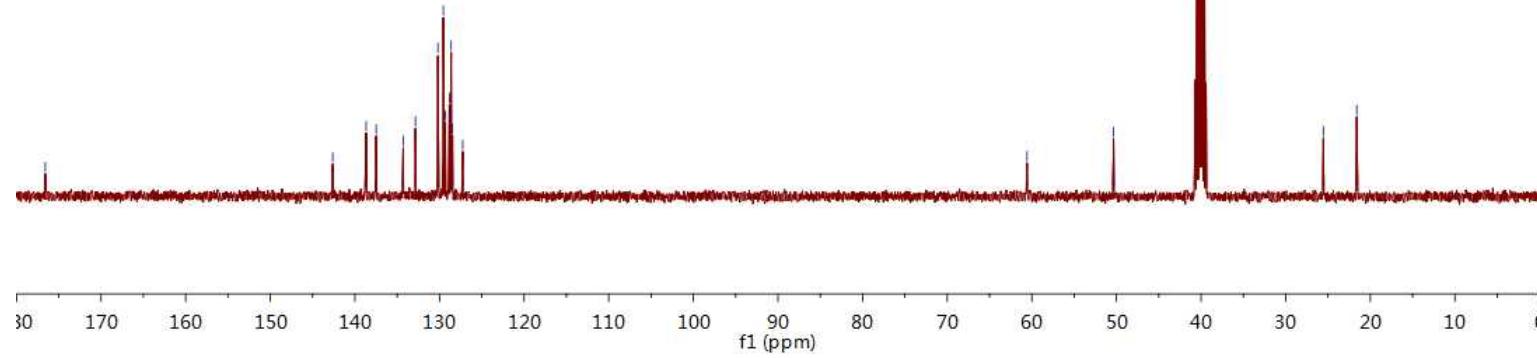
-60.57

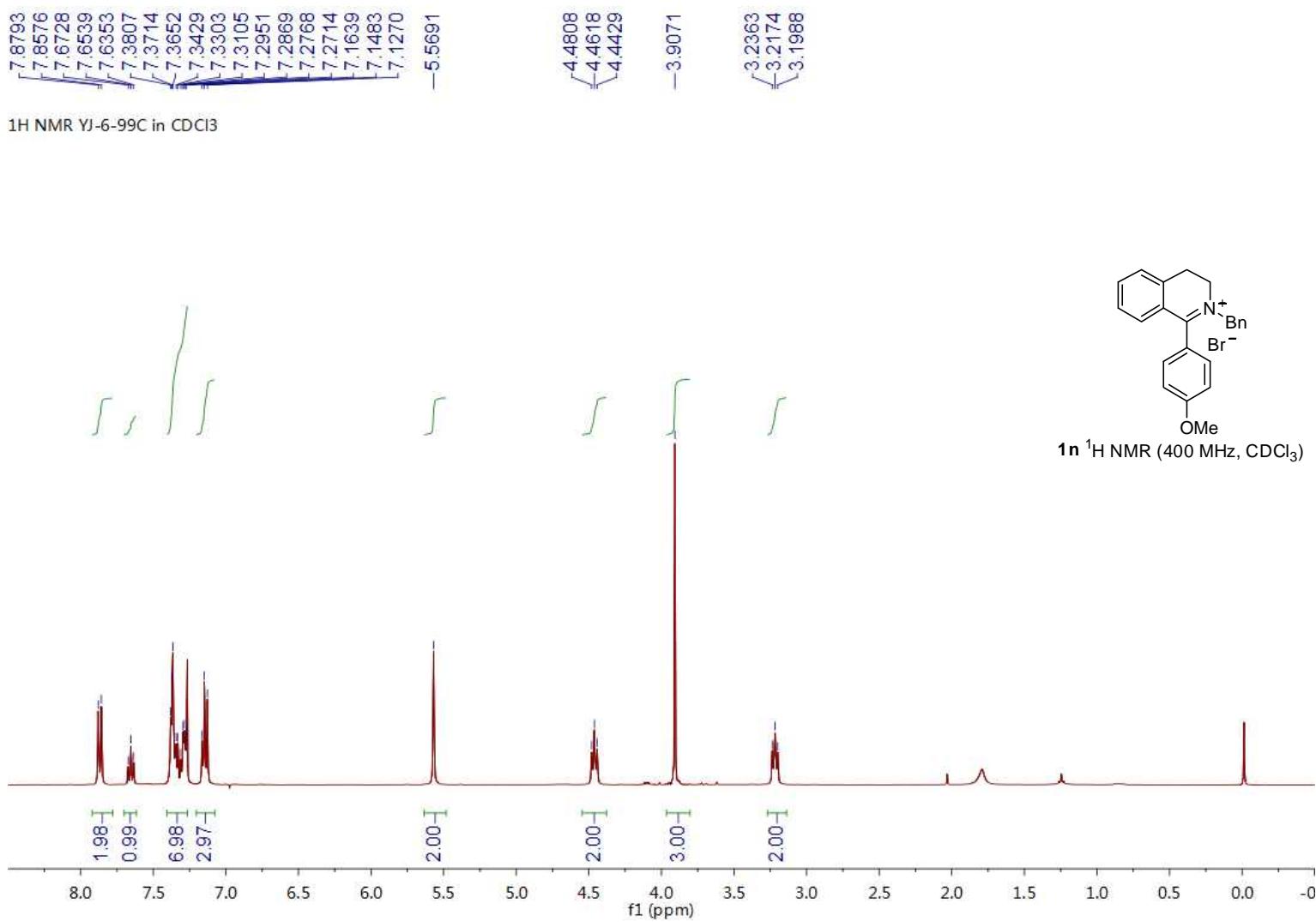
-50.33

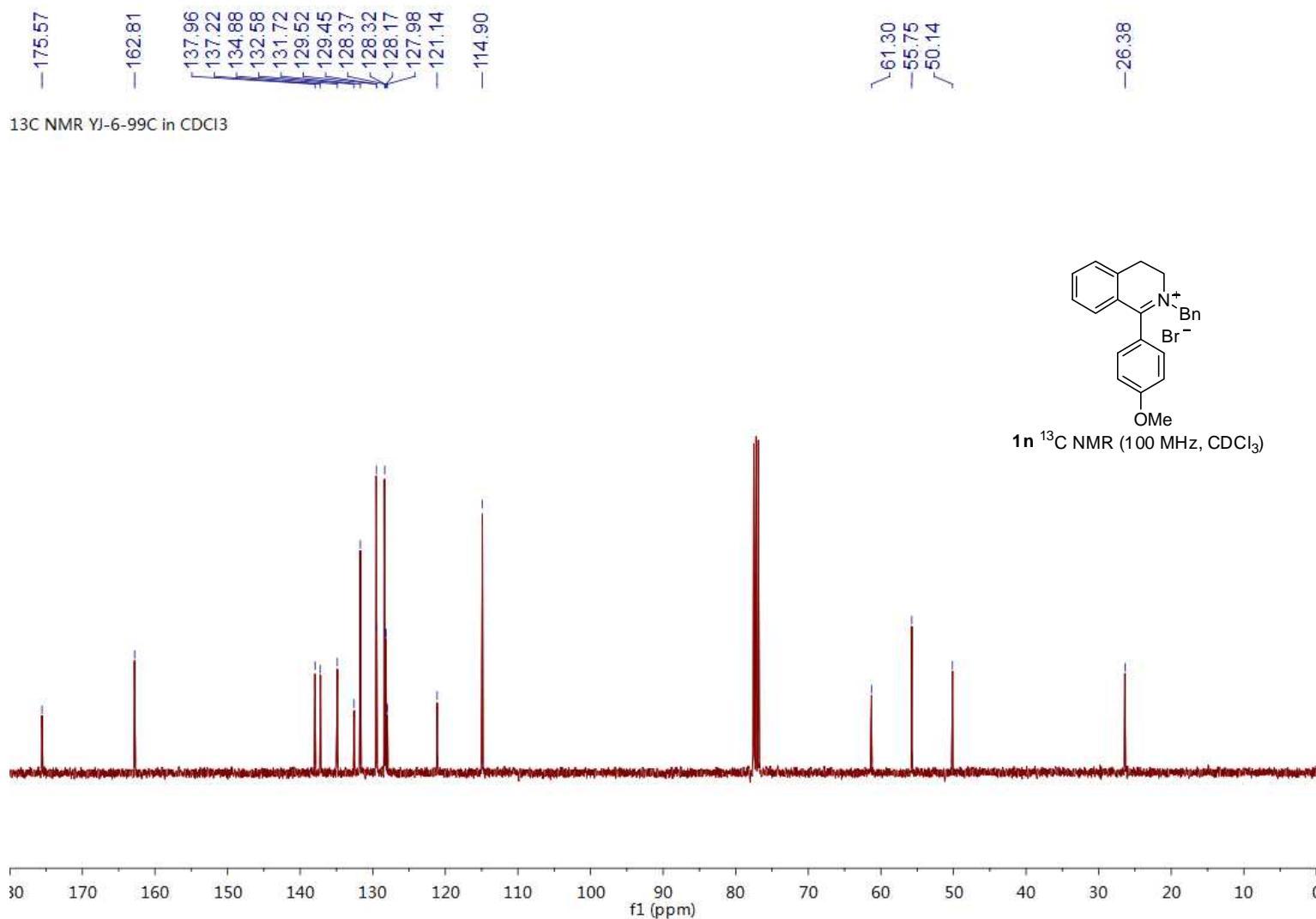
-25.55
-21.60



1m ^{13}C NMR (100 MHz, $\text{d}^6\text{-DMSO}$)





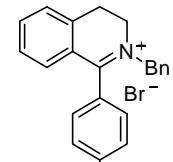


7.9999
7.9789
7.6921
7.6731
7.6539
7.6329
7.4150
7.4061
7.3966
7.3908
7.3597
7.3400
7.3279
7.3211
7.3026
7.2940
7.2847
7.0943
5.9346

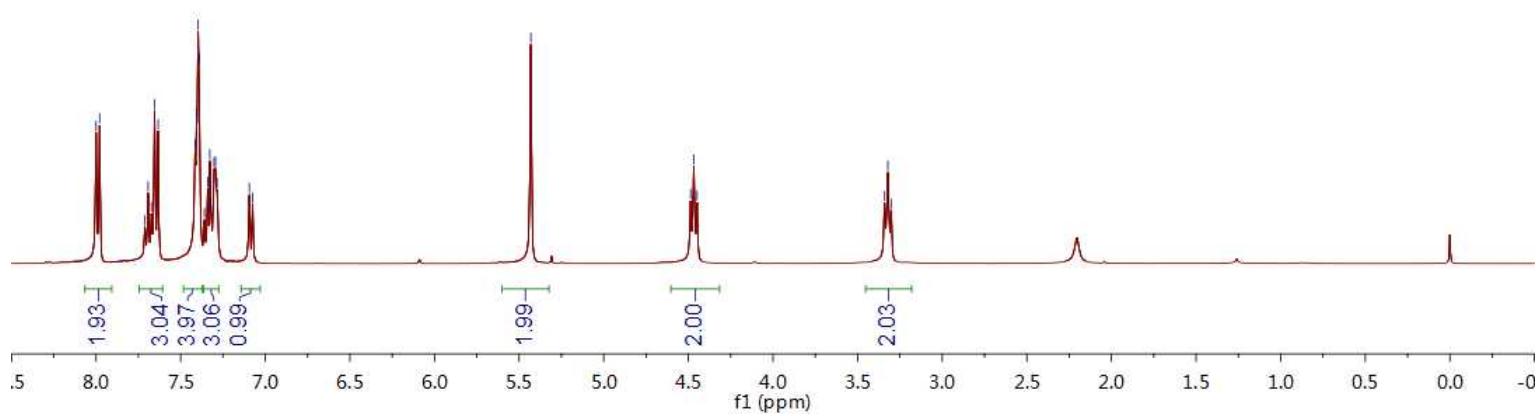
^1H NMR YJ-6-99B in CDCl_3

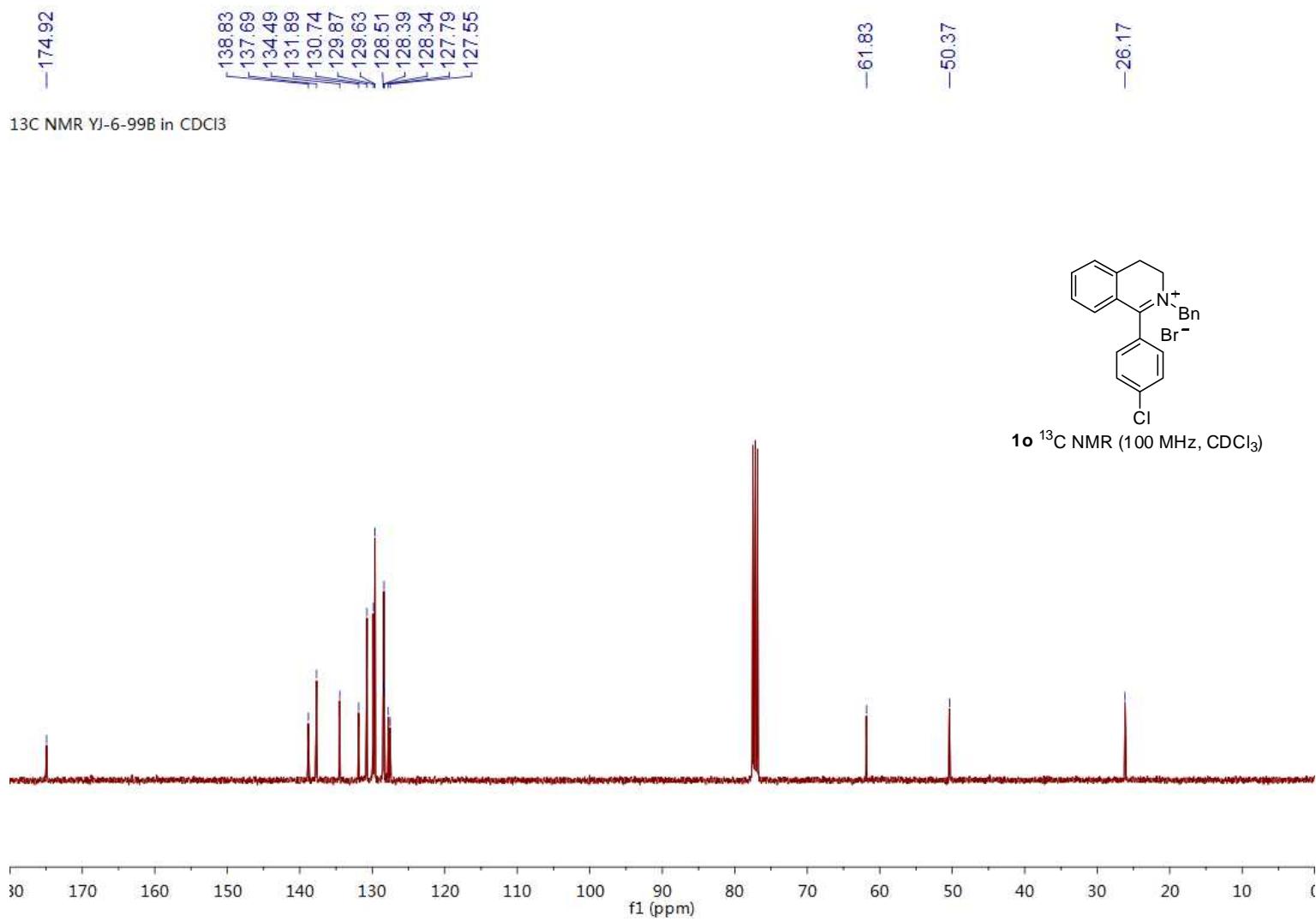
4.4868
4.4680
4.4484

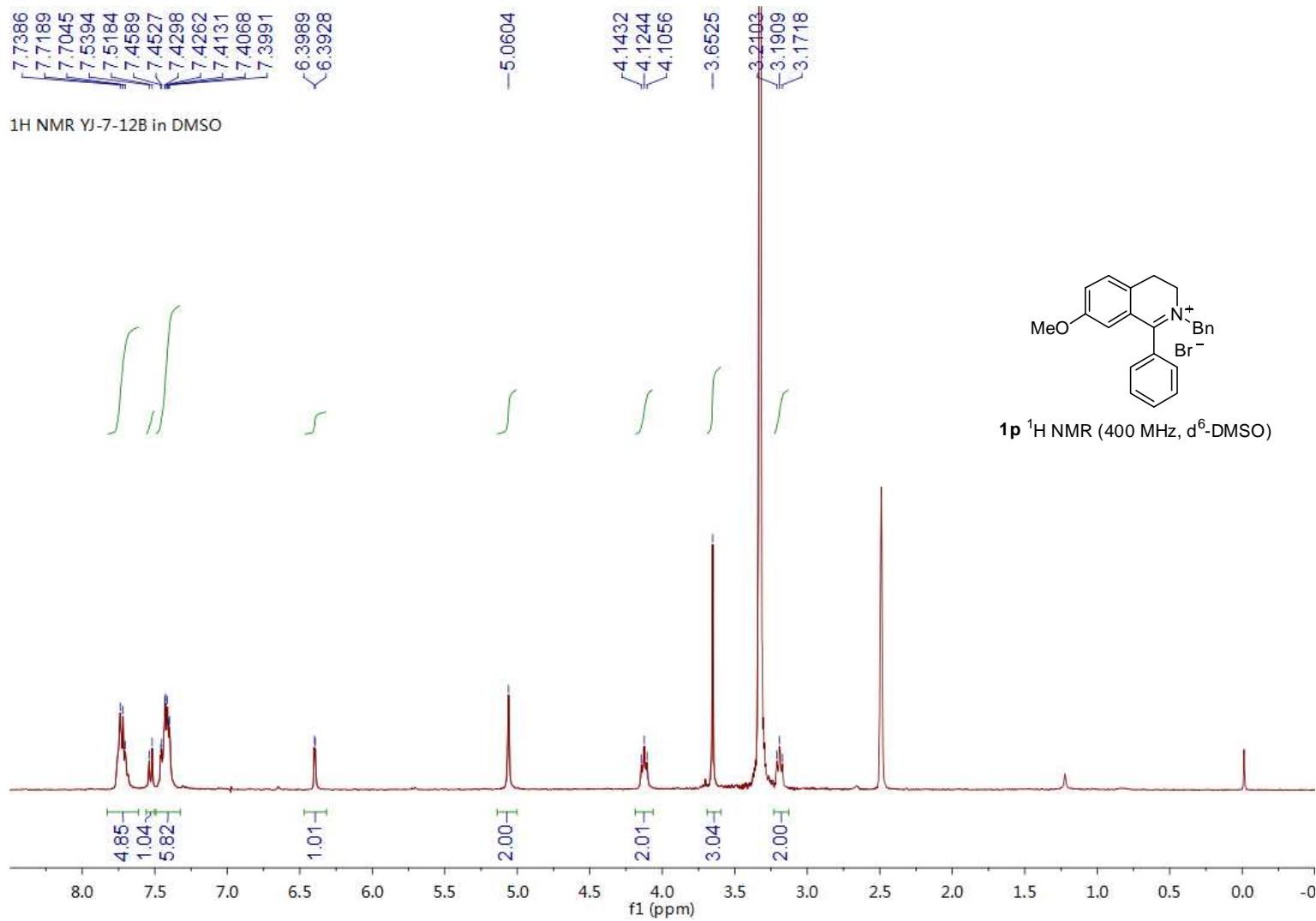
3.3402
3.3211
3.3020

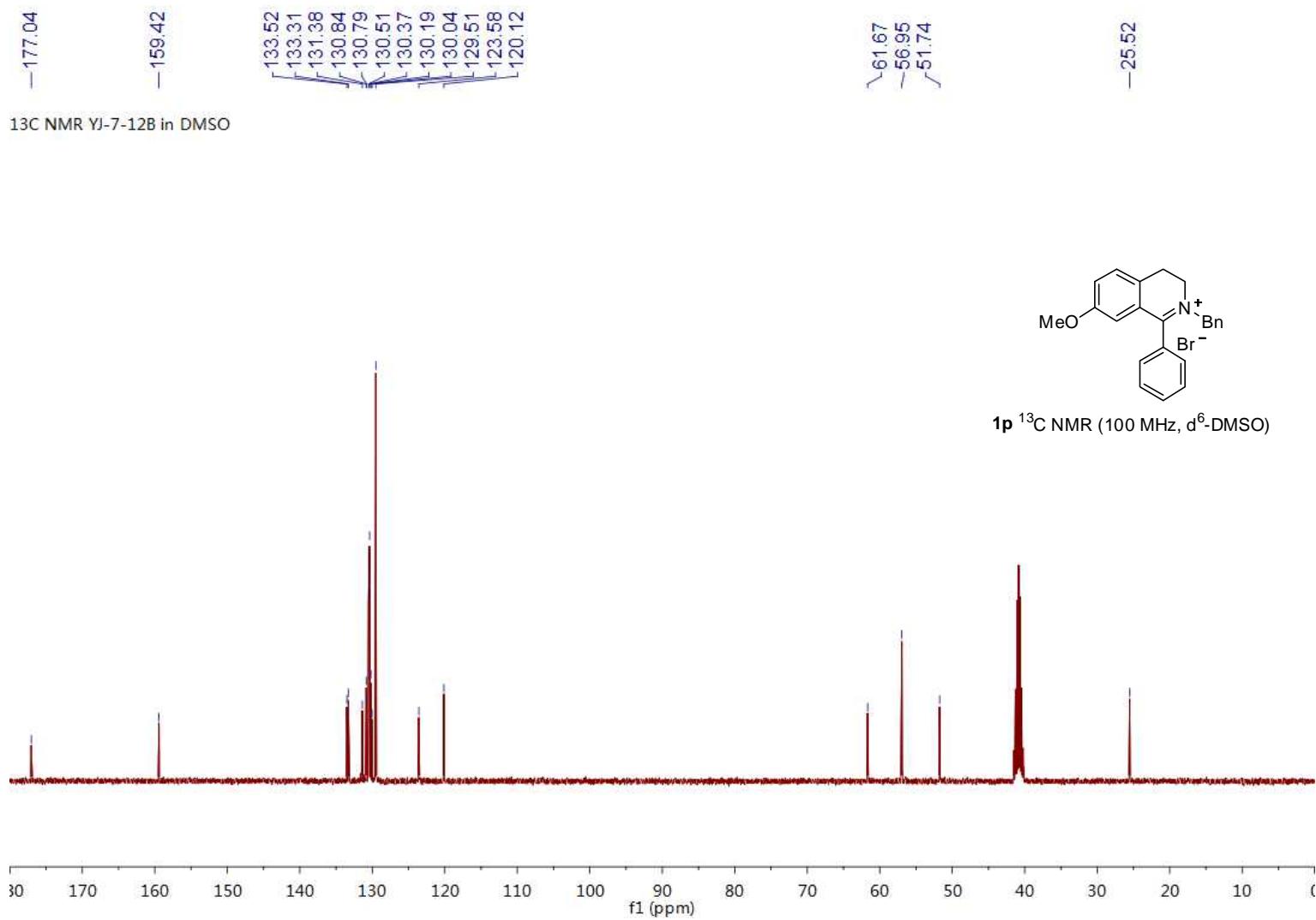


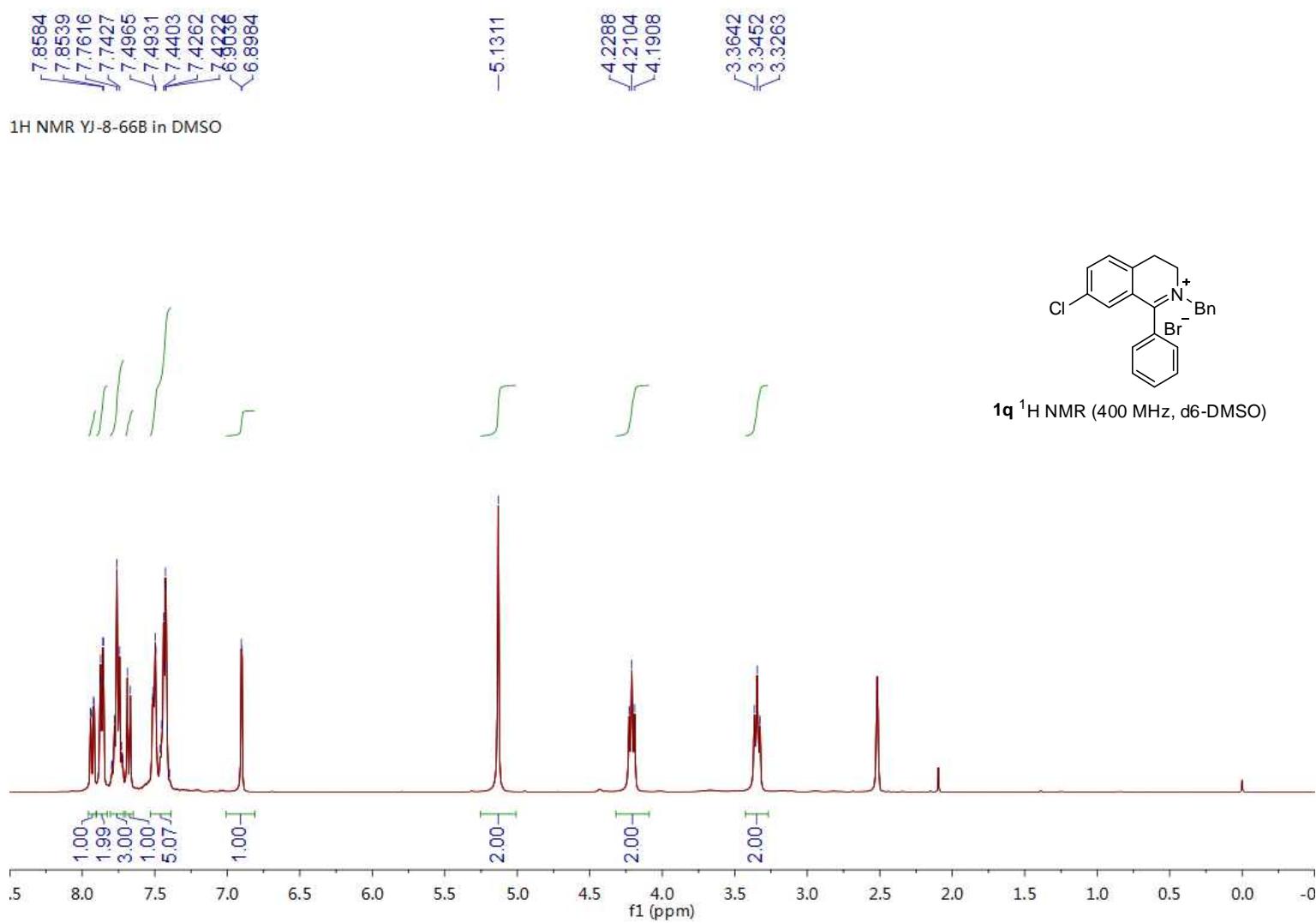
1o ^1H NMR (400 MHz, CDCl_3)

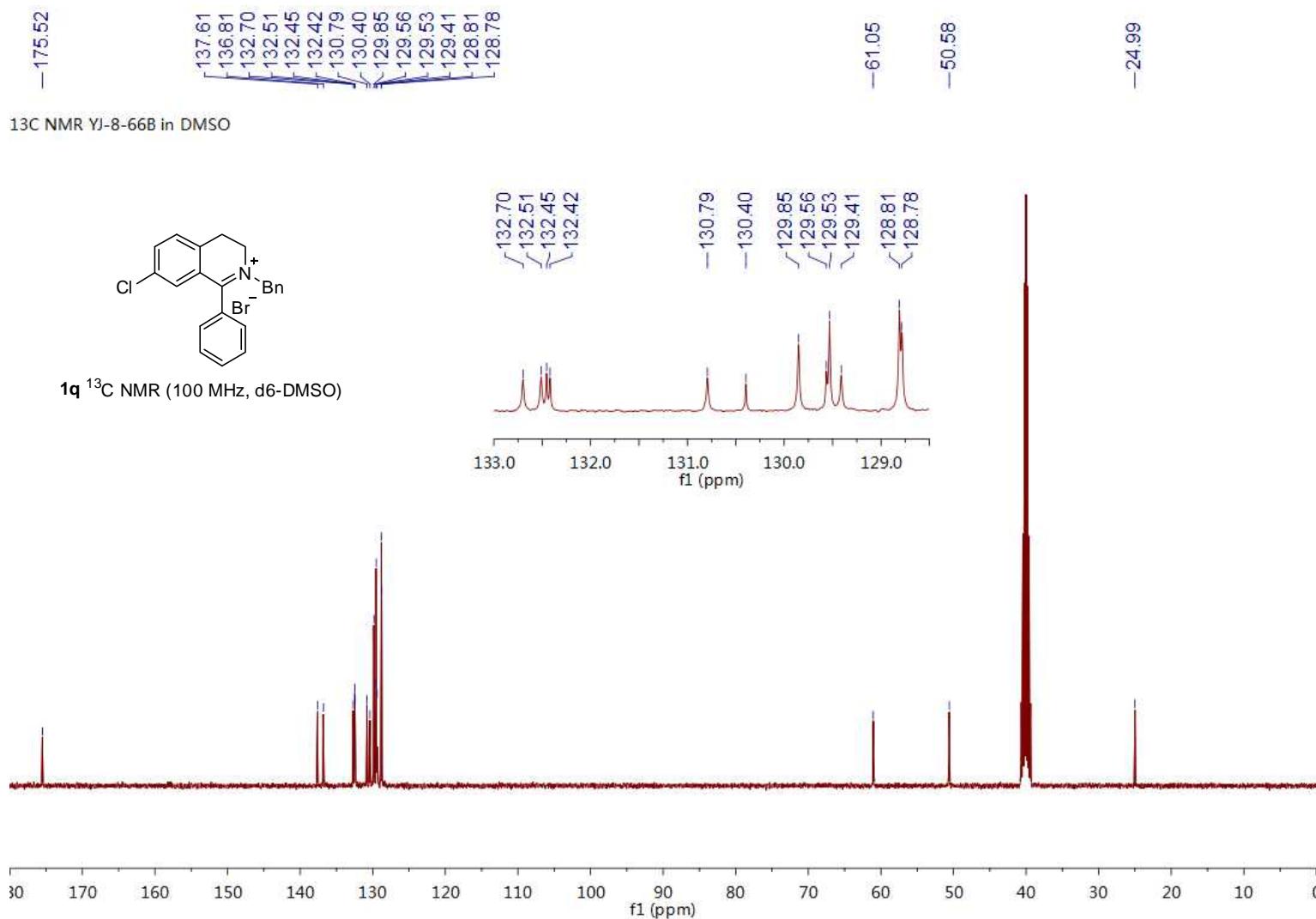










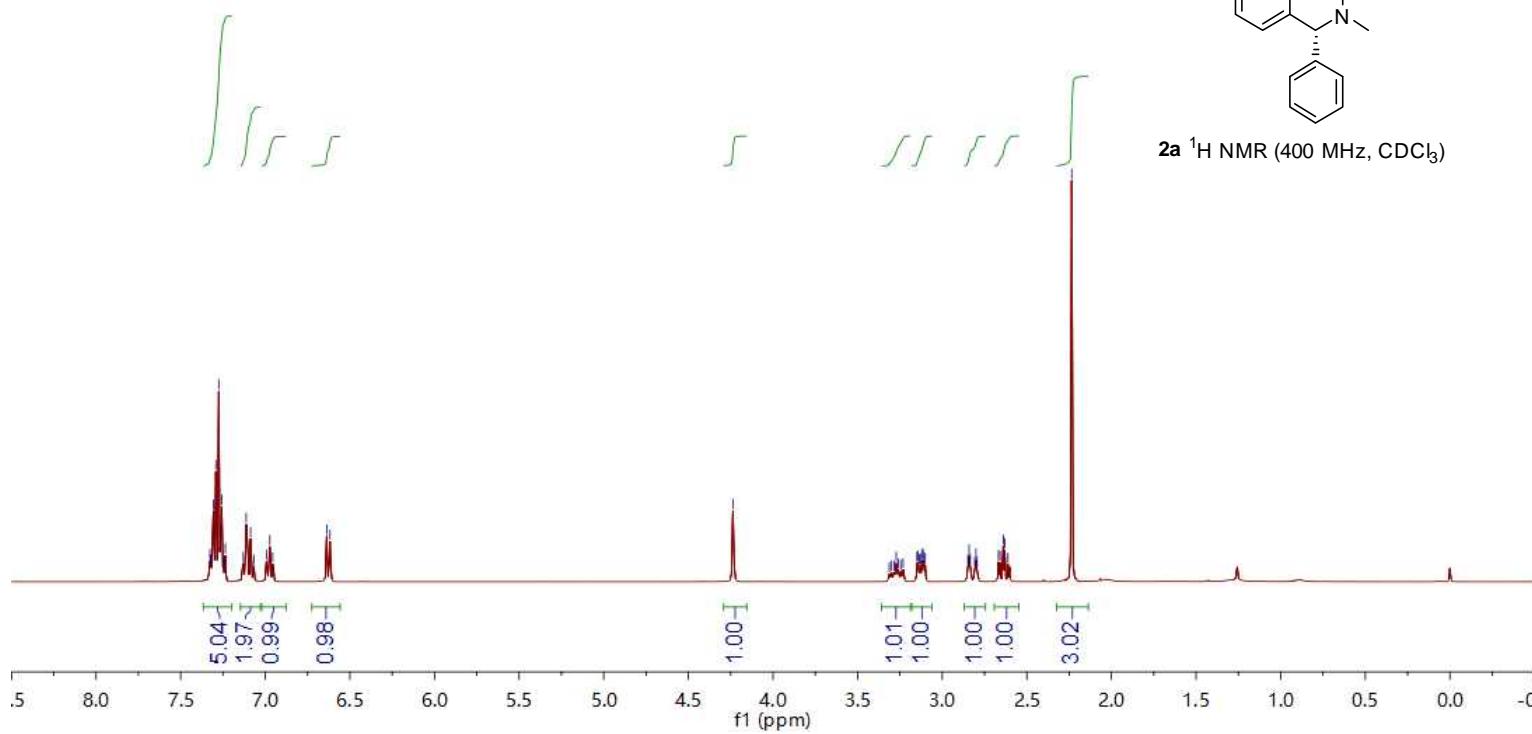
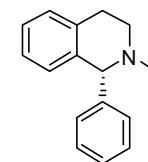


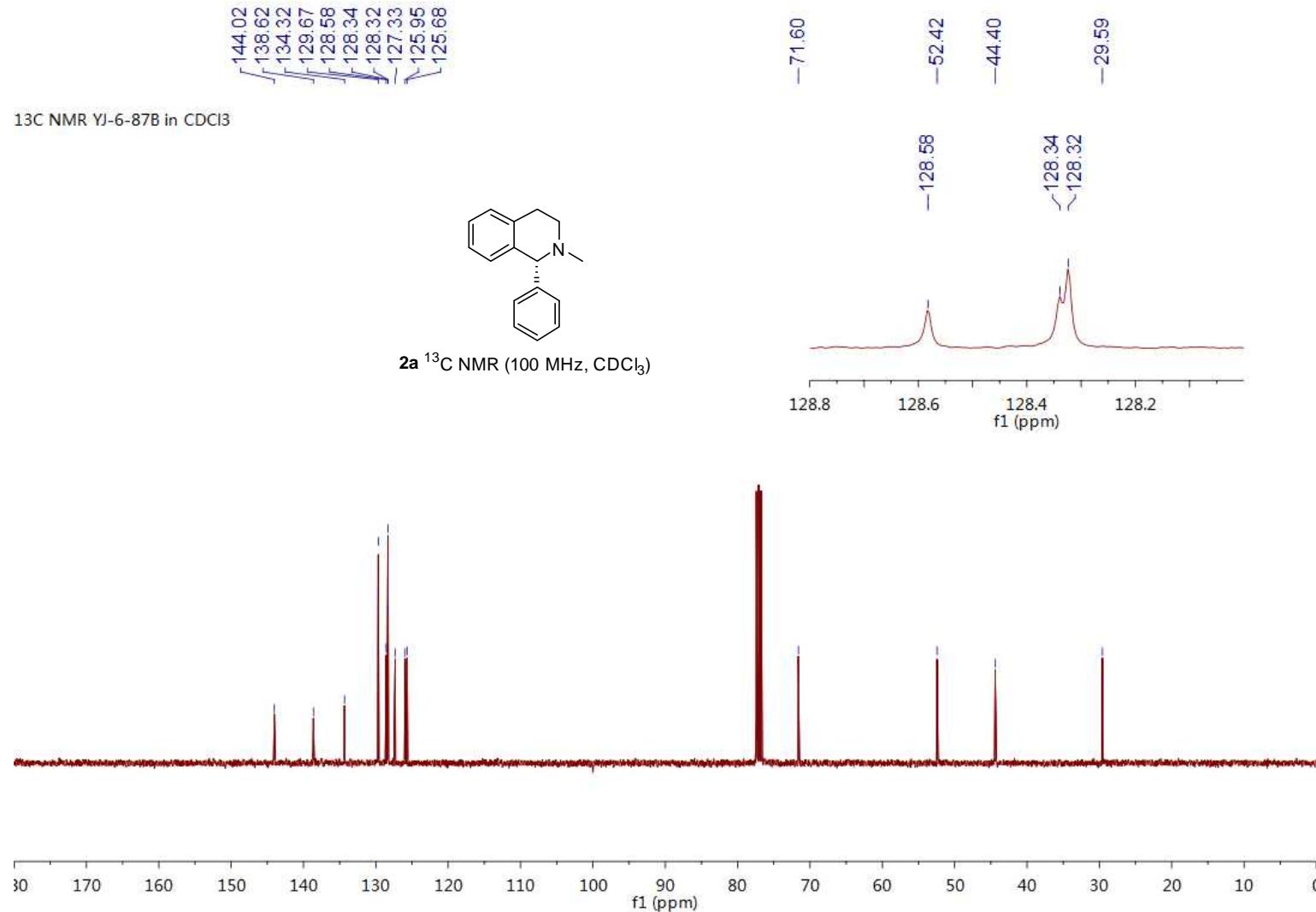
7.3270
7.3071
7.2922
7.2748
7.2641
7.2570
7.2356
7.1128
7.0868
6.9927
6.9370
6.6175

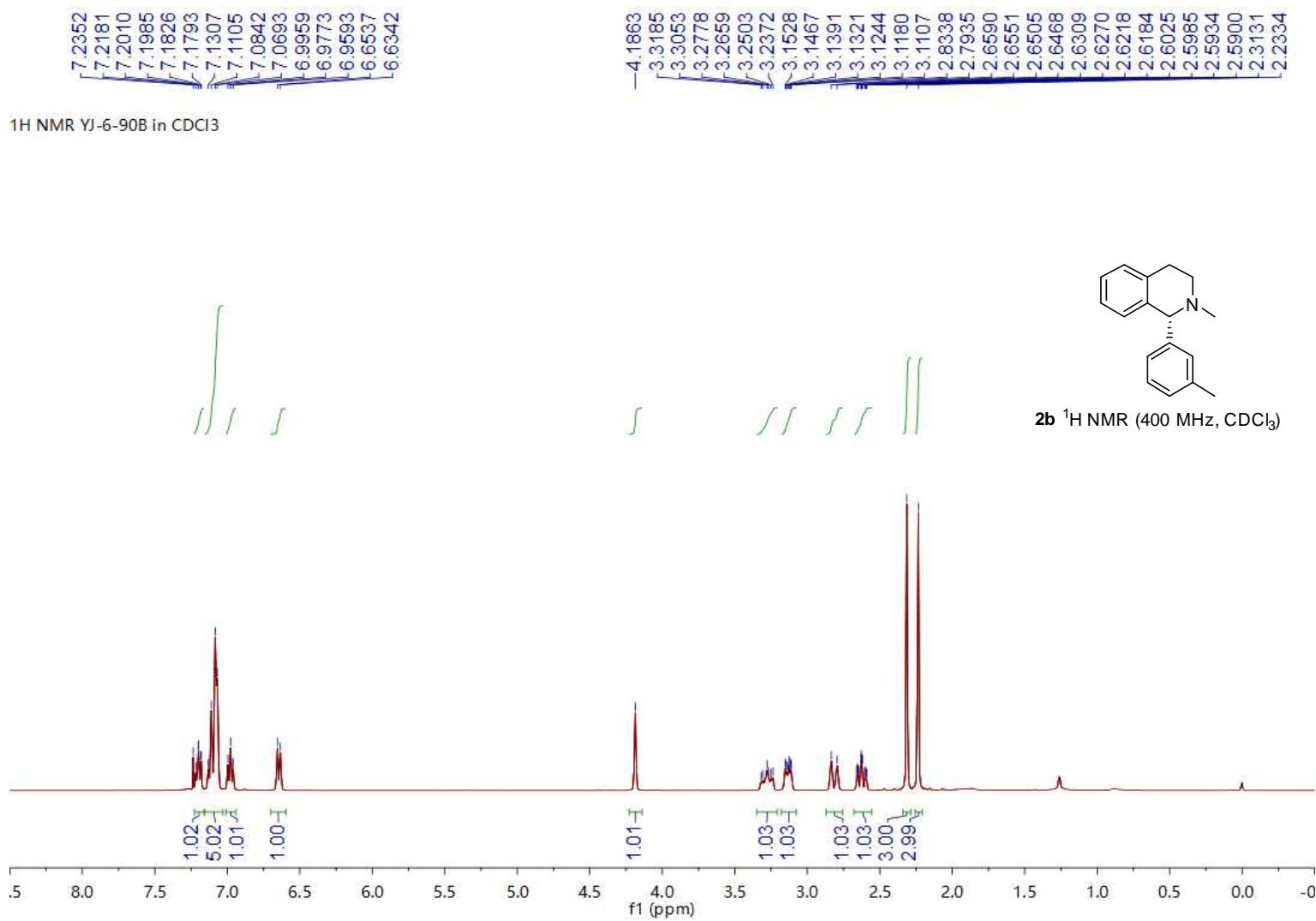
1H NMR YJ-6-87B in CDCl₃

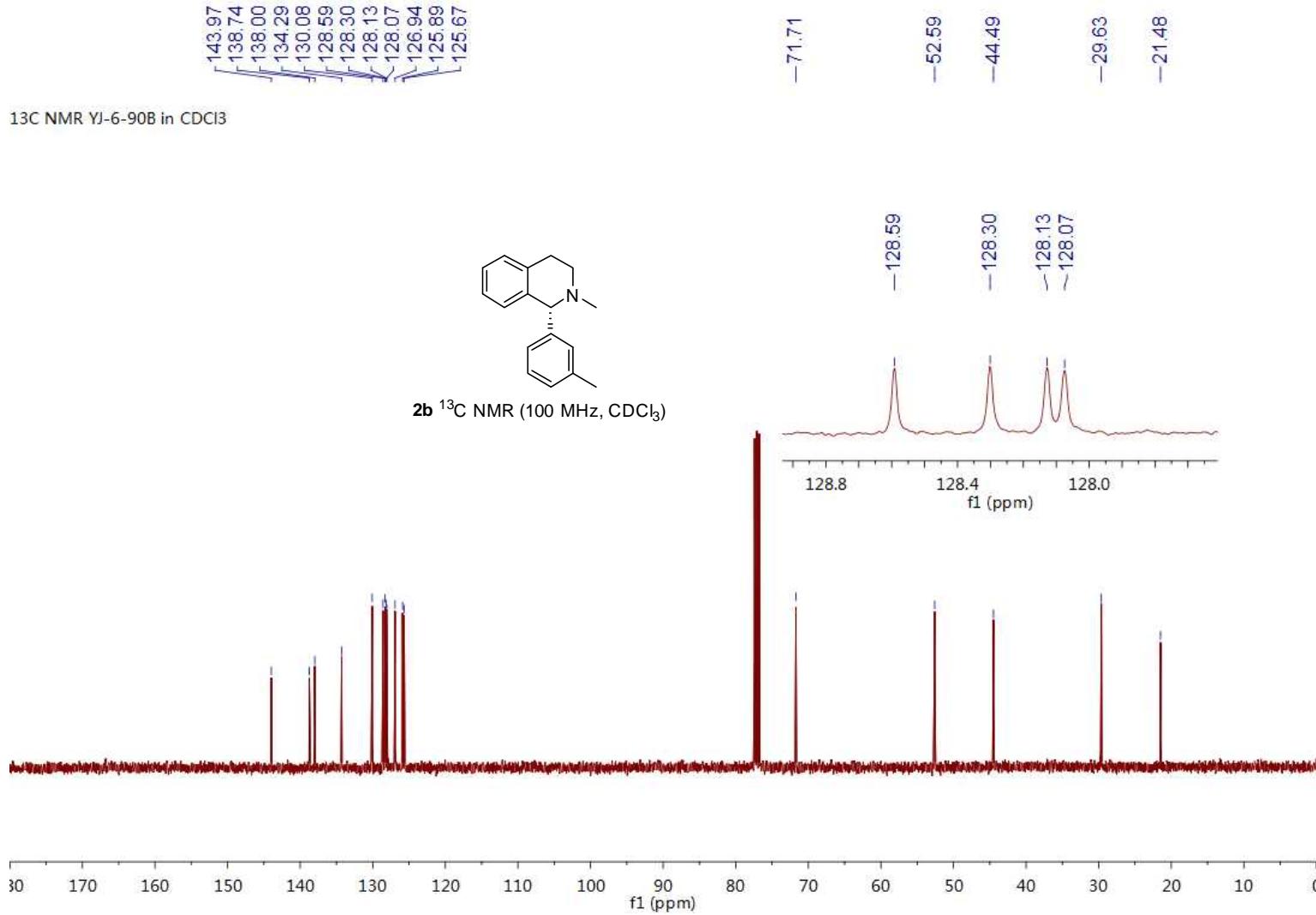
-4.2362
3.2715
3.1490
3.1423
3.1206
3.1136
3.1068
2.8410
2.8004
2.6677
2.6583
2.6393
2.2348

2a ¹H NMR (400 MHz, CDCl₃)



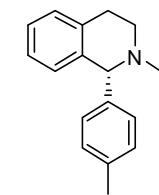
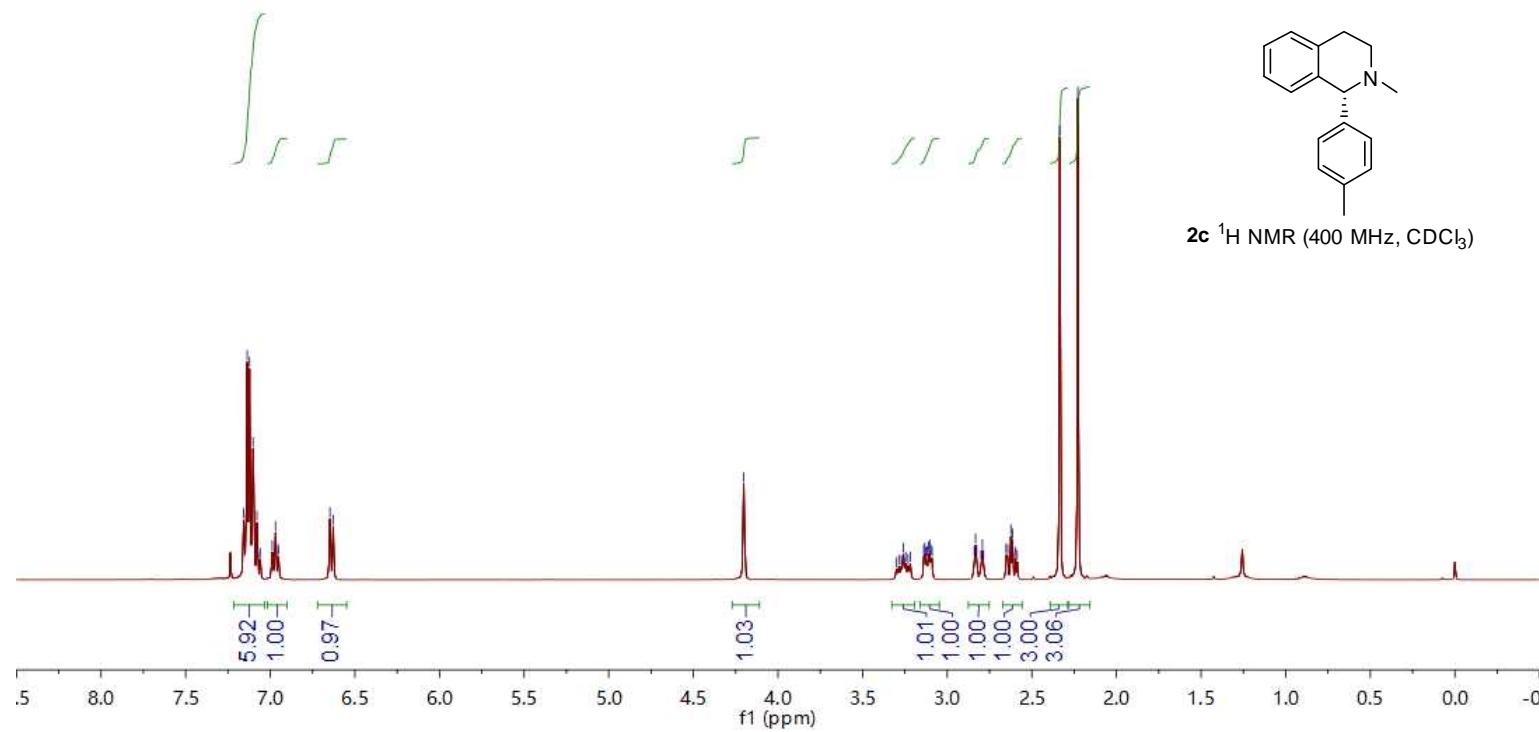




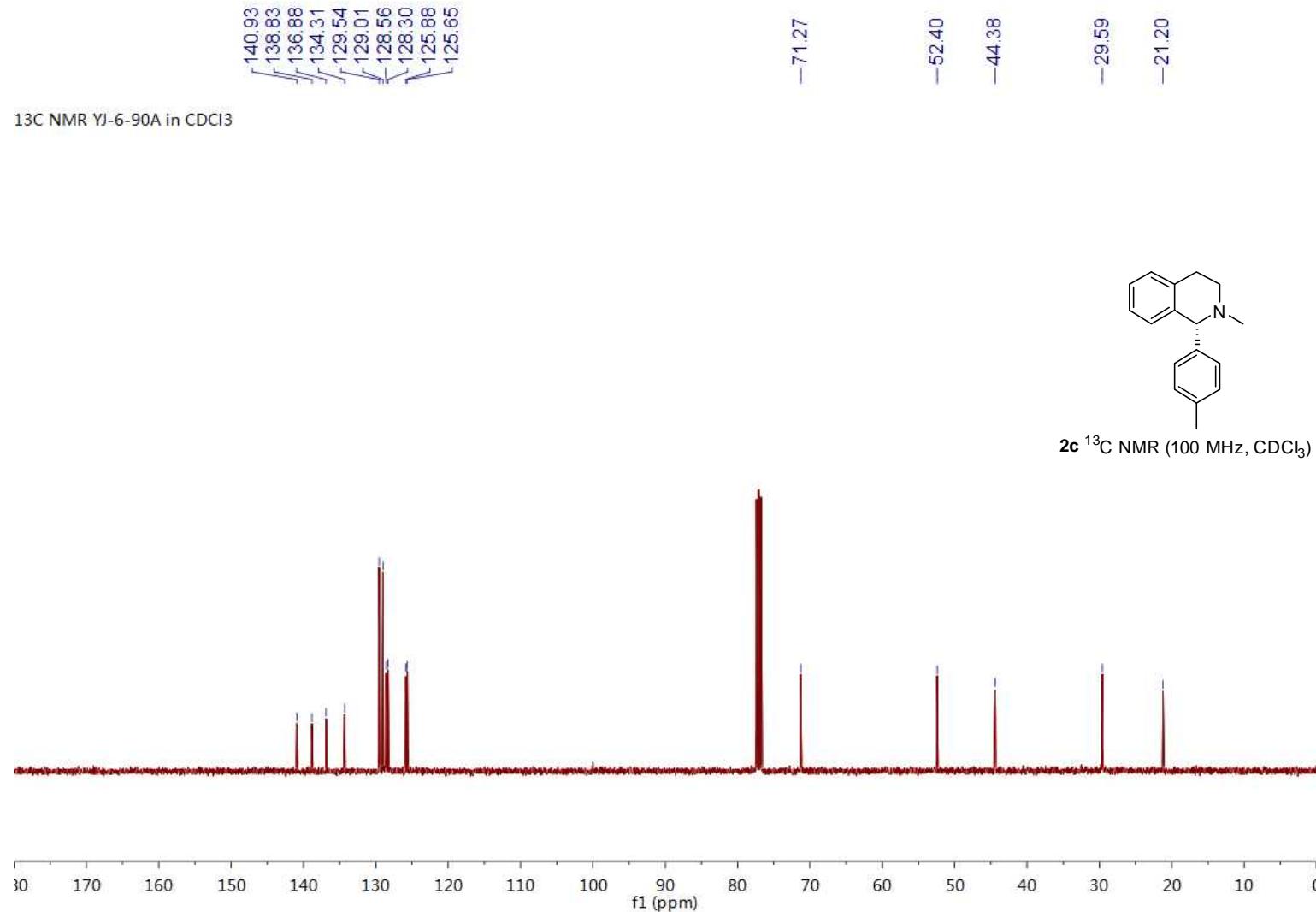


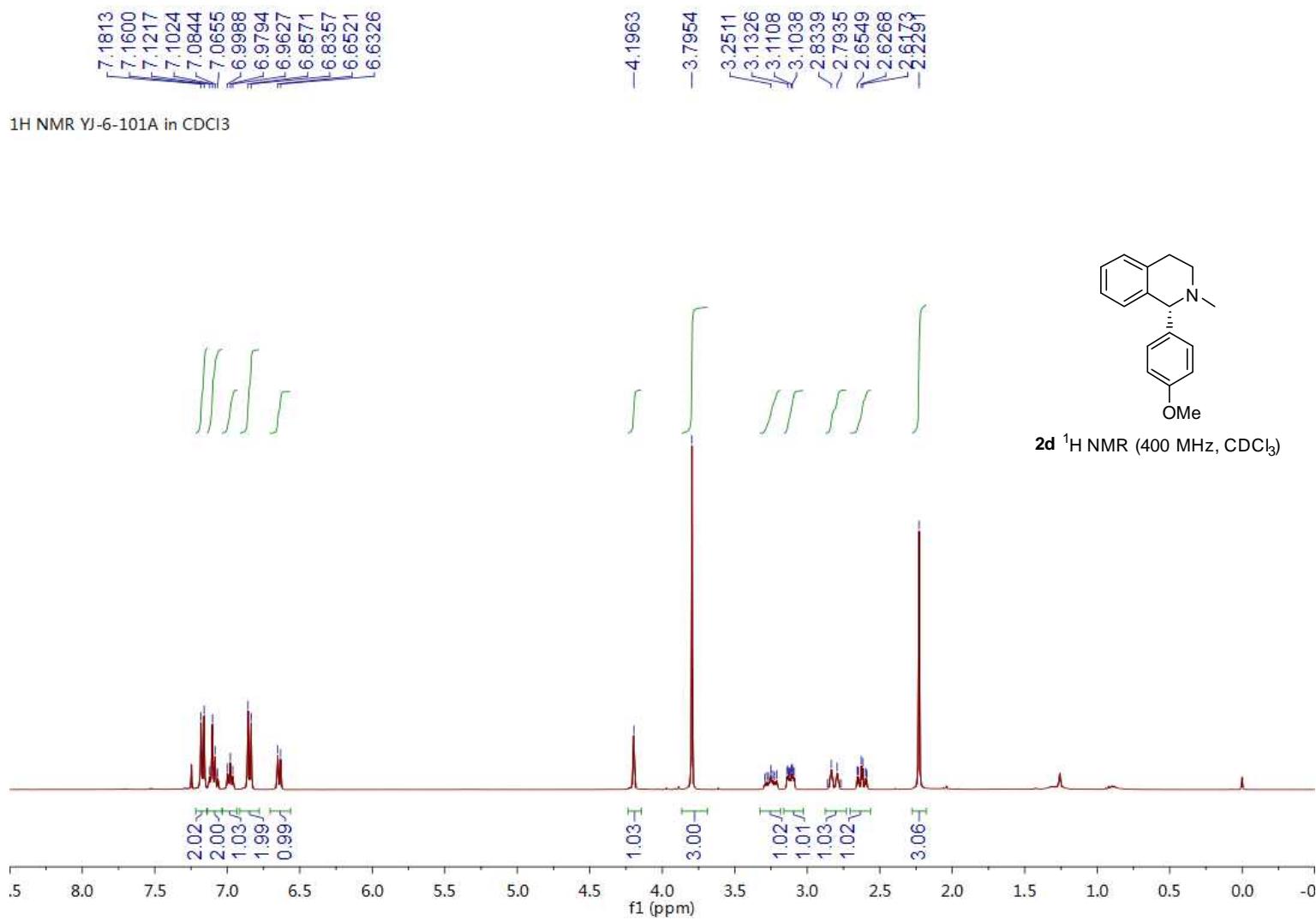


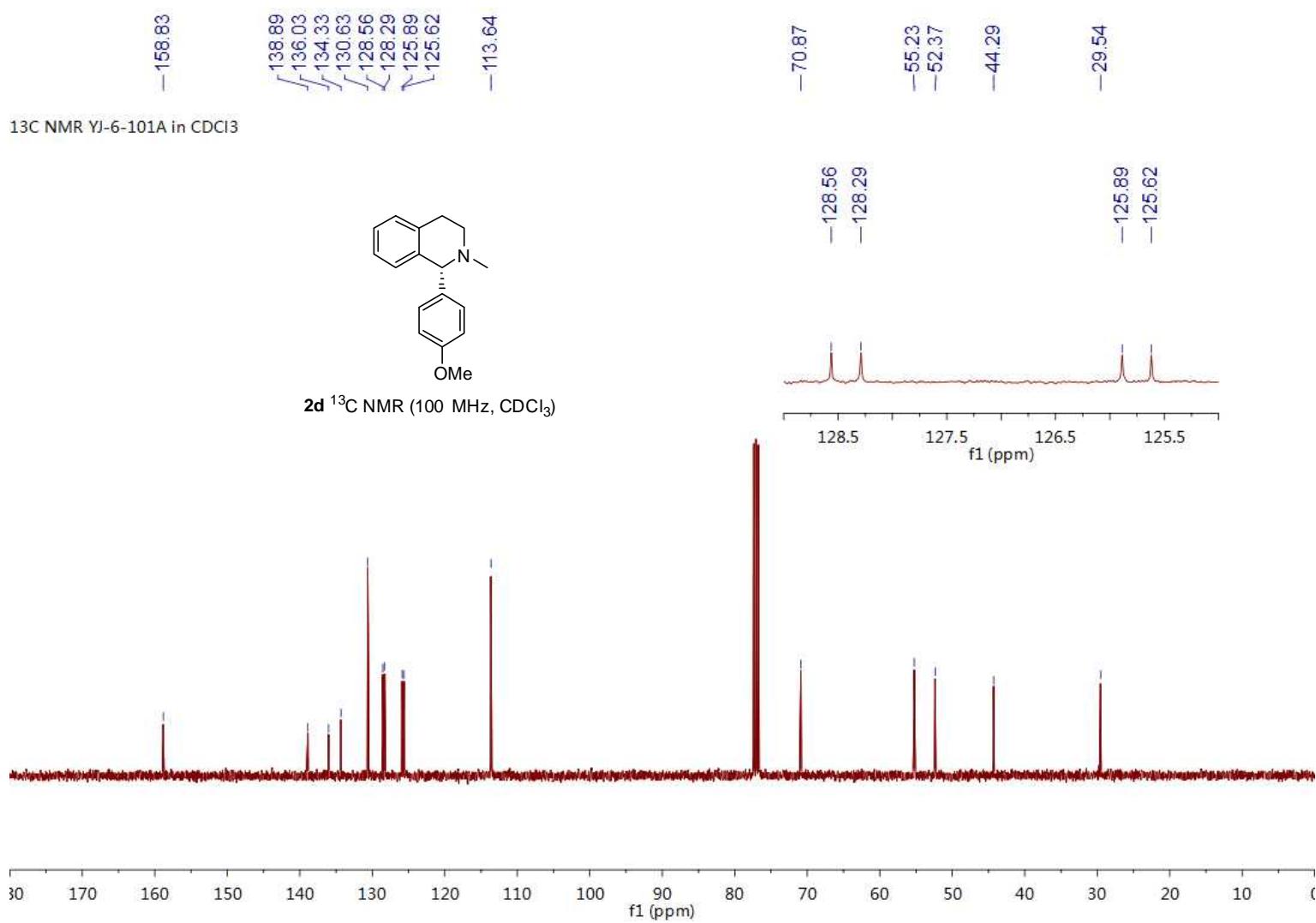
¹H NMR YJ-6-90A in CDCl₃



2c ^1H NMR (400 MHz, CDCl_3)



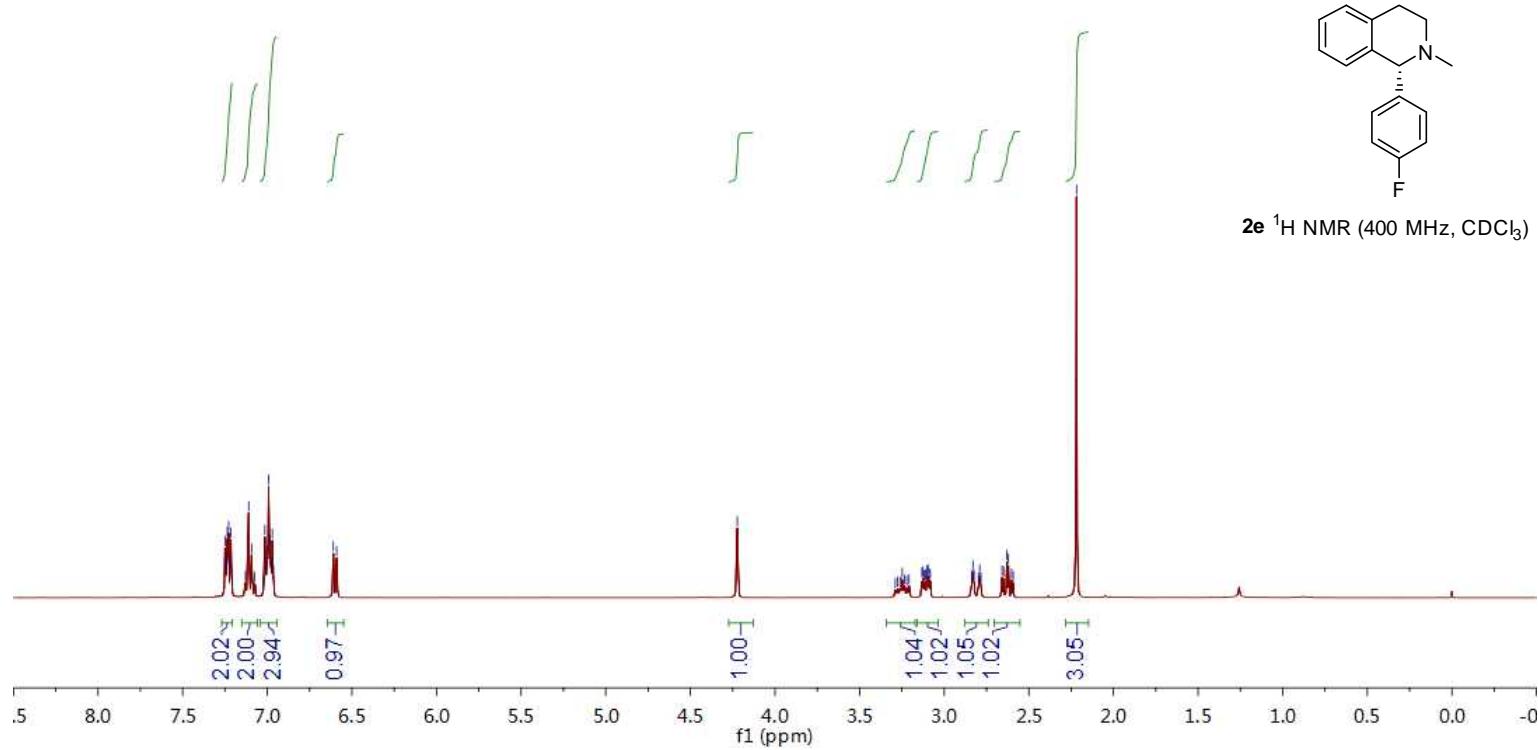


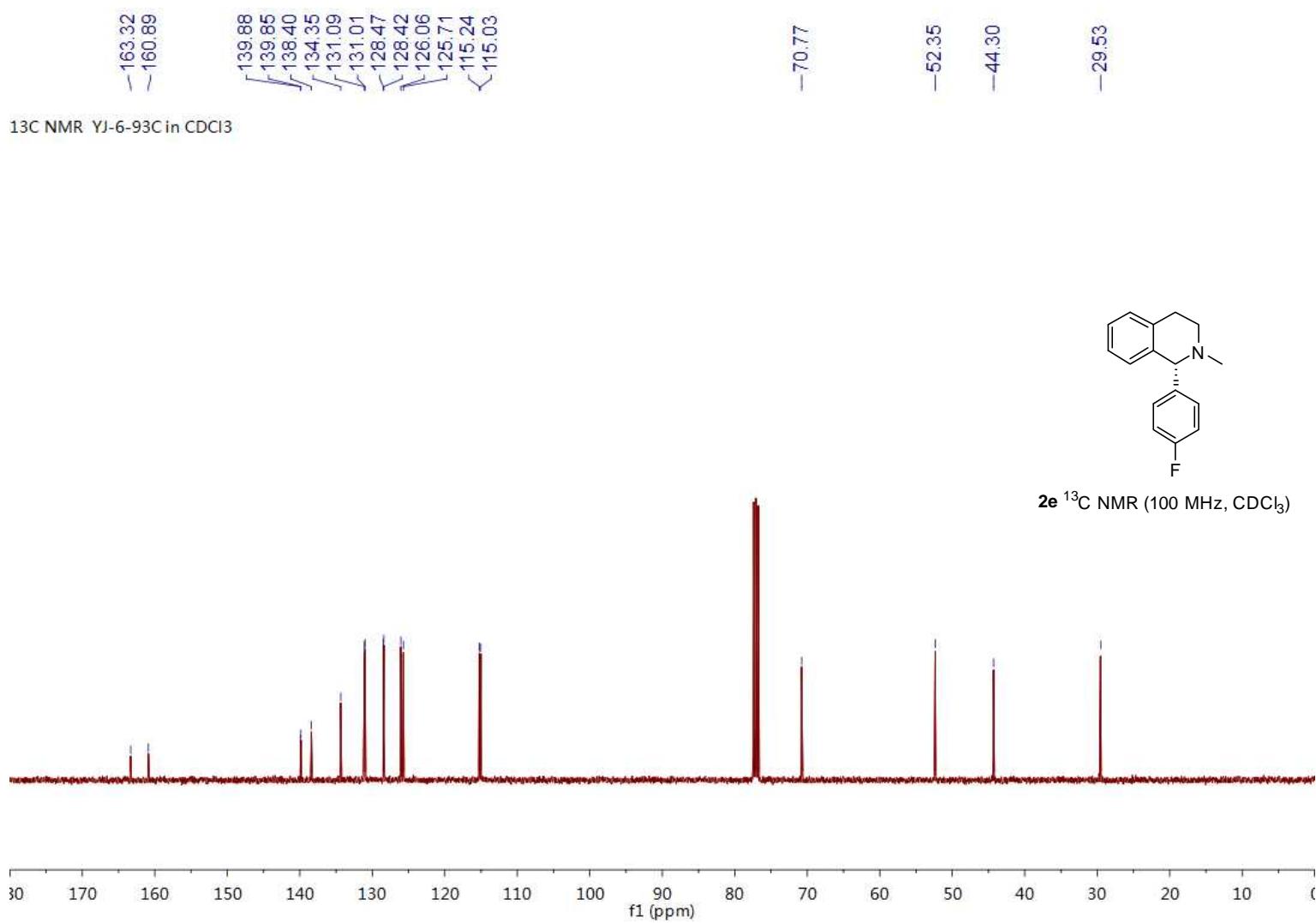


7.2498
7.2360
7.2284
7.2145
7.1099
7.0927
7.0141
6.9976
6.9924
6.9867
6.9795
6.6097
6.5902

¹H NMR YJ-6-93C in CDCl₃

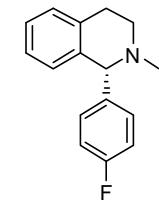
-4.2226
3.2481
3.1316
3.1249
3.1029
3.0962
3.0892
2.8308
2.7904
2.6583
2.6490
2.6300
2.6286



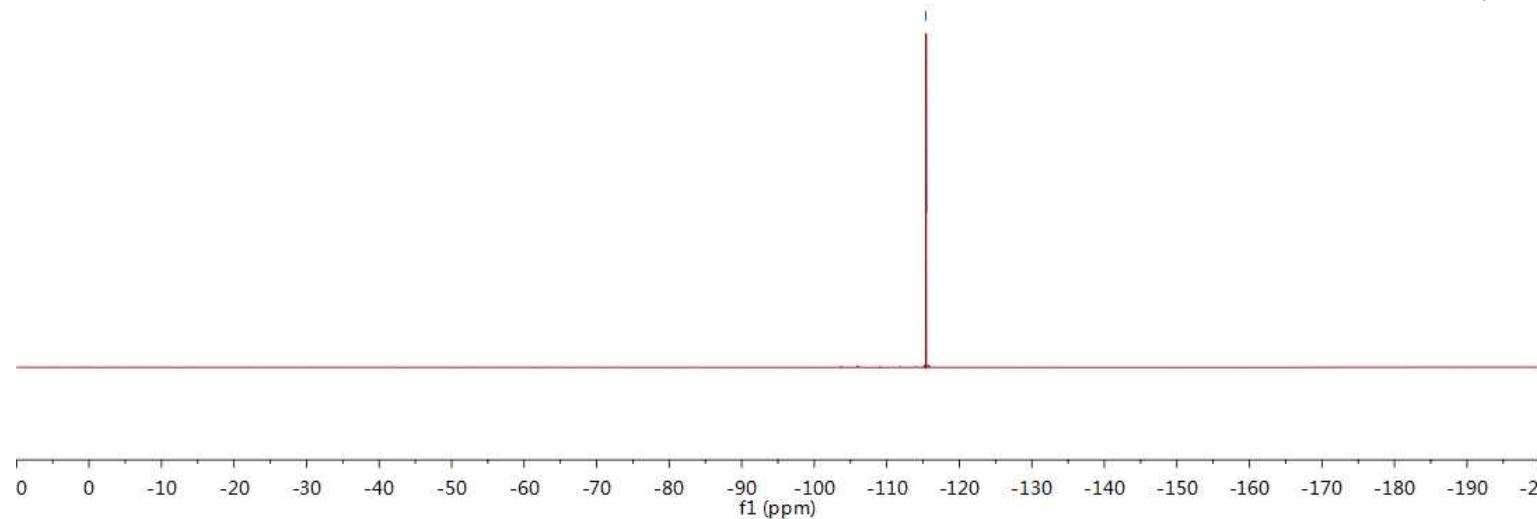


¹⁹F NMR YJ-6-93C in CDCl₃

—115.41



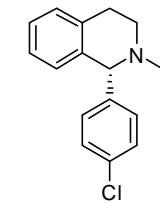
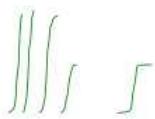
2e ¹⁹F NMR (376 MHz, CDCl₃)



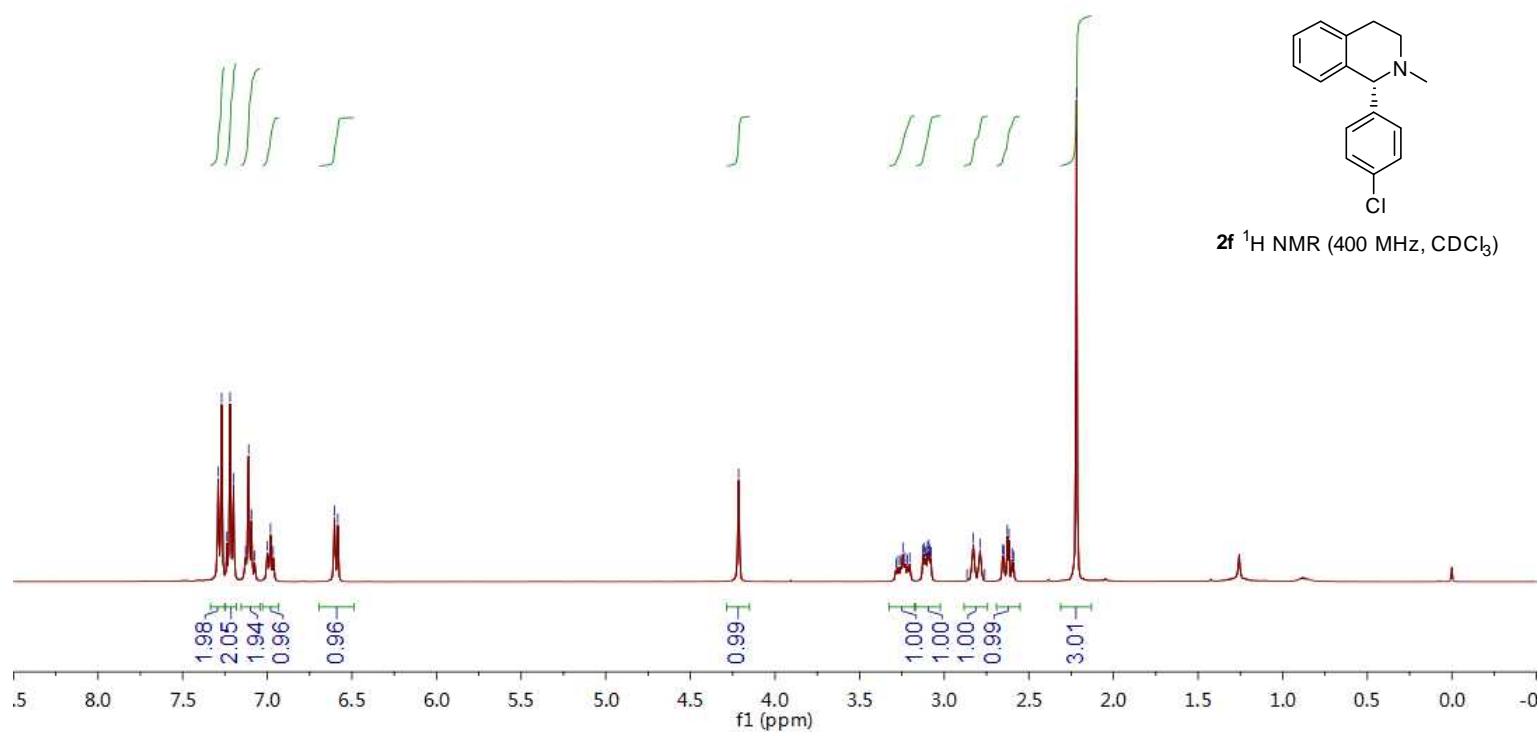
7.2907
7.2698
7.2385
7.2198
7.1989
7.1278
7.1100
7.0931
7.0005
6.9808
6.9647
6.6828
6.5833

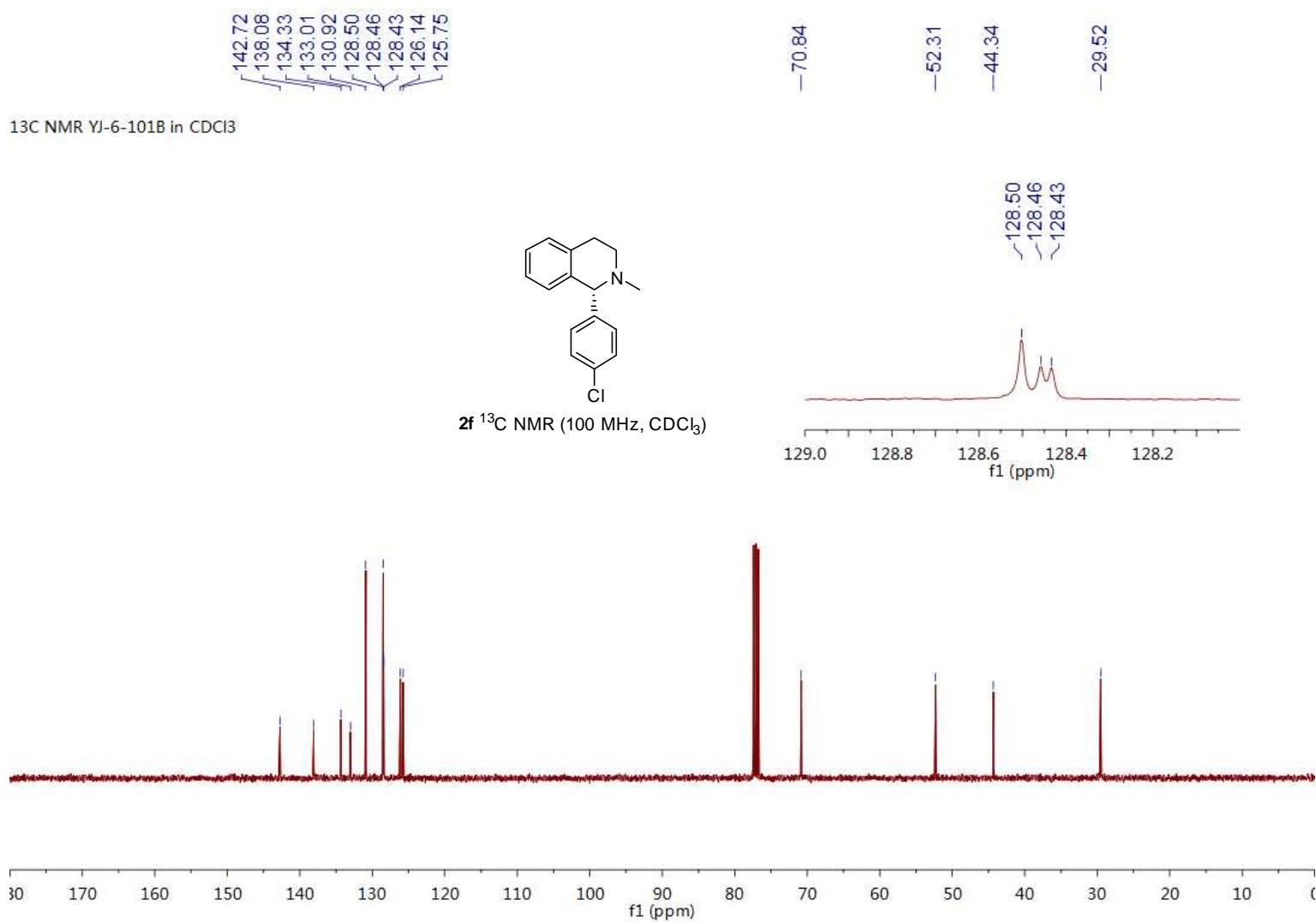
1H NMR YJ-6-101B in CDCl₃

-4.2148
3.2442
3.1264
3.1200
3.0979
3.0912
3.0843
2.8284
2.7880
2.6554
2.6462
2.6272
2.6179



2f ¹H NMR (400 MHz, CDCl₃)

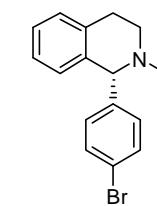




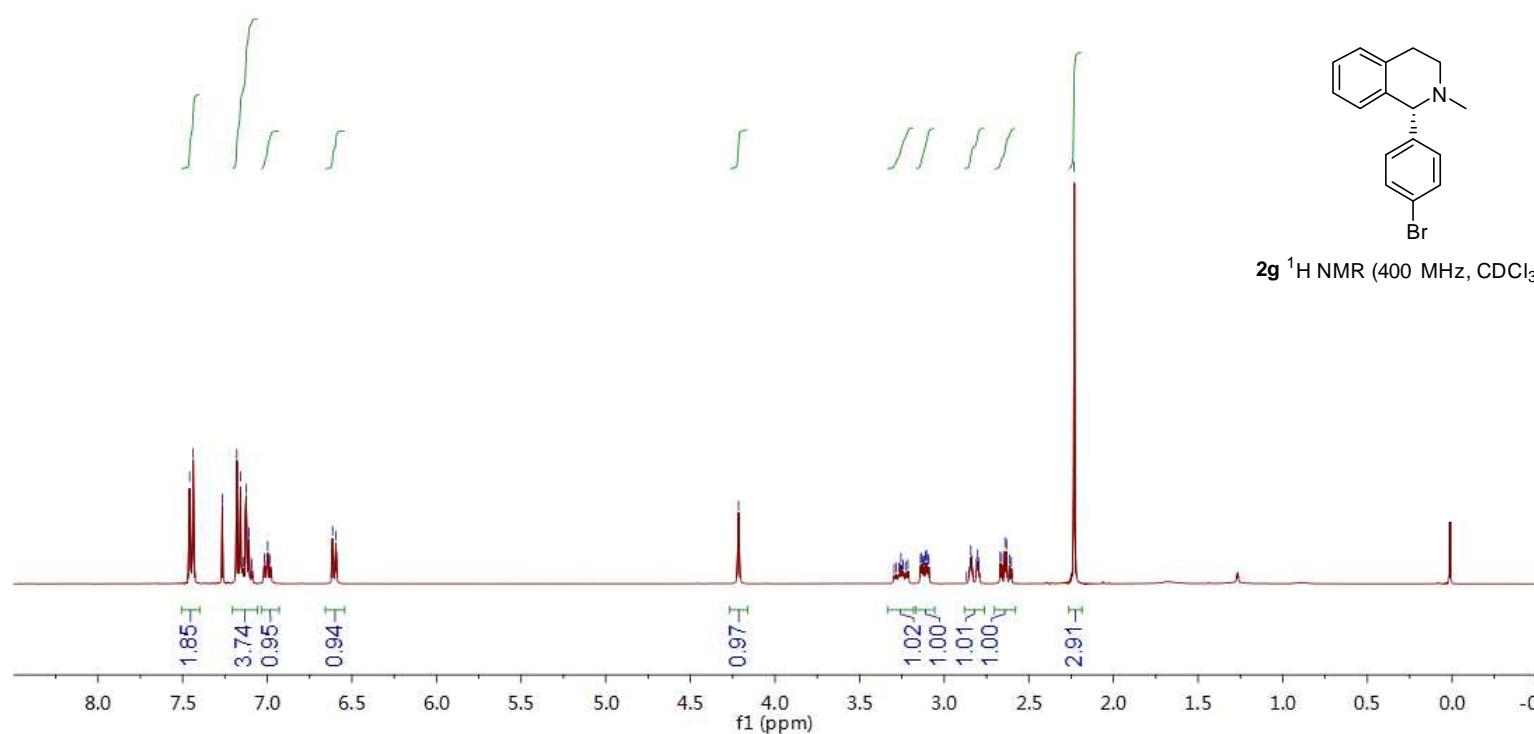
7.4586
7.4377
7.2657
7.1797
7.1589
7.1425
7.1255
7.1090
7.0173
6.9973
6.9818
6.6139
6.5944

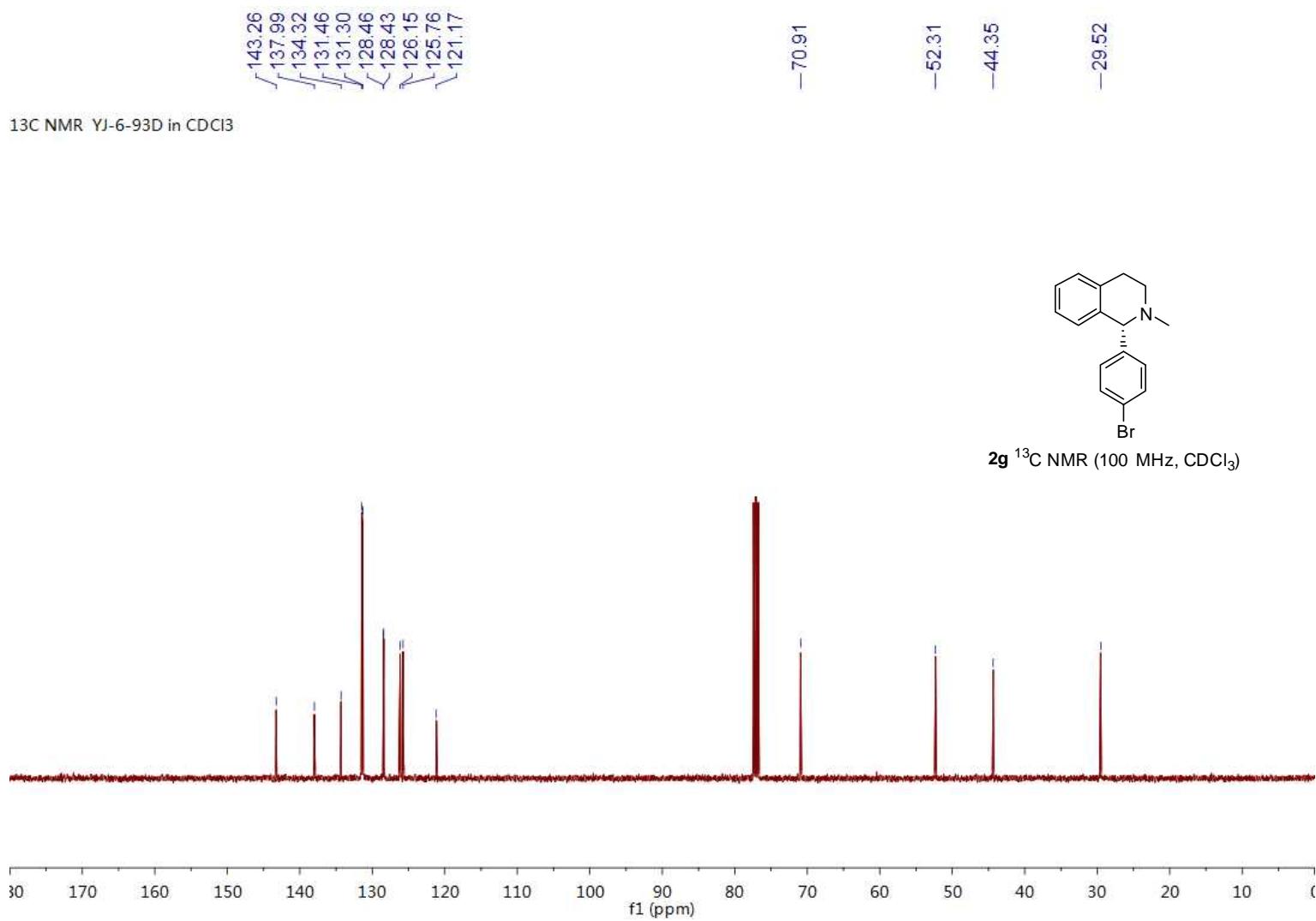
1H NMR YJ-6-93D in CDCl₃

-4.2155
3.2552
3.1331
3.1111
3.1045
3.0975
2.8426
2.8022
2.6692
2.6598
2.6409
2.2317



2g ¹H NMR (400 MHz, CDCl₃)

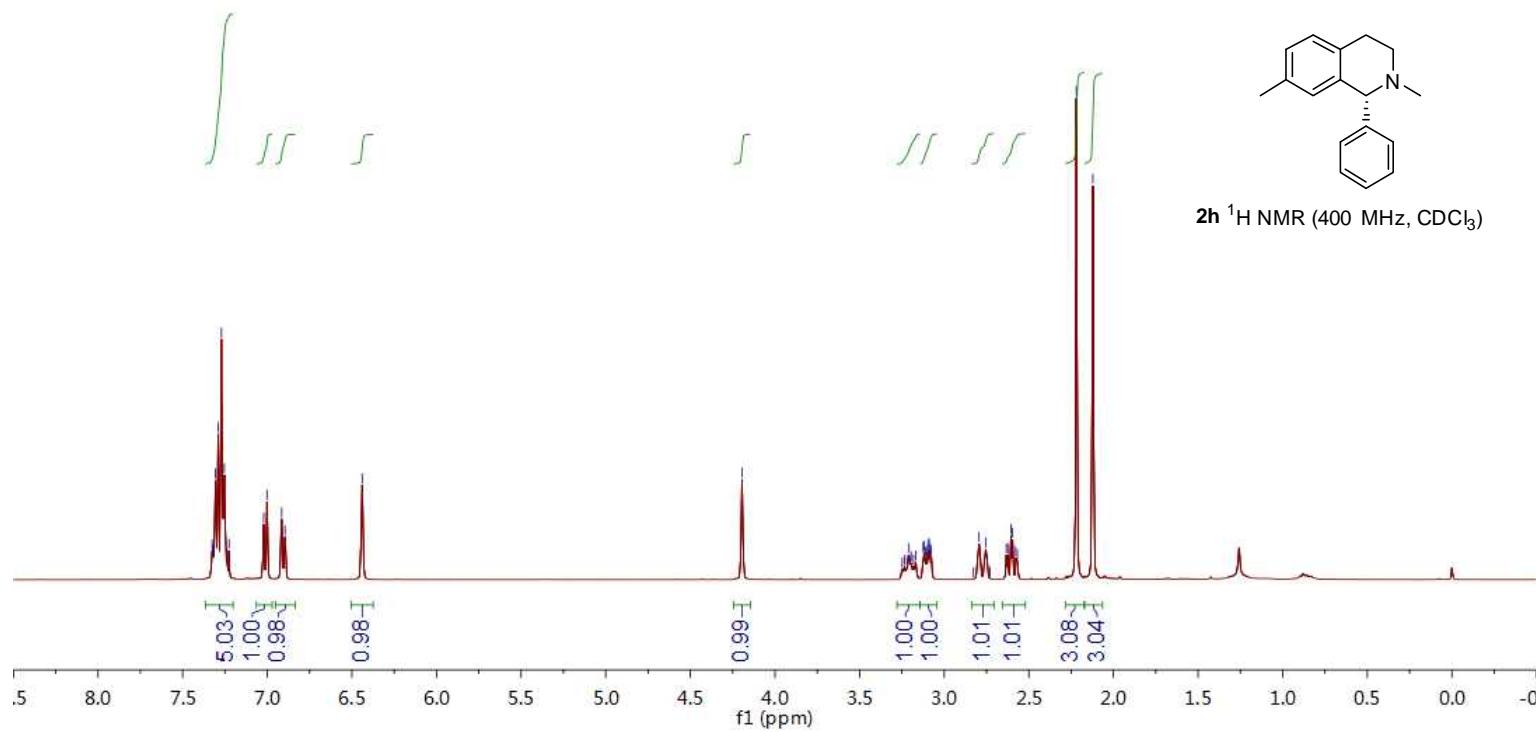


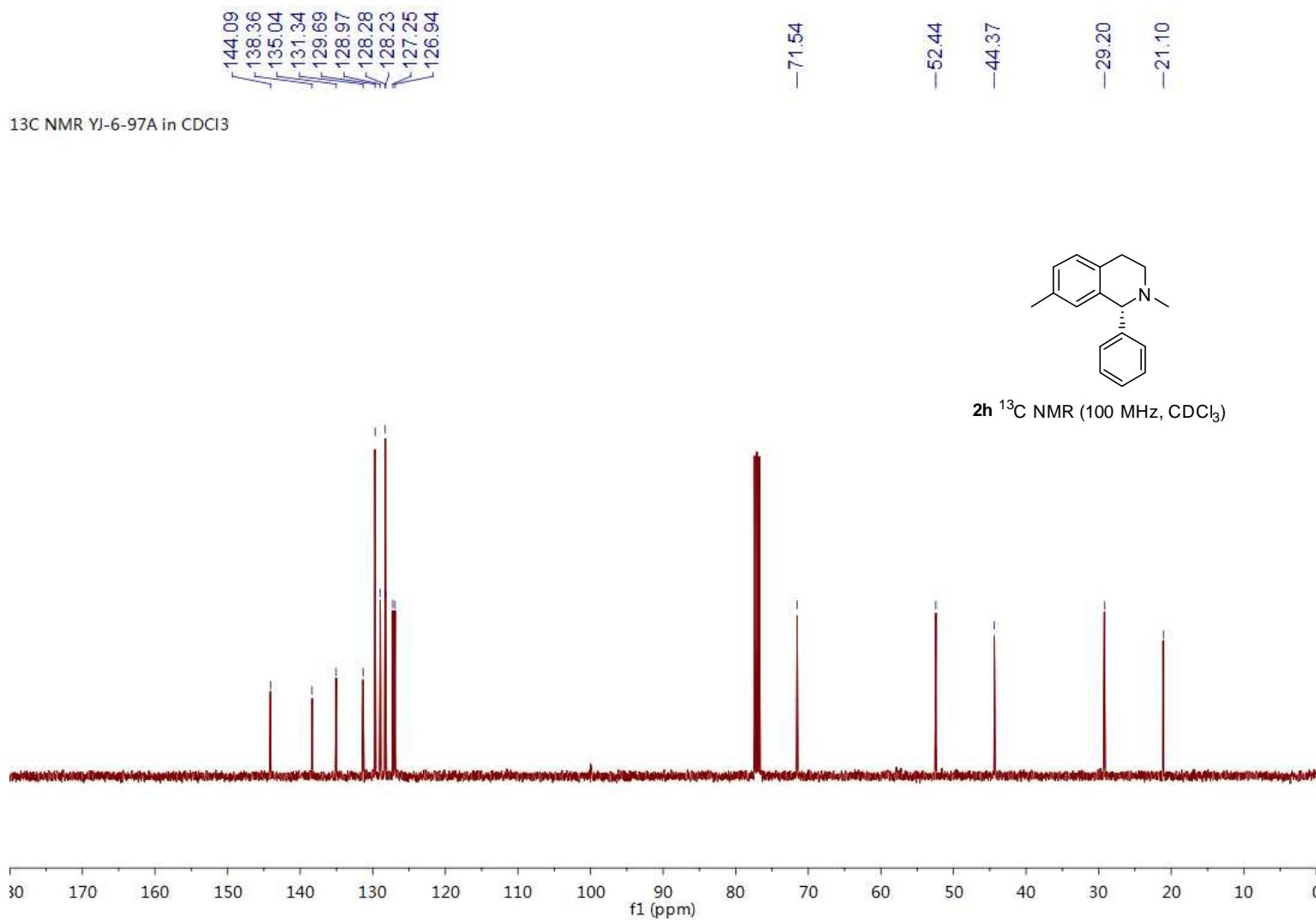


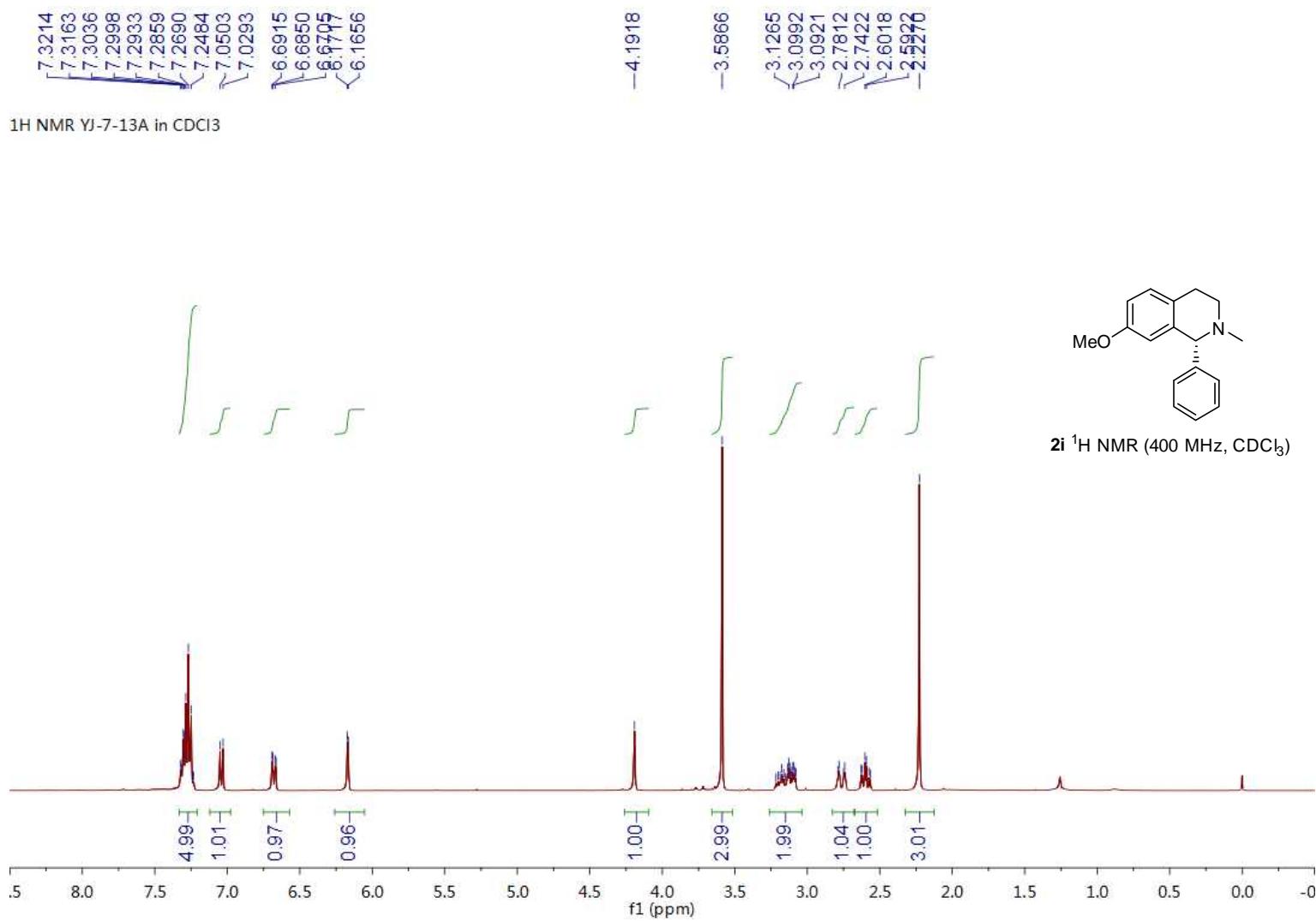
7.3257
7.3214
7.3178
7.3057
7.2903
7.2710
7.2536
7.2262
7.0212
7.0019
6.9147
6.8954

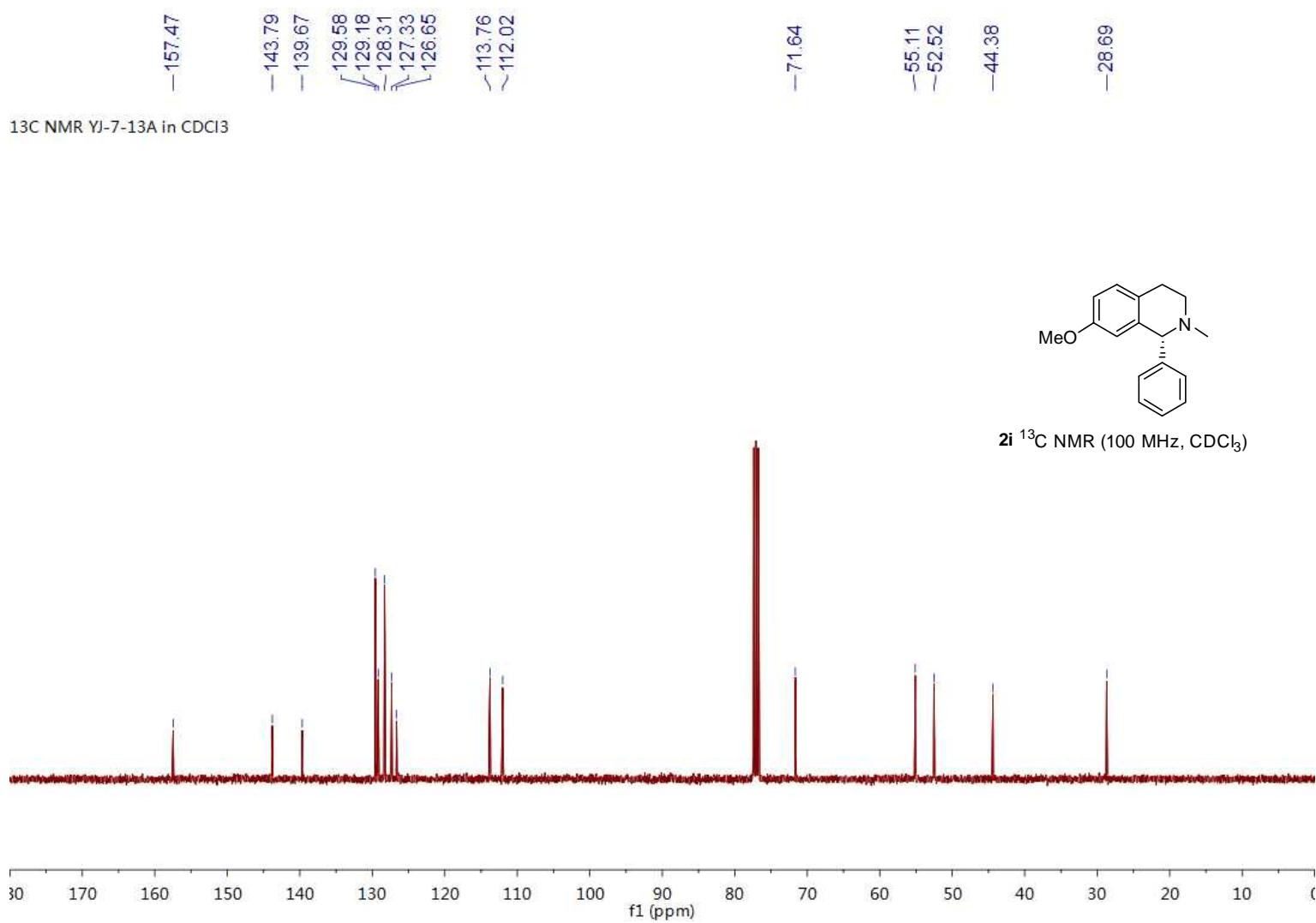
-4.1948
3.2078
3.1244
3.1177
3.0962
3.0890
3.0831
2.7940
2.7542
2.6331
2.6053
2.5958
2.2128
-2.1218

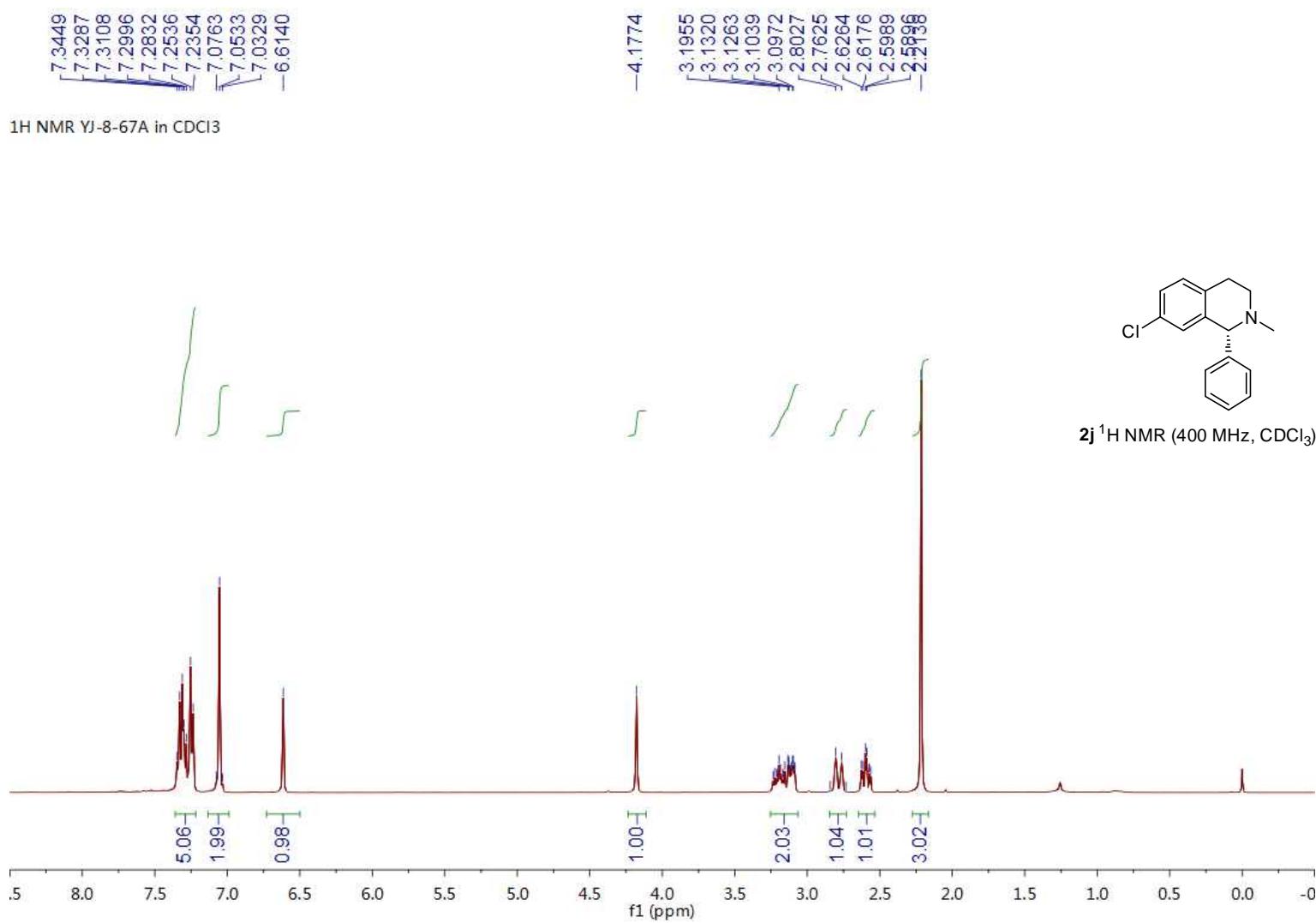
1H NMR YJ-6-97A in CDCl₃

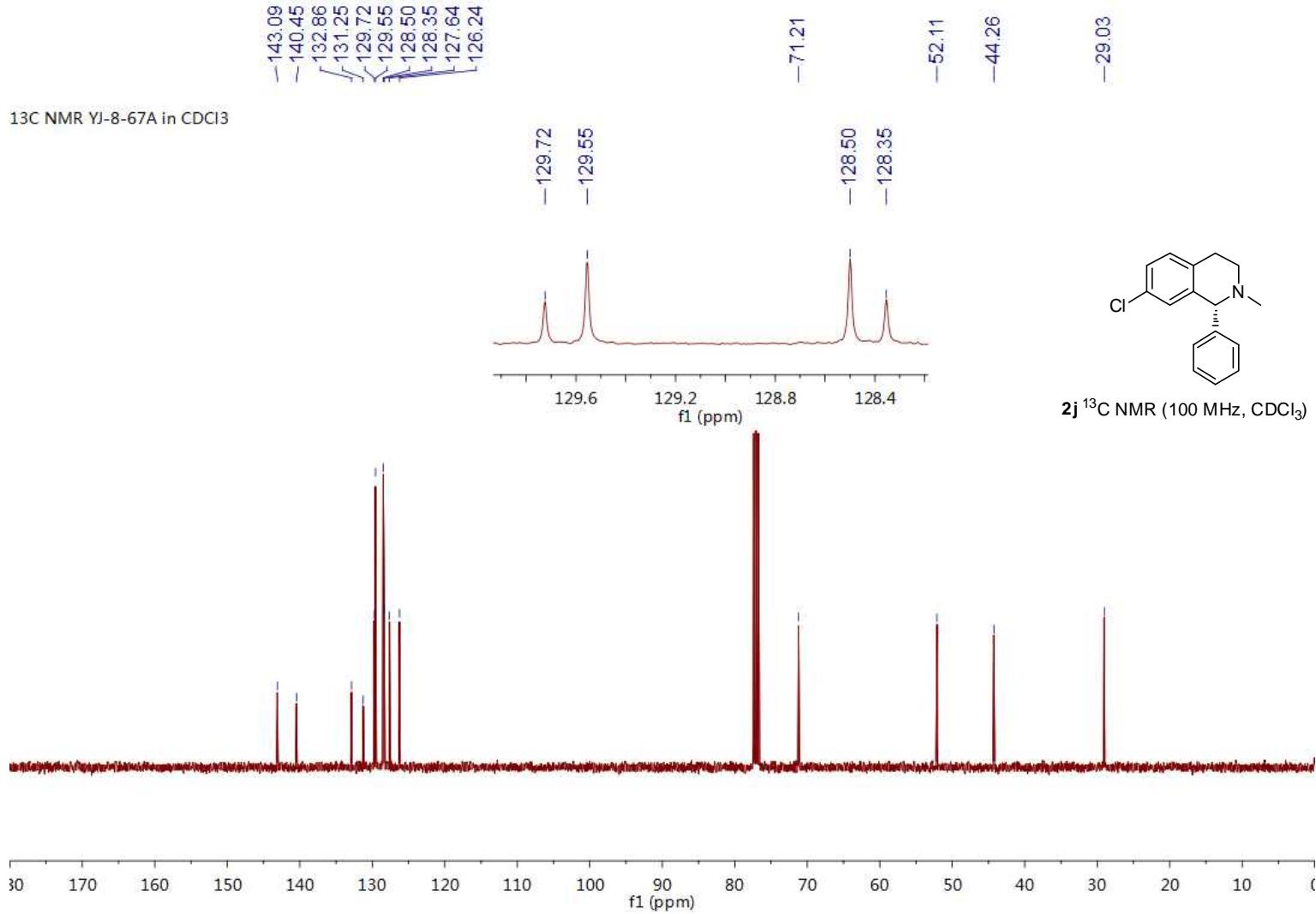


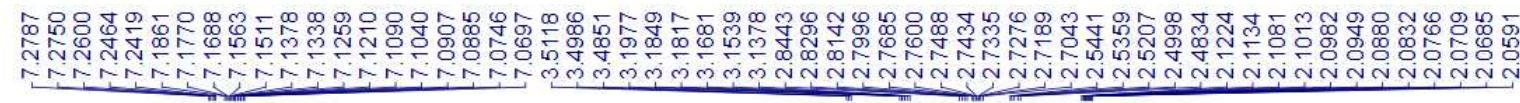




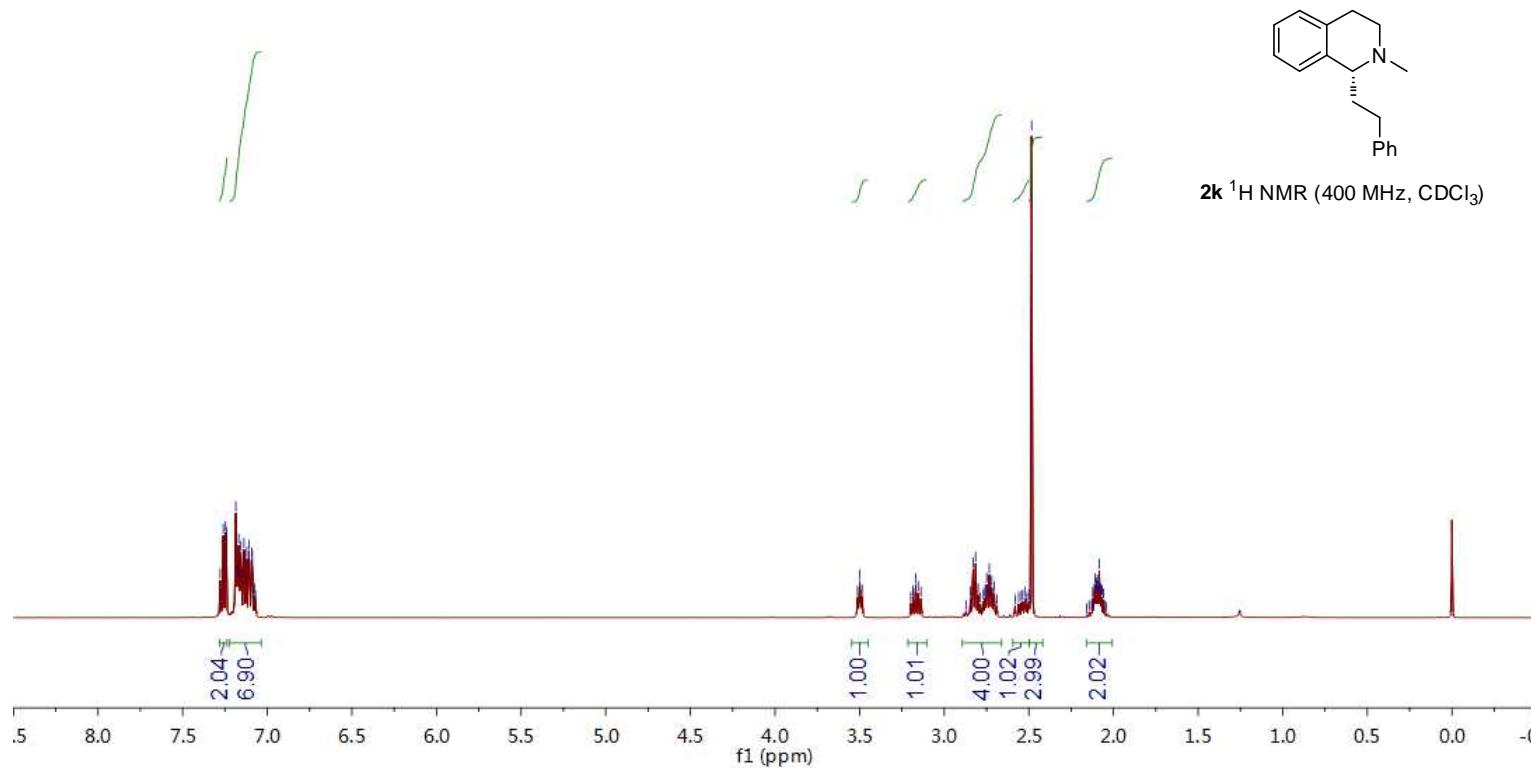


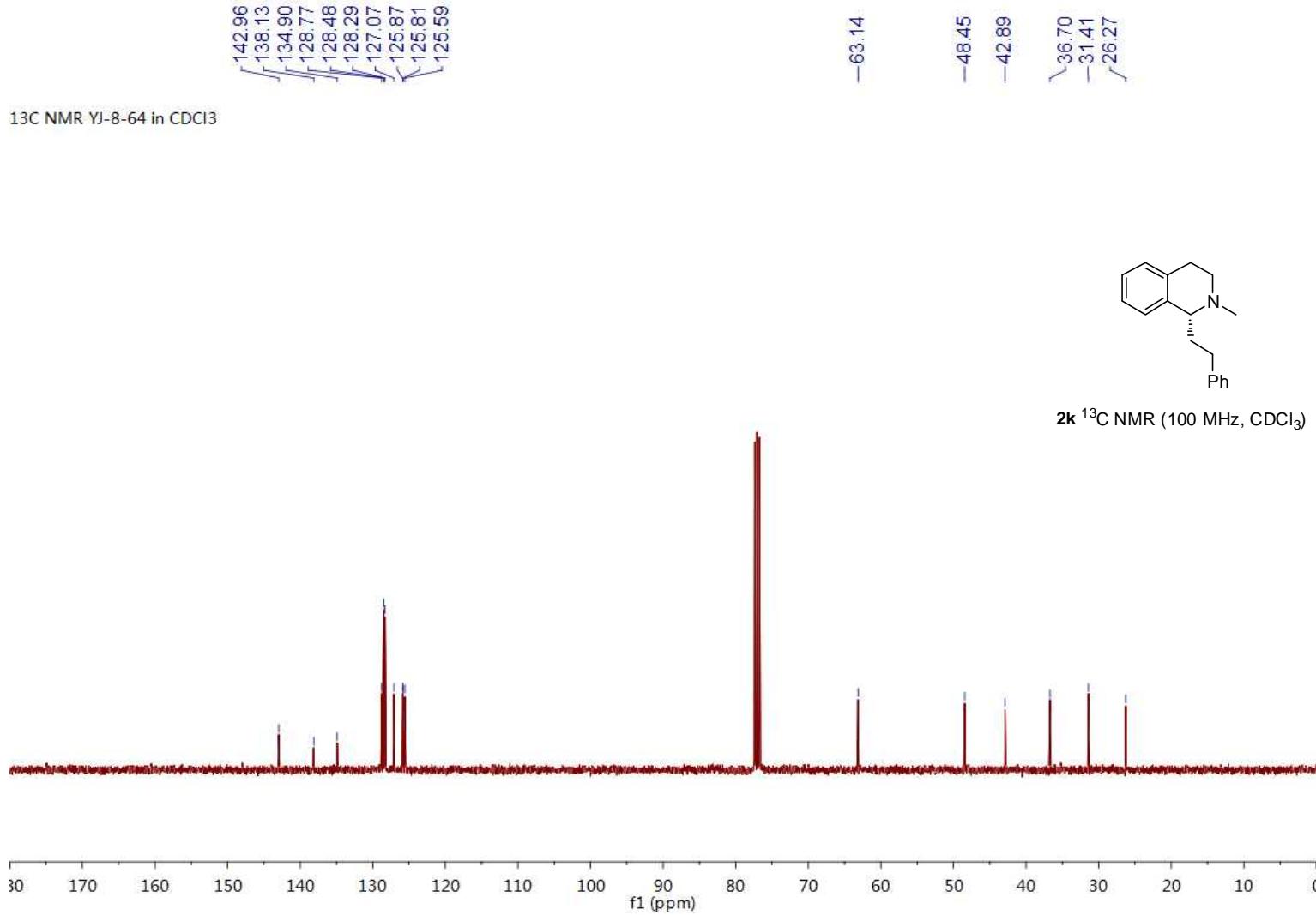






^1H NMR YJ-8-64 in CDCl_3



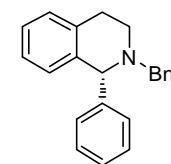


7.3897
 7.3713
 7.3363
 7.3182
 7.2978
 7.2803
 7.2620
 7.2495
 7.2369
 7.2309
 7.2151
 7.1989
 7.1205
 7.1061
 7.0891
 7.0699
 7.0177
 6.9975
 6.9823
 6.9771
 6.7333
 6.7139

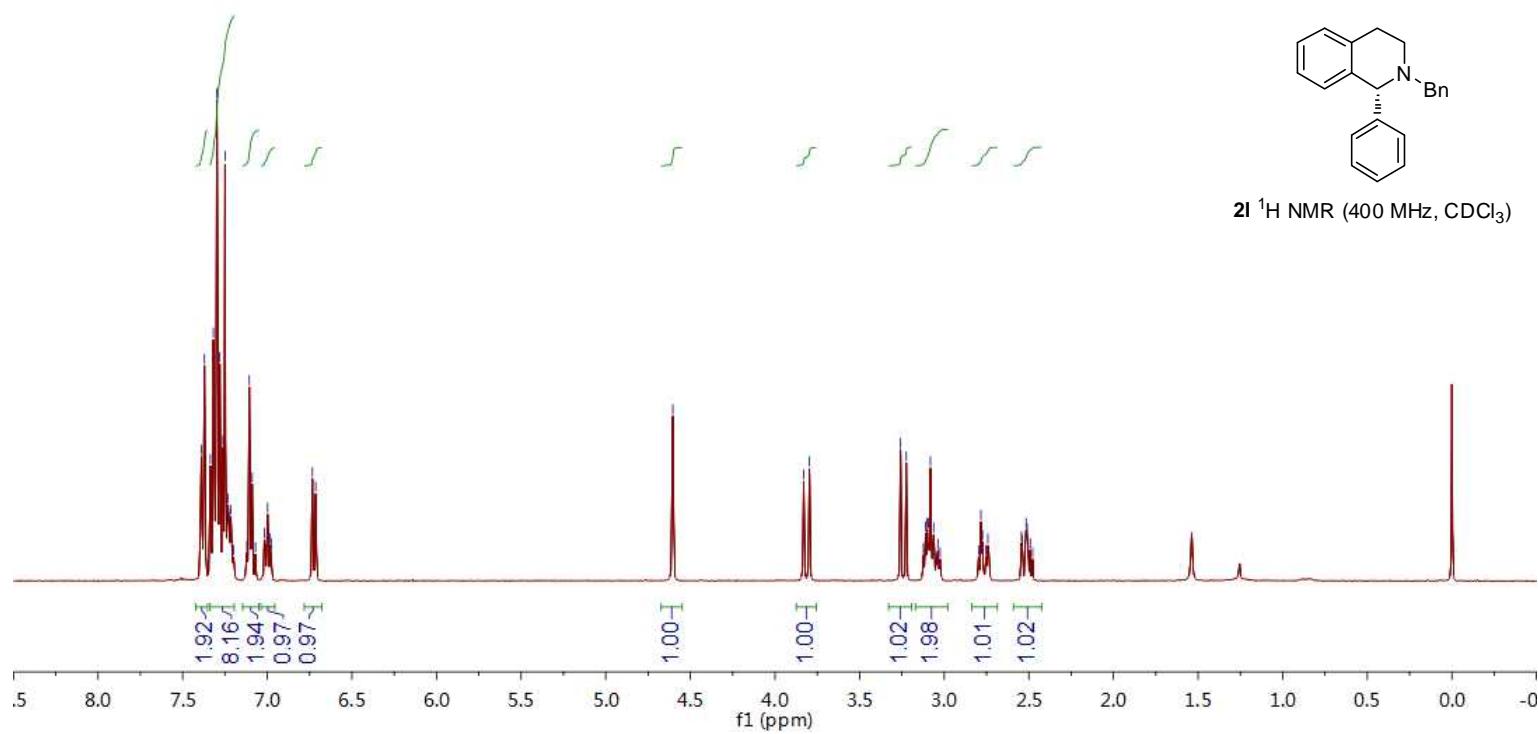
1H NMR YJ-6-67B in CDCl₃

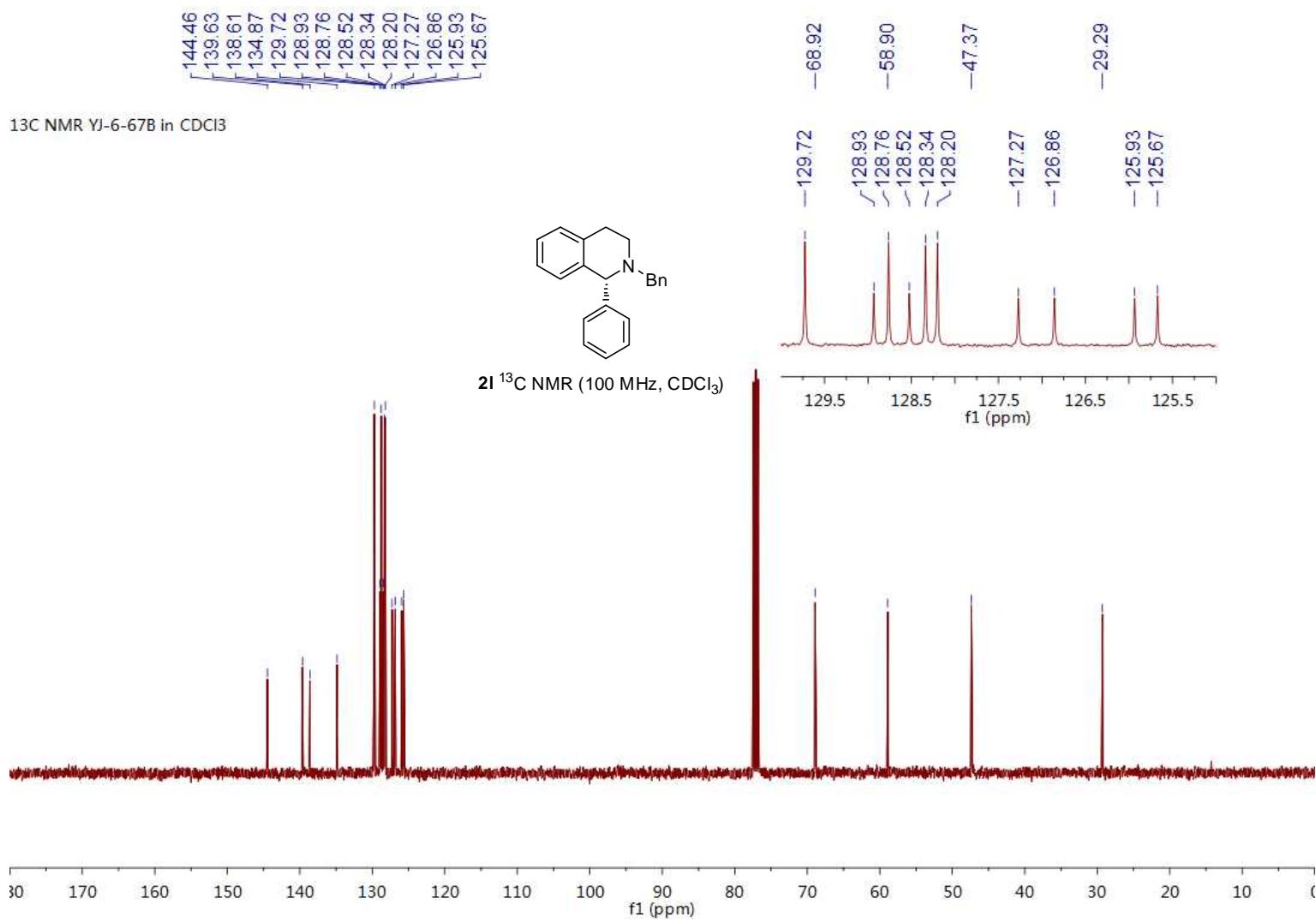
-4.6038

3.8302
 3.7963
 3.2581
 3.2242
 3.1227
 3.1098
 3.1004
 3.0937
 3.0814
 3.0623
 3.0473
 3.0374
 3.0242
 2.7970
 2.7833
 2.7740
 2.7528
 2.7420
 2.5436
 2.5385
 2.5161
 2.5097
 2.4917
 2.4776



2l ¹H NMR (400 MHz, CDCl₃)



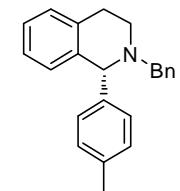


7.3086
 7.2917
 7.2827
 7.2637
 7.2424
 7.2140
 7.1969
 7.1826
 7.1239
 7.1043
 7.0853
 7.0690
 7.0496
 7.0018
 6.9963
 6.9665
 6.9611
 6.9818
 6.7386
 6.7192

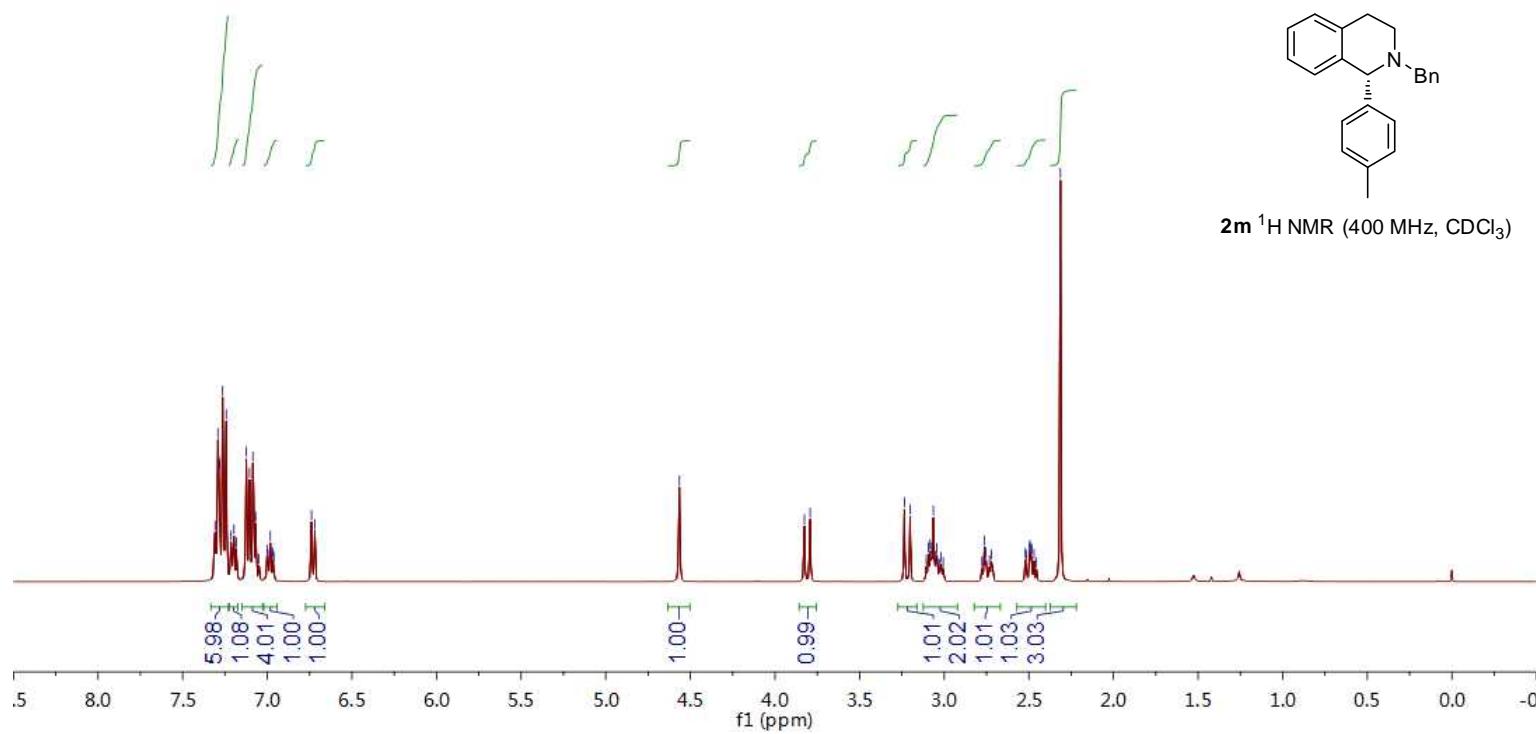
¹H NMR YJ-7-4A in CDCl₃

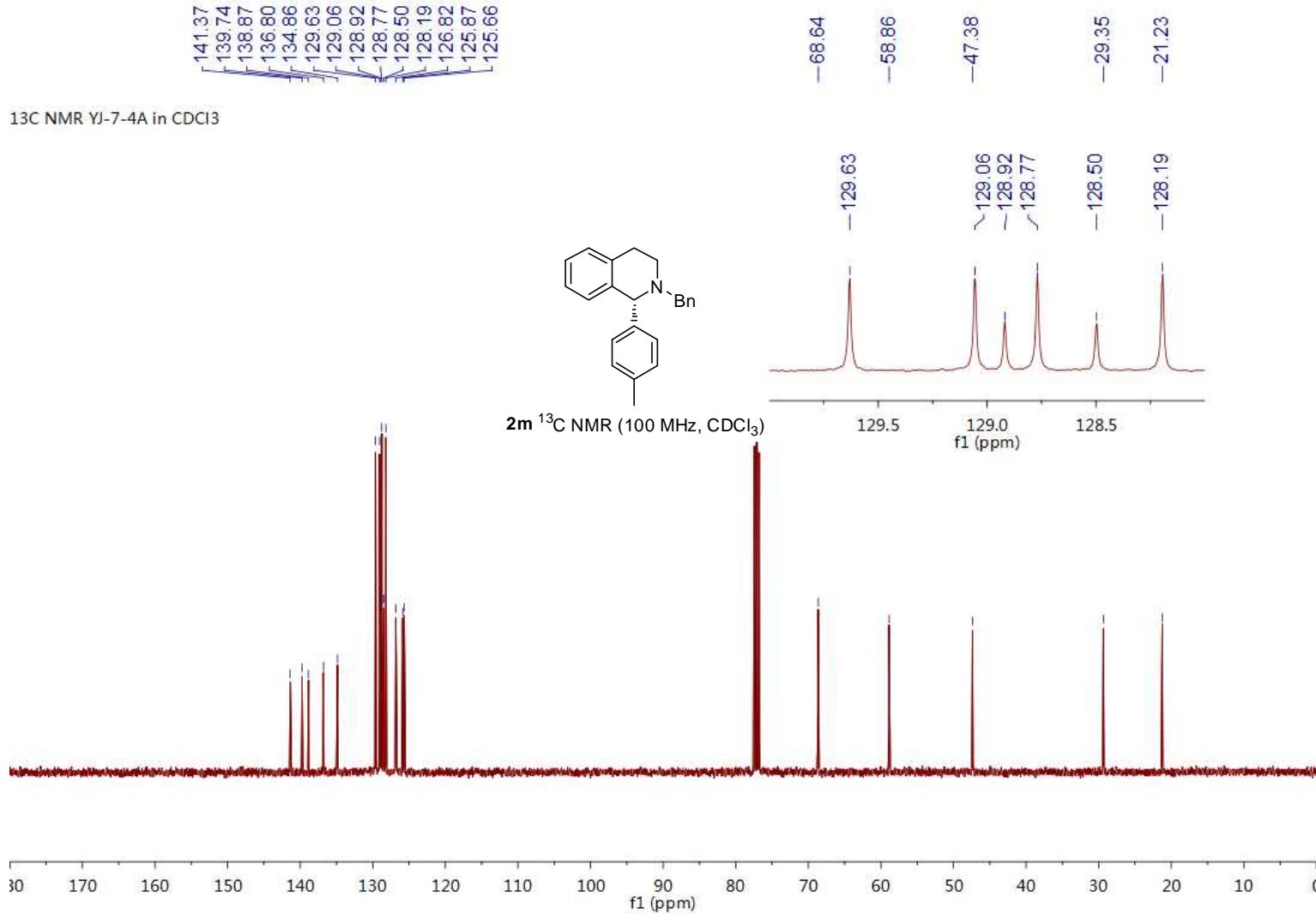
-4.5649

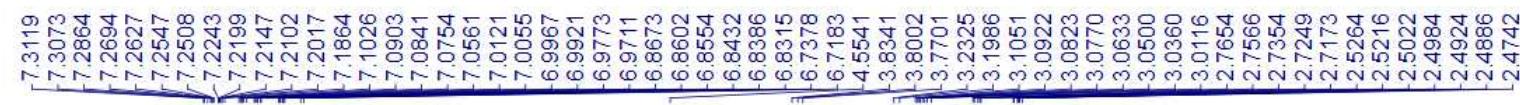
3.8261
 3.7922
 3.2354
 3.2014
 3.1055
 3.0926
 3.0831
 3.0744
 3.0641
 3.0504
 3.0435
 3.0291
 3.0191
 3.0060
 2.7765
 2.7627
 2.7538
 2.7324
 2.7214
 2.5205
 2.5153
 2.4960
 2.4929
 2.4867
 2.4832
 2.4688
 2.4547
 2.3131



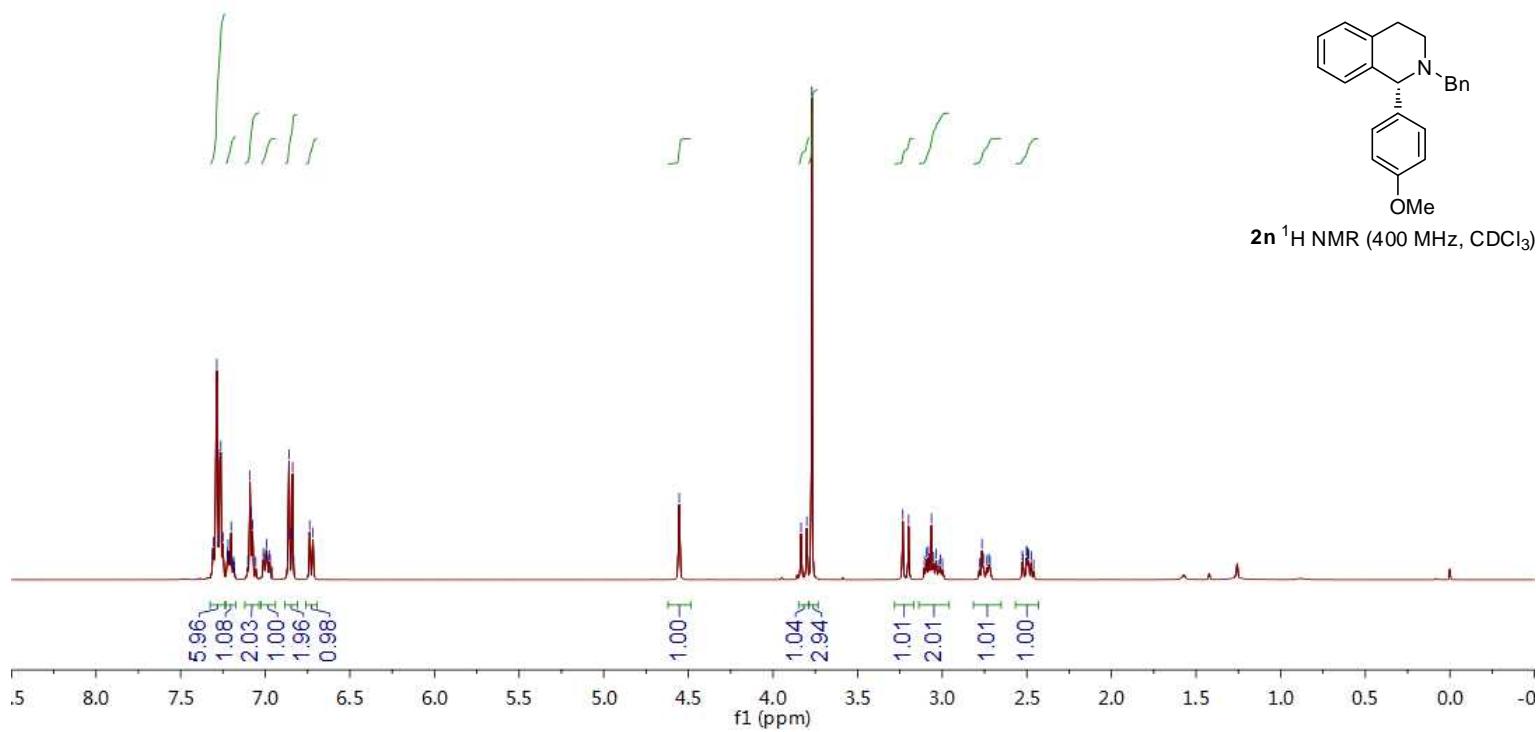
2m ¹H NMR (400 MHz, CDCl₃)

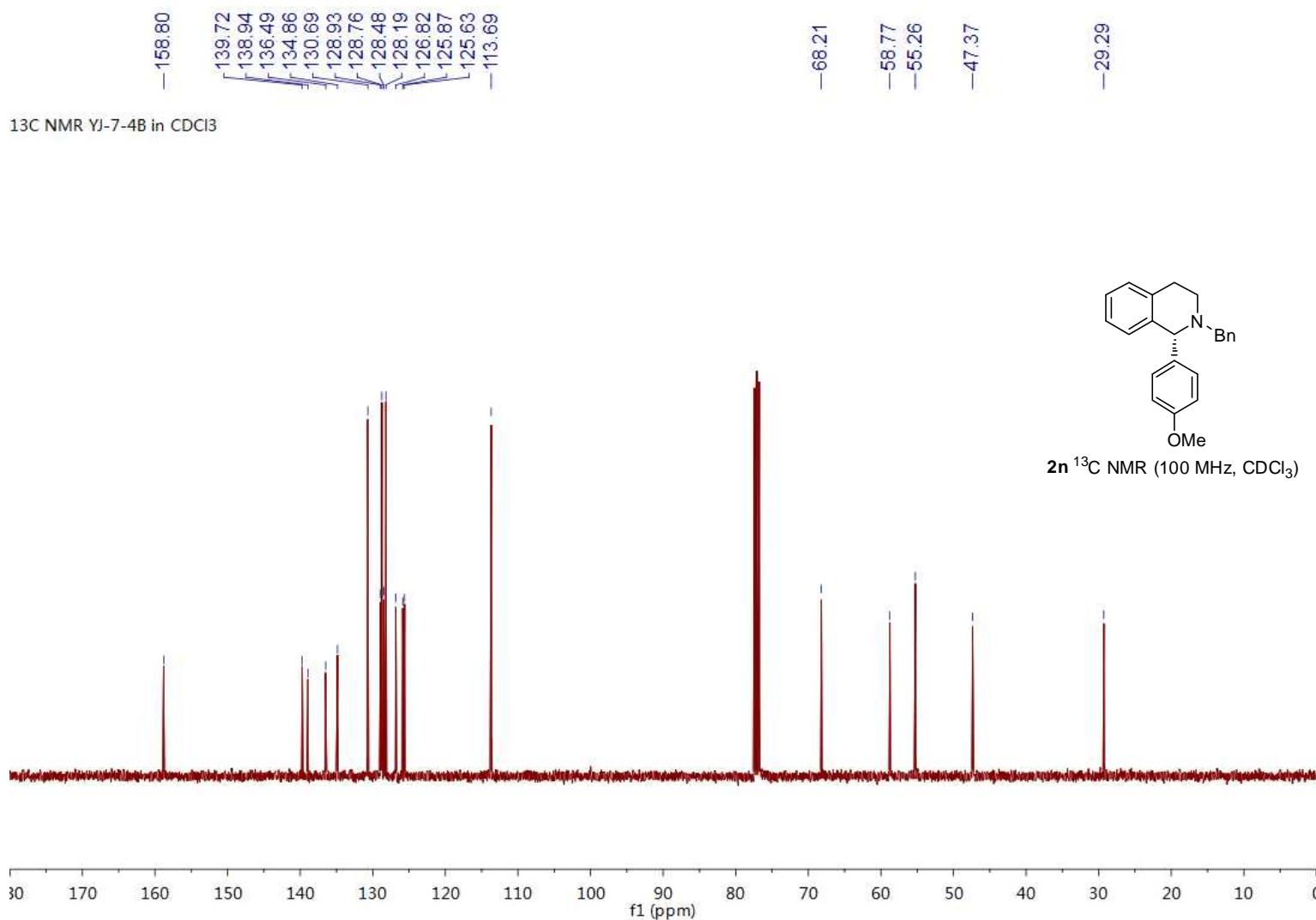


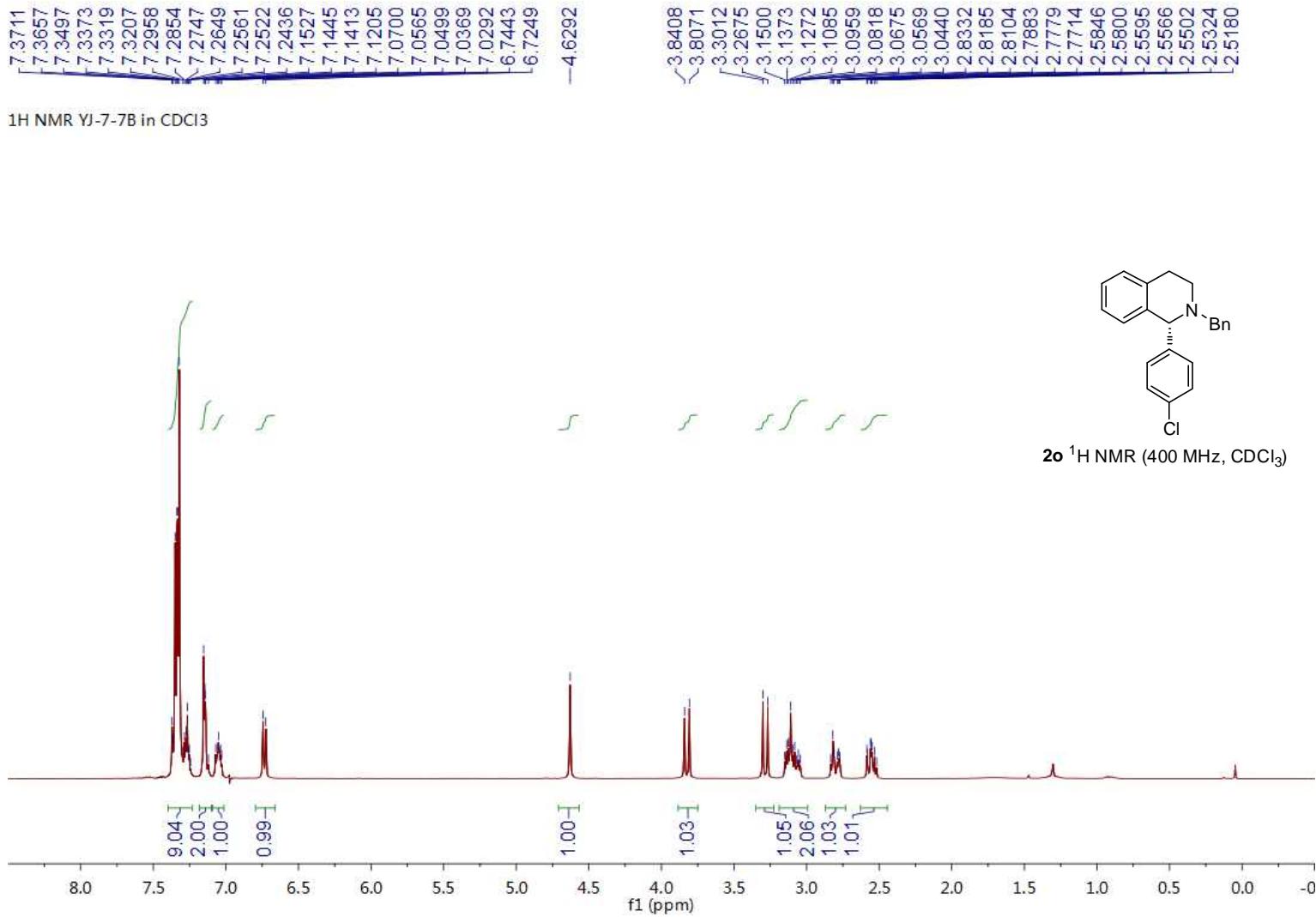


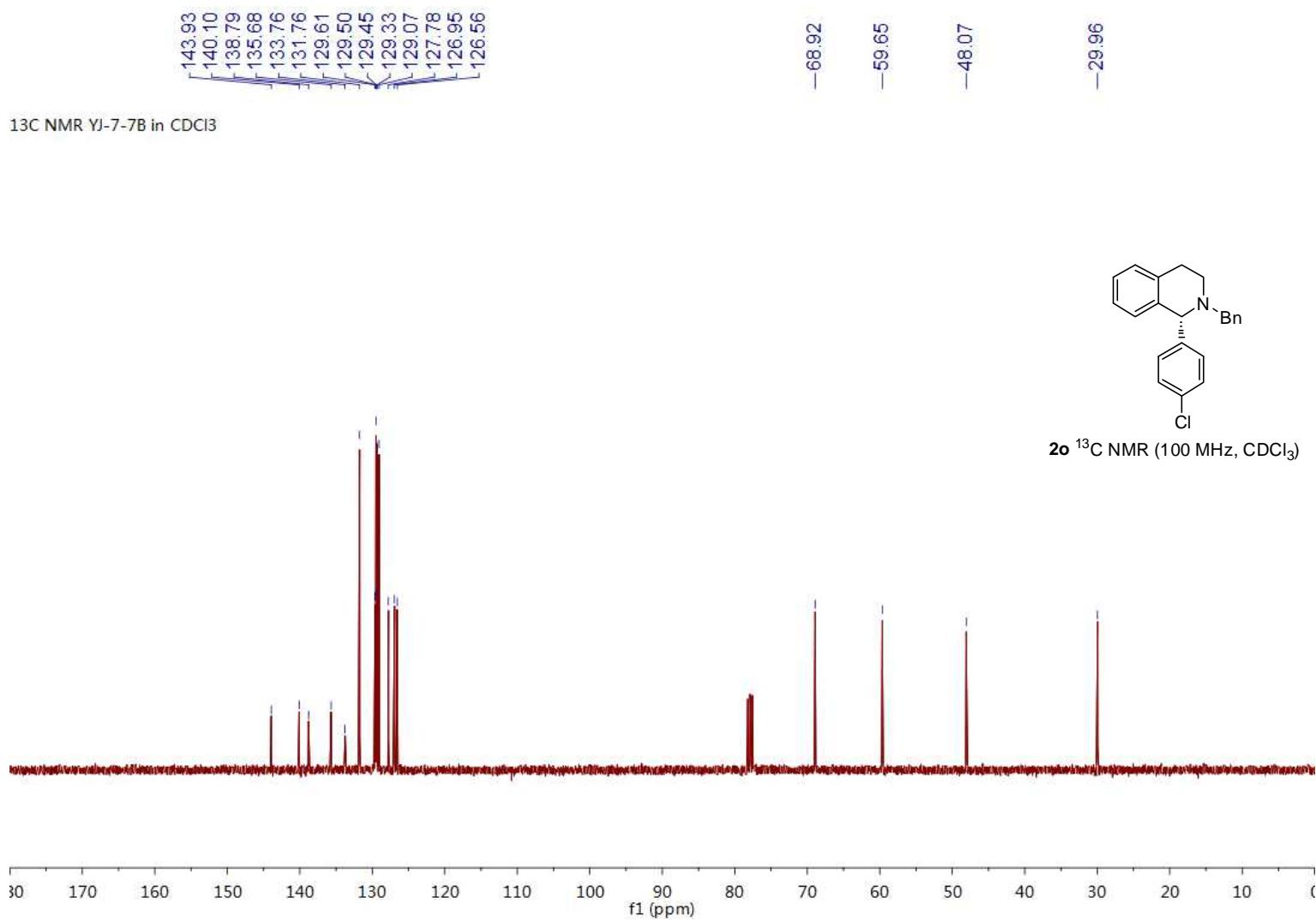


¹H NMR YJ-7-4B in CDCl₃



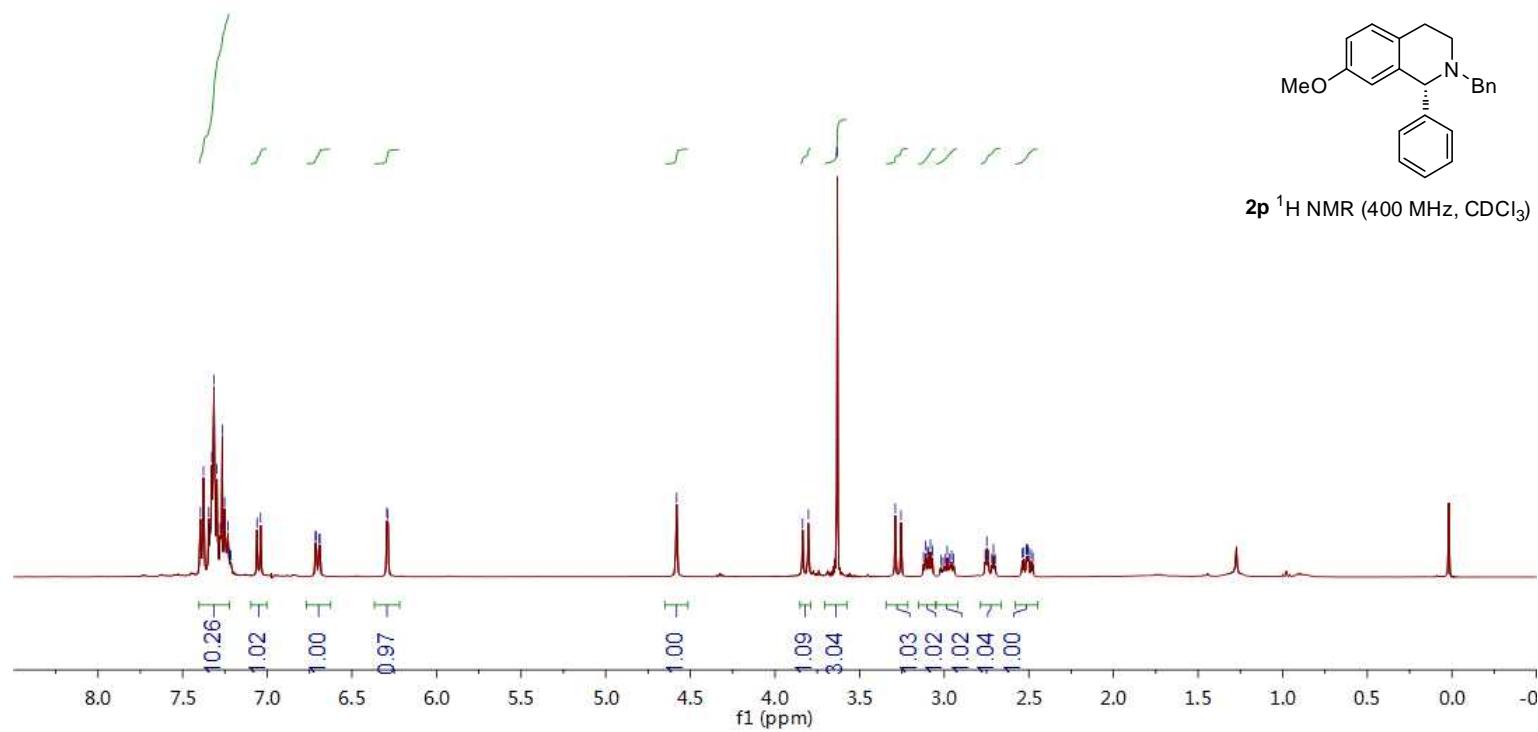


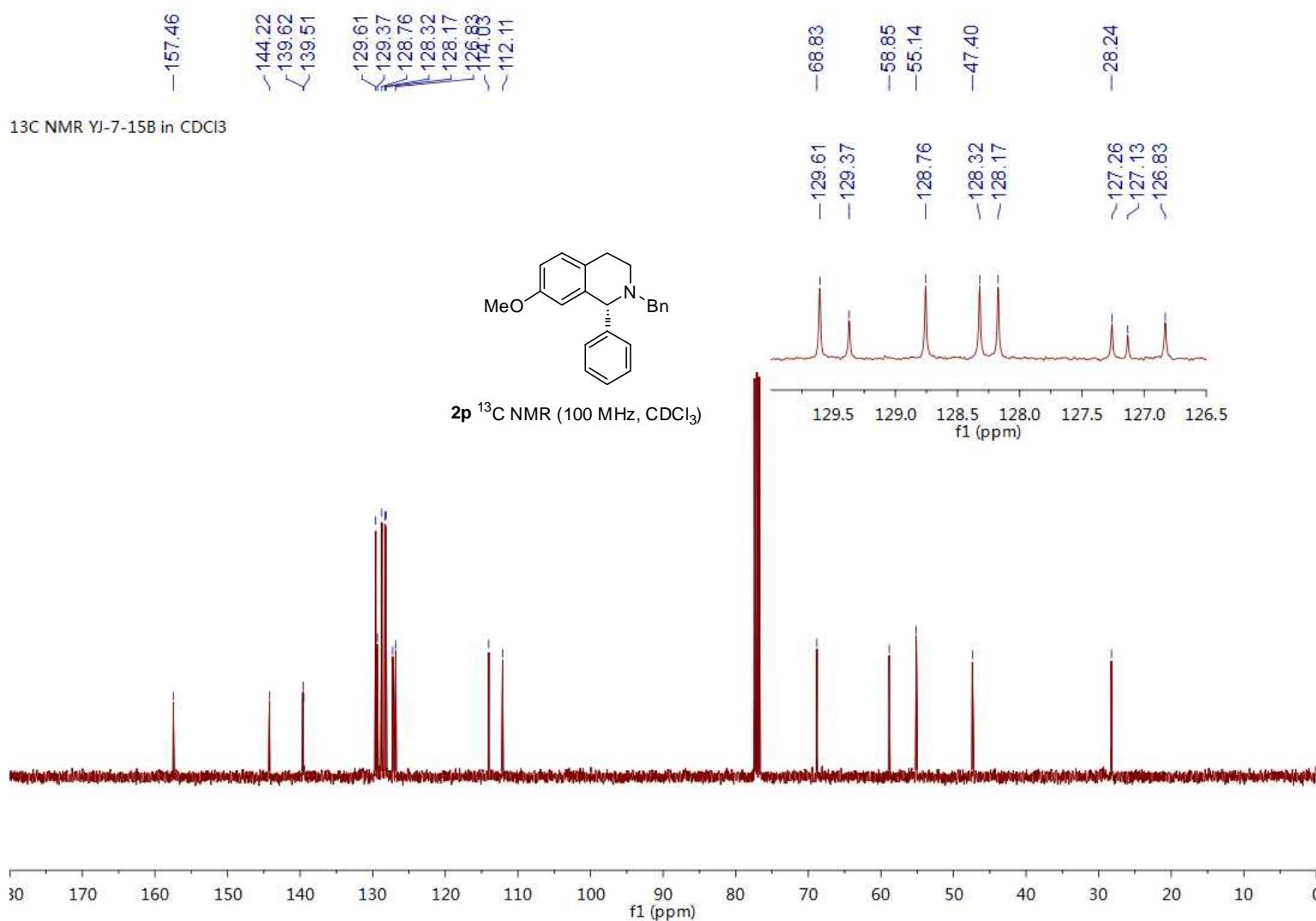


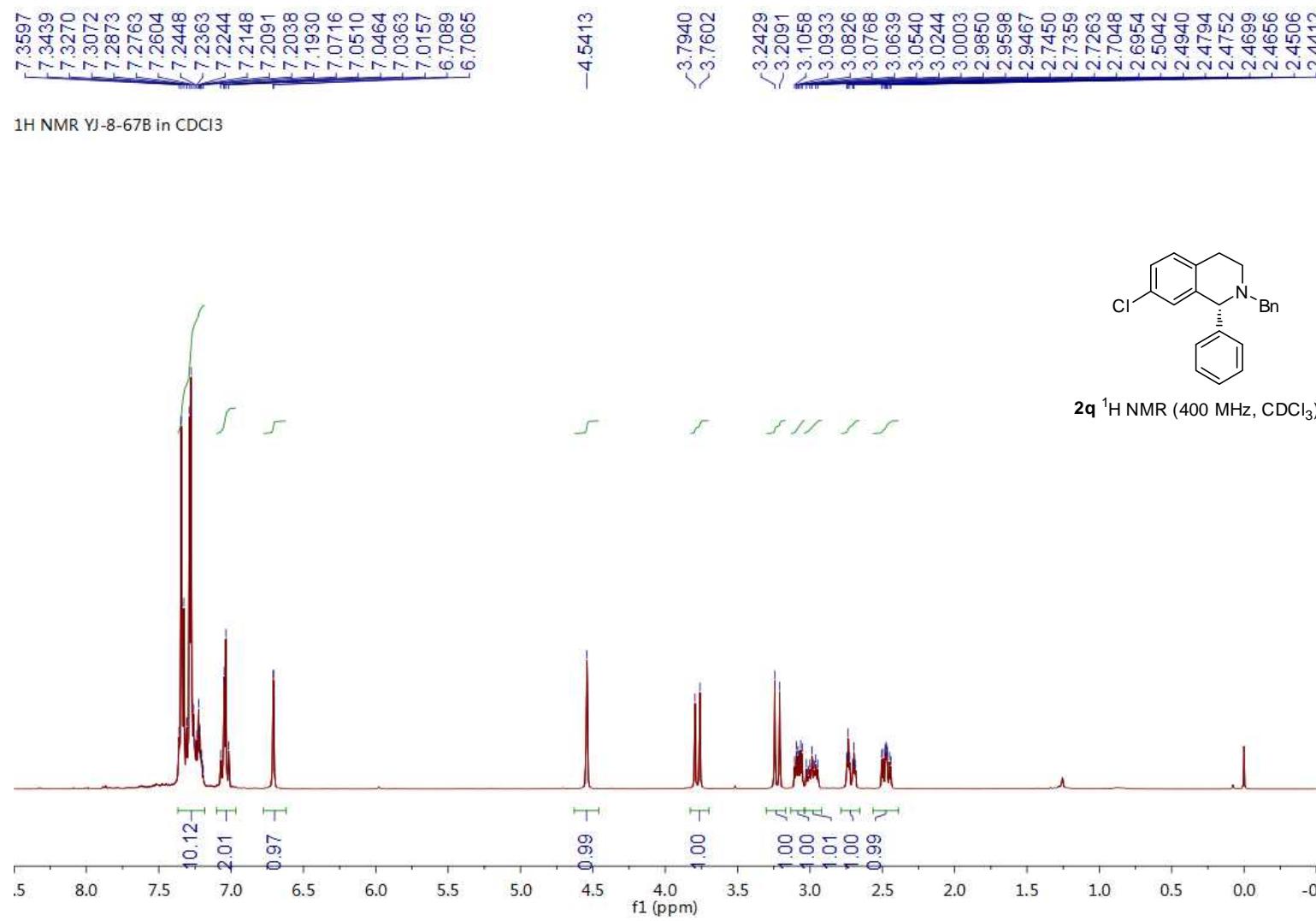


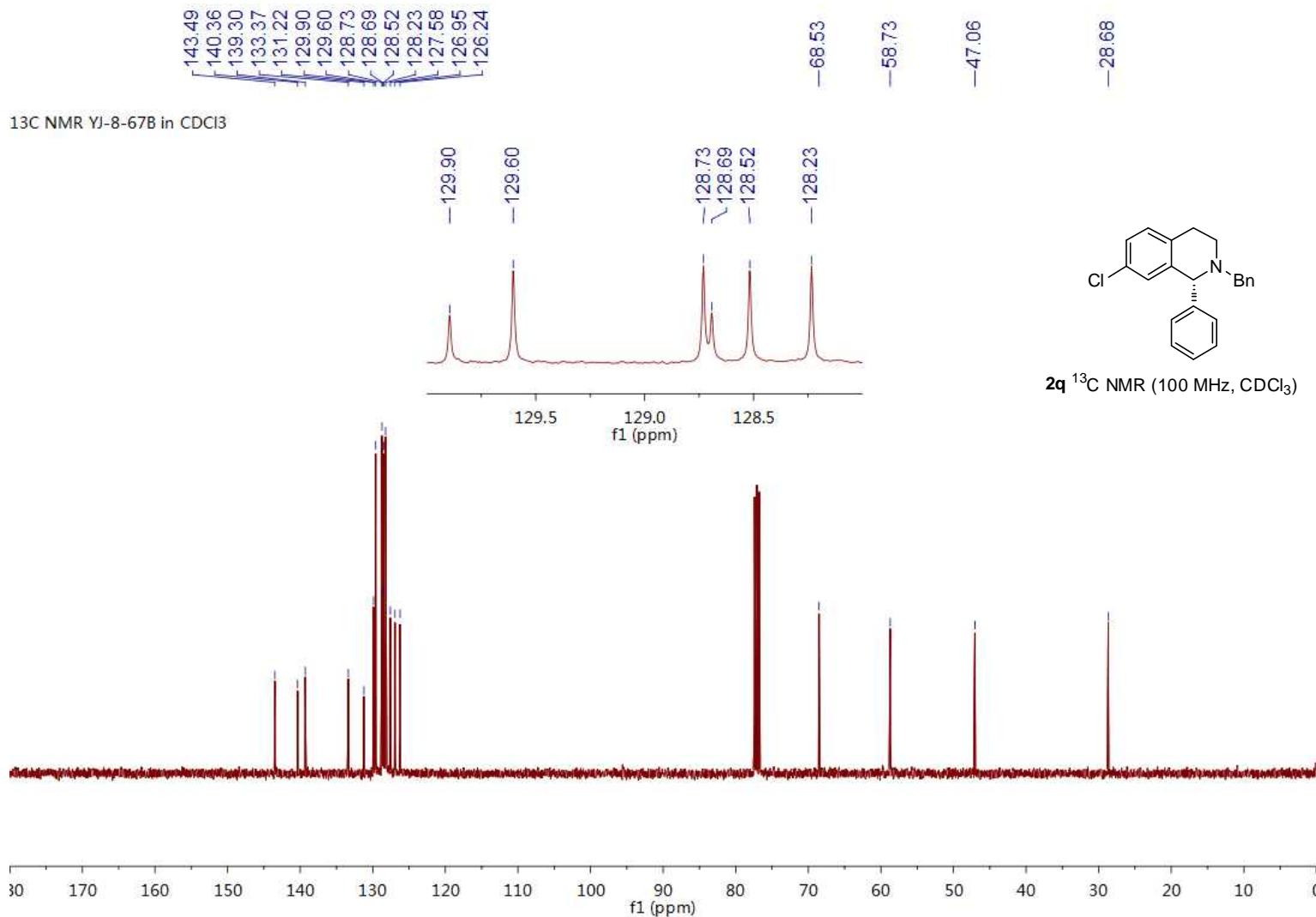


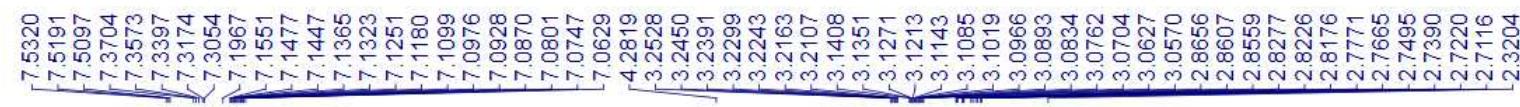
1H NMR YJ-7-15B in CDCl₃



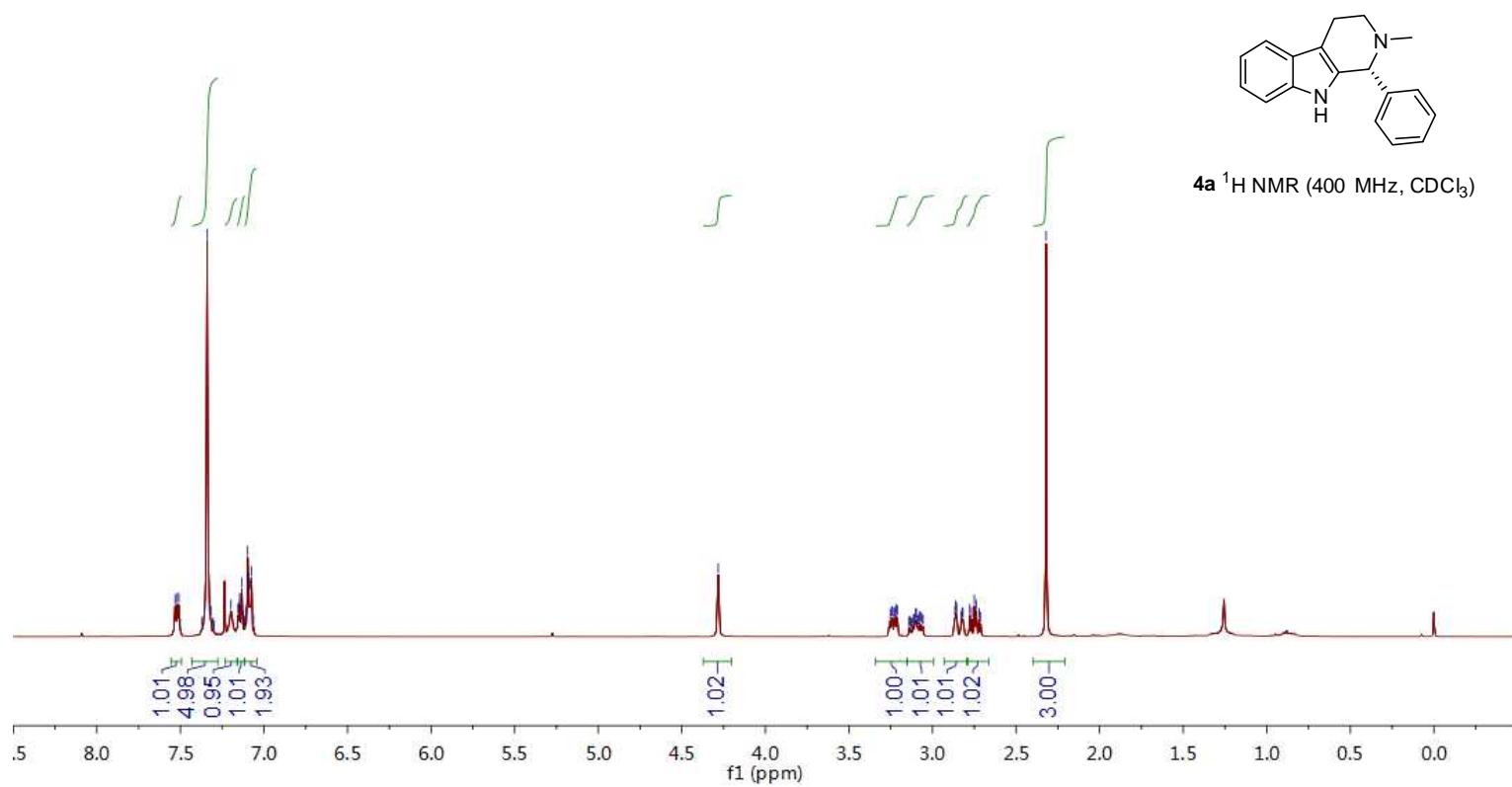


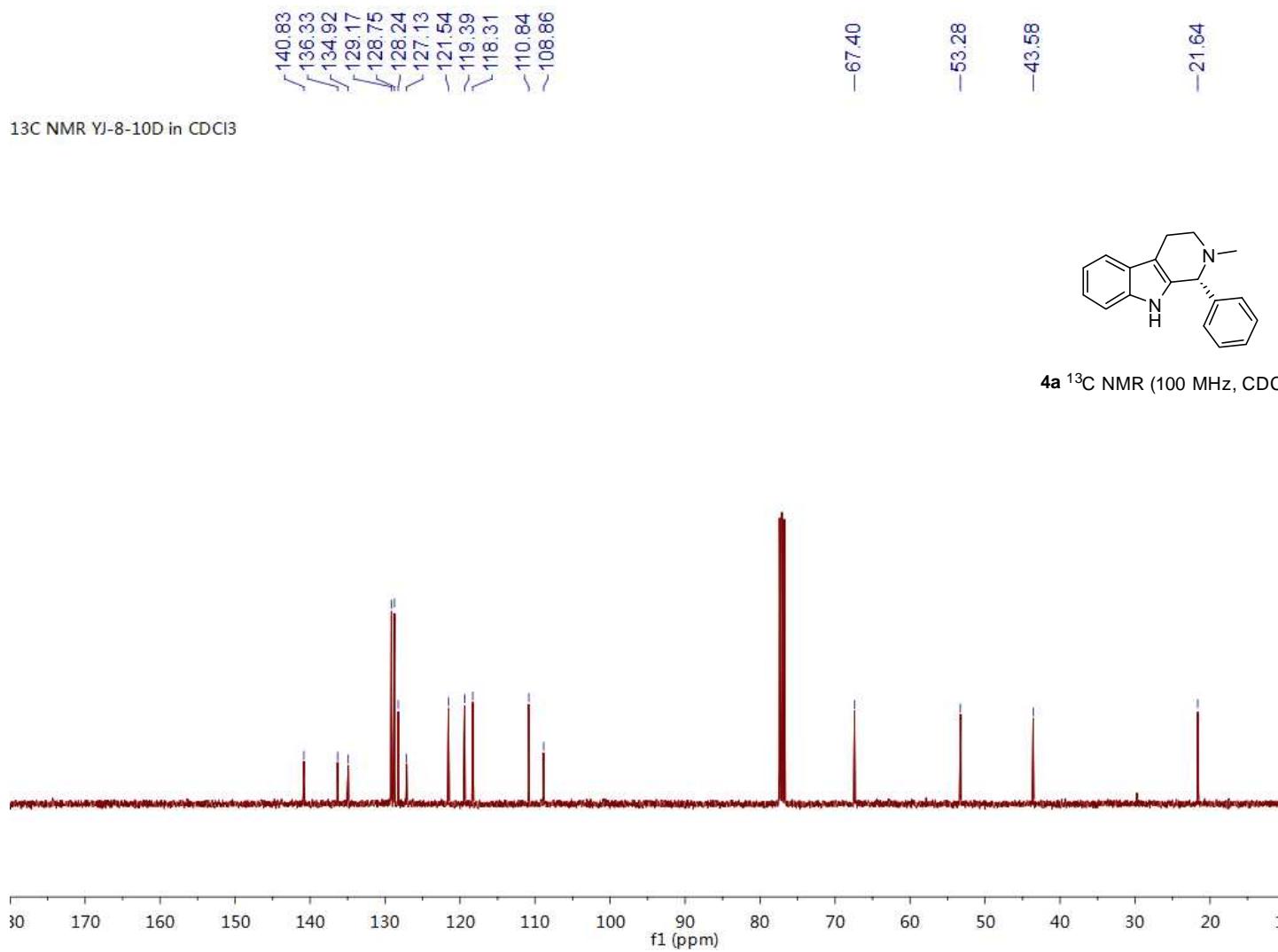


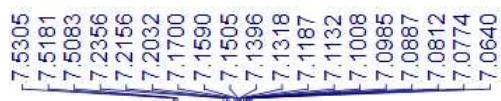




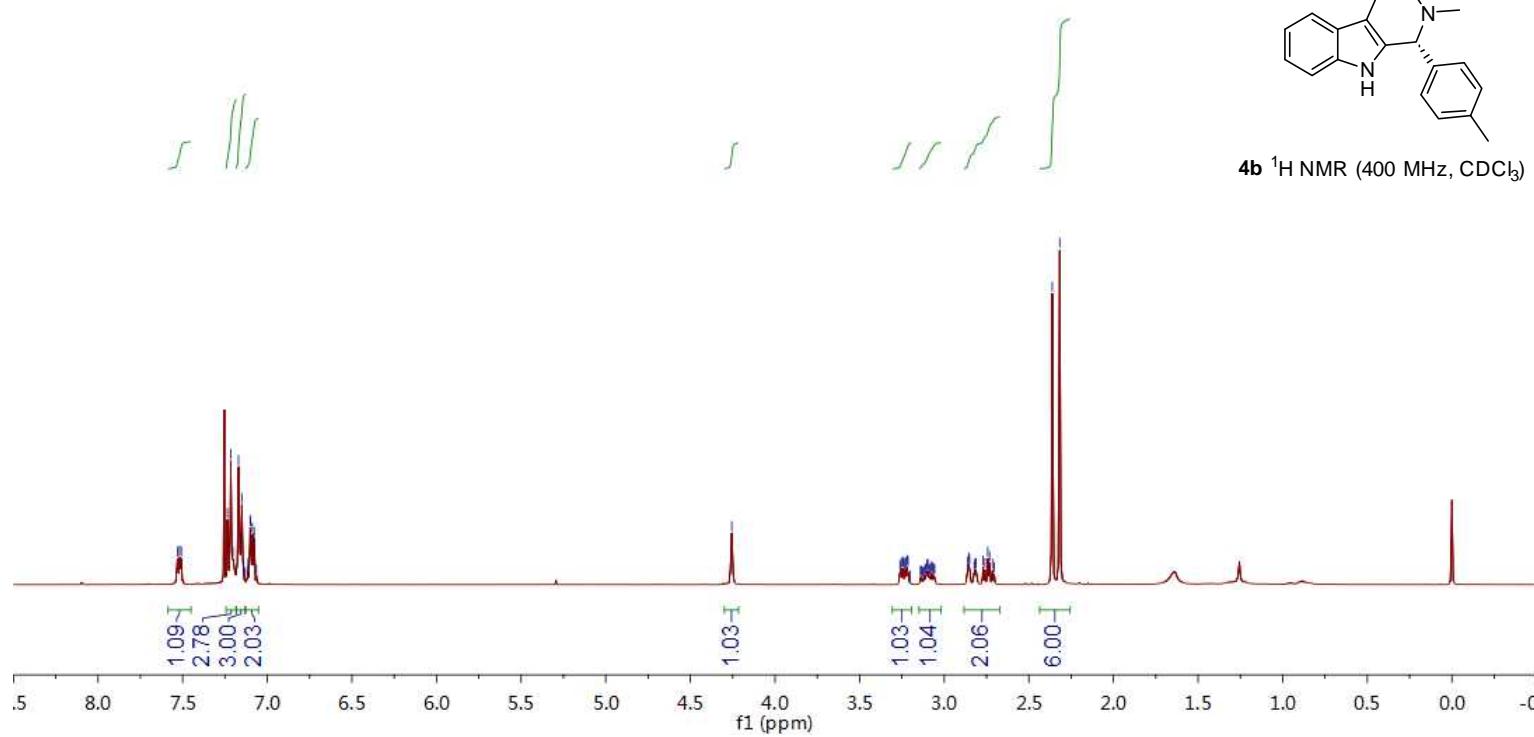
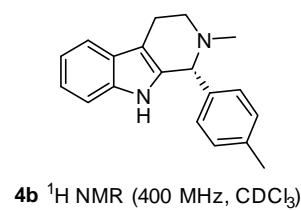
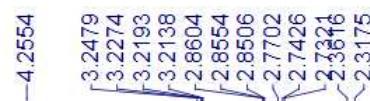
¹H NMR YJ-8-10D in CDCl₃

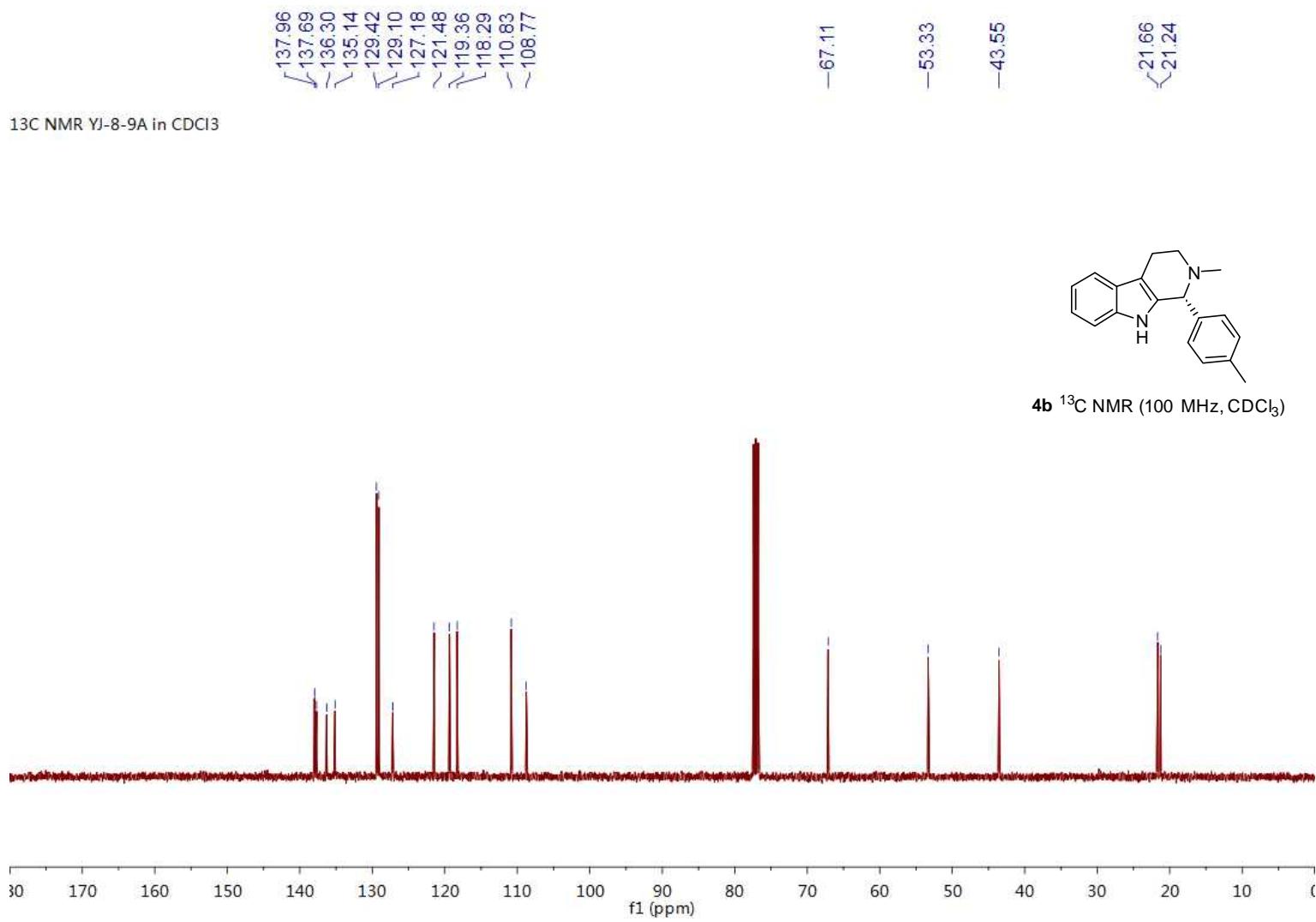


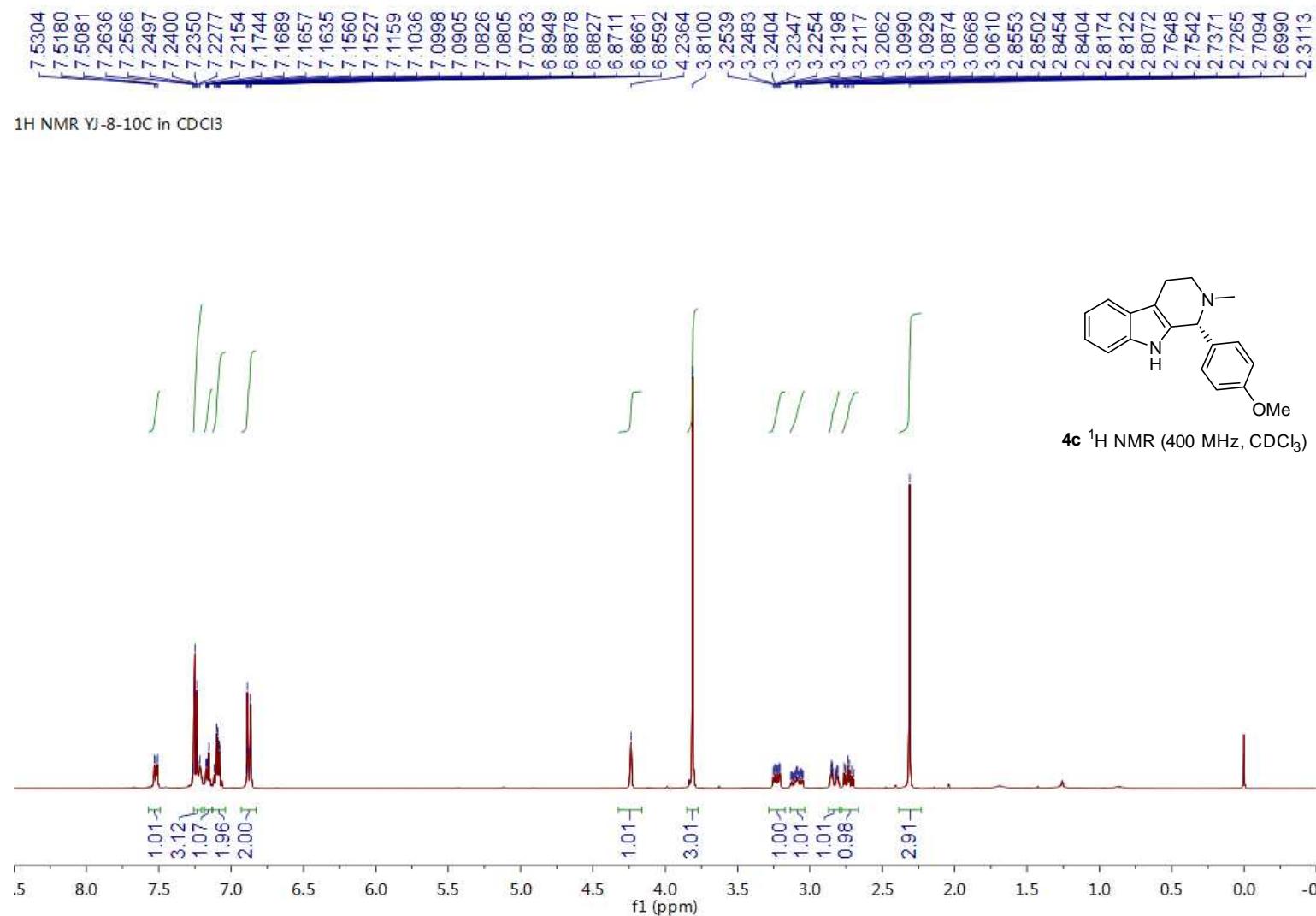


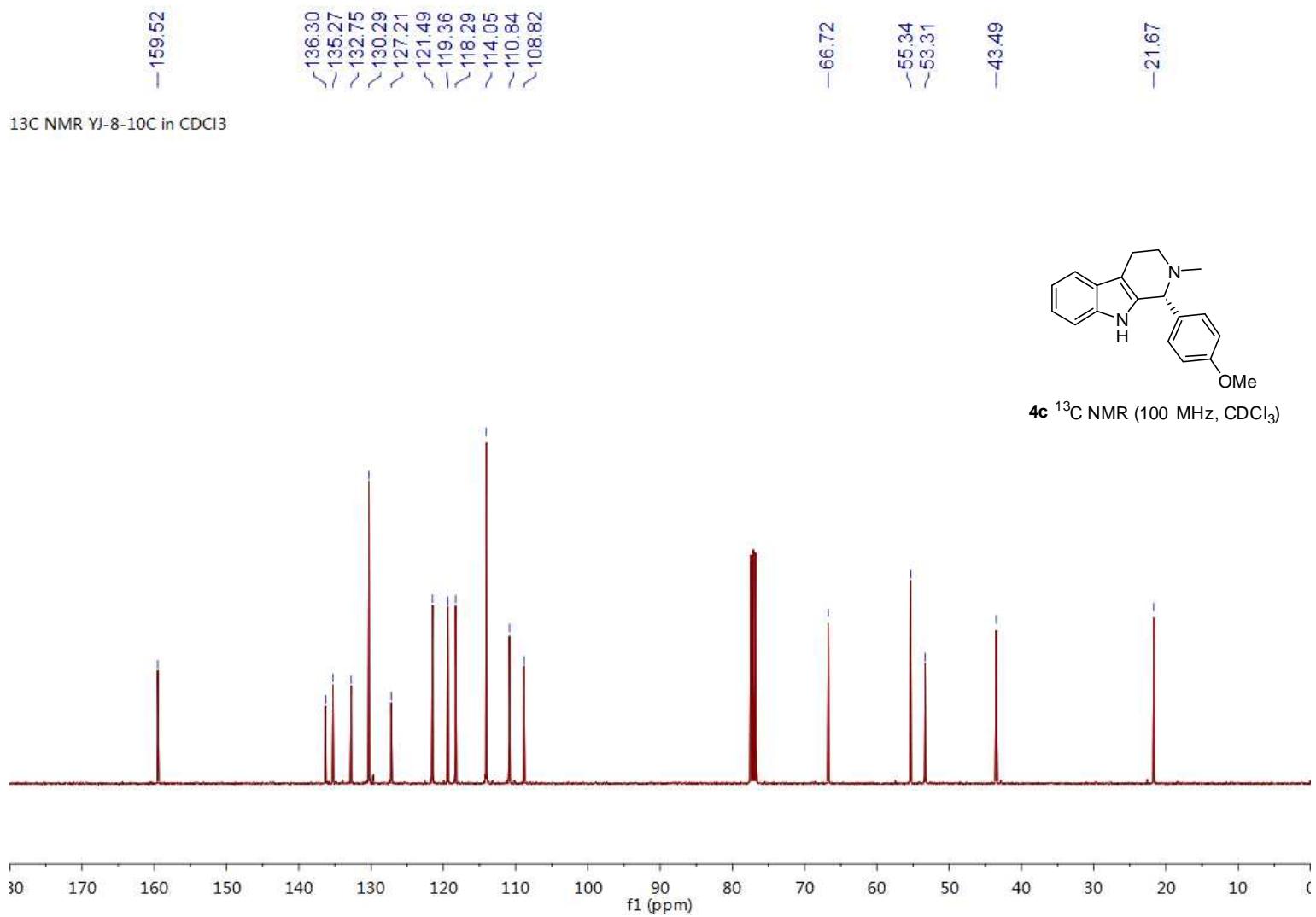


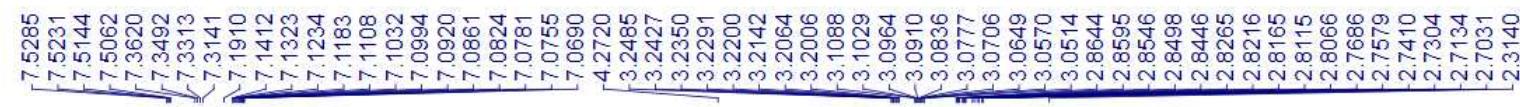
^1H NMR YJ-8-9A in CDCl_3



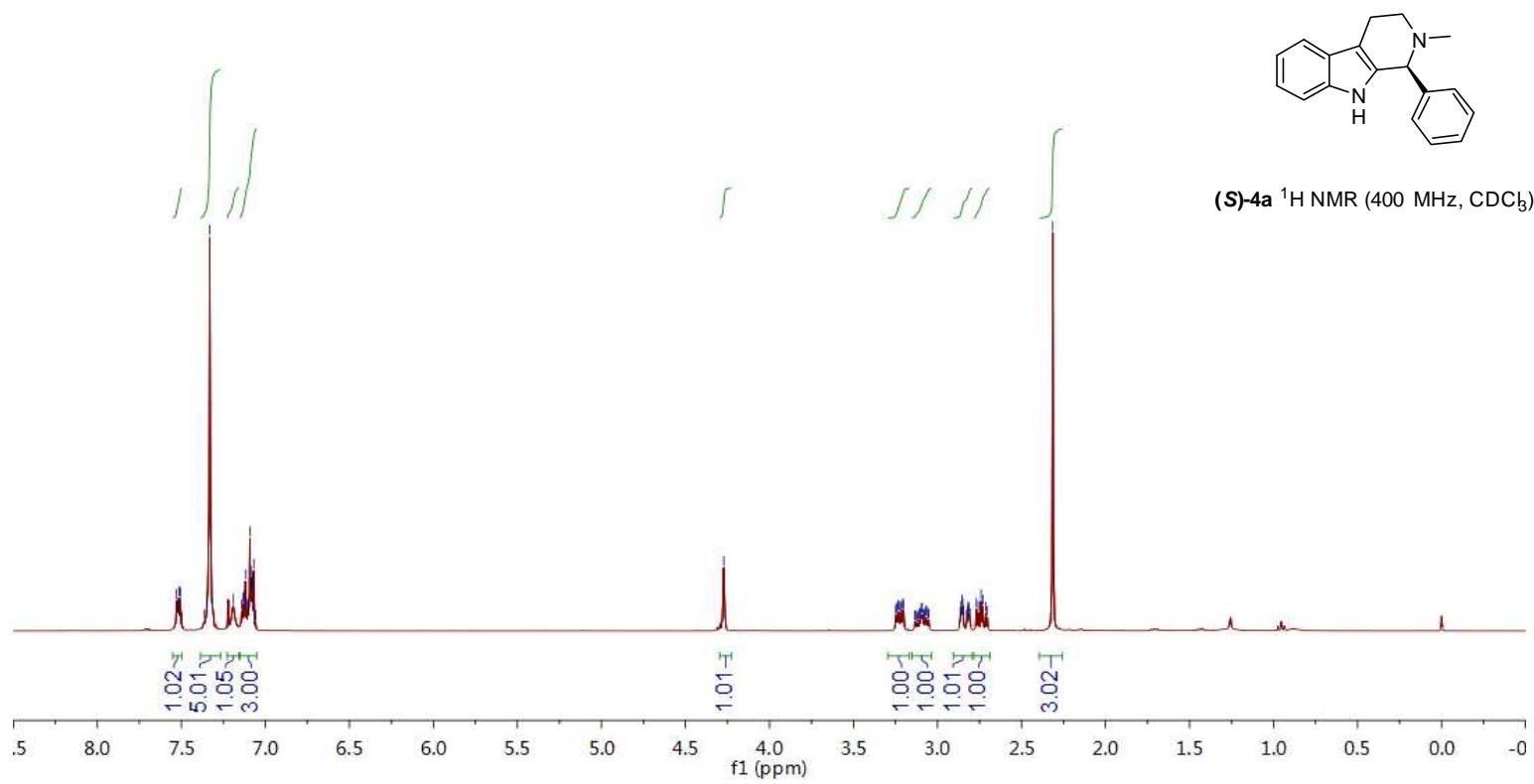






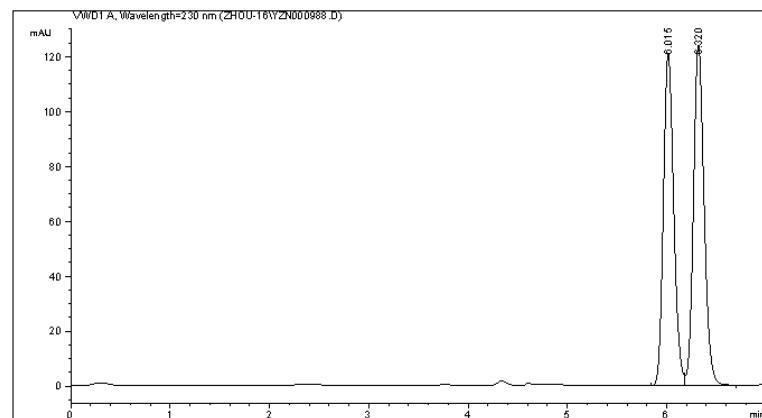


¹H NMR YJ-8-25 in CDCl₃



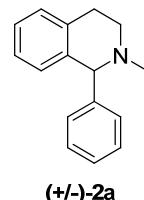
Data File C:\CHEM32\1\DATA\ZHOU-16\YZN000988.D
Sample Name: YJ-6-87+-

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 4/12/2016 9:47:05 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 4/12/2016 9:45:42 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 7:47:19 PM
(modified after loading)
Sample Info : UD-H, Hex/i-PrOH = 98/2, 0.7 mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```



Signal 1: VWD1 A, Wavelength=230 nm

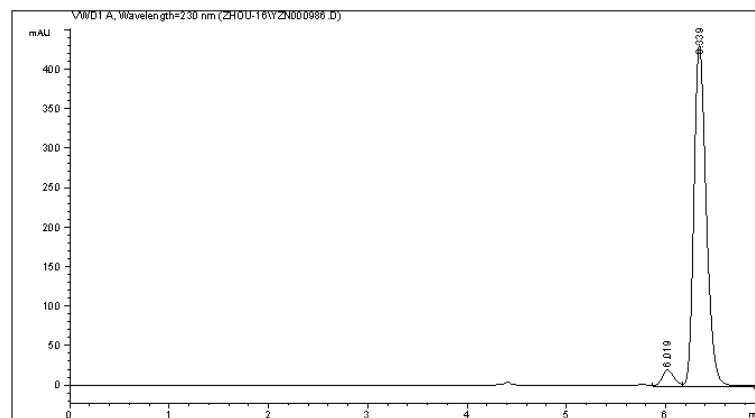
Peak RetTime	Type	Width	Area	Height	Area	
# [min]		[min]	[mAU]	*s	[mAU]	%
1	6.015	BV	0.1118	872.79413	121.08356	48.6320
2	6.320	VB	0.1143	921.89819	124.11383	51.3680

Totals : 1794.69232 245.19739

```
=====
*** End of Report ***
```

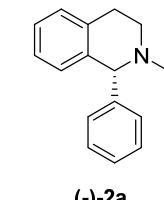
Data File C:\CHEM32\1\DATA\ZHOU-16\YZN000988.D
Sample Name: YJ-6-87B

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 4/12/2016 9:28:57 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 4/12/2016 9:18:21 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 7:47:19 PM
(modified after loading)
Sample Info : UD-H, Hex/i-PrOH = 98/2, 0.7 mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```



Signal 1: VWD1 A, Wavelength=230 nm

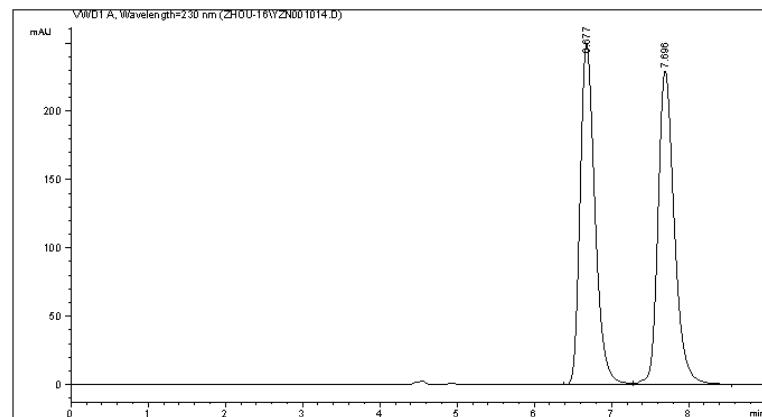
Peak RetTime	Type	Width	Area	Height	Area	
# [min]		[min]	[mAU]	*s	[mAU]	%
1	6.019	VV	0.1283	168.95340	20.17665	4.3576
2	6.339	VB	0.1307	3708.29297	432.08823	95.6424

Totals : 3877.24637 452.26488

```
=====
*** End of Report ***
```

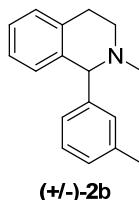
Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001014.D
Sample Name: YJ-6-90B+-

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 4/14/2016 9:56:55 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 4/14/2016 9:55:26 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 7:50:25 PM
(modified after loading)
Sample Info : UJ-H, Hex/i-PrOH = 95/5, 0.7 mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```



Signal 1: VWD1 A, Wavelength=230 nm

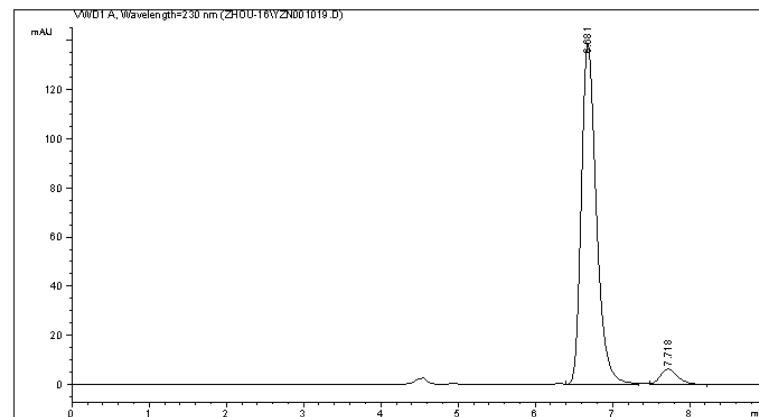
Peak RetTime	Type	Width	Area	Height	Area	
# [min]		[min]	[mAU]	*s [mAU]	%	
1	6.677	VB	0.2050	3323.76807	249.38417	49.7525
2	7.696	VB	0.2231	3356.83643	229.42706	50.2475

Totals : 6680.60449 478.81123

```
=====
*** End of Report ***
```

Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001019.D
Sample Name: YJ-6-90B

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 4/14/2016 11:01:42 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 4/14/2016 10:58:07 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 7:50:25 PM
(modified after loading)
Sample Info : UJ-H, Hex/i-PrOH = 95/5, 0.7 mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

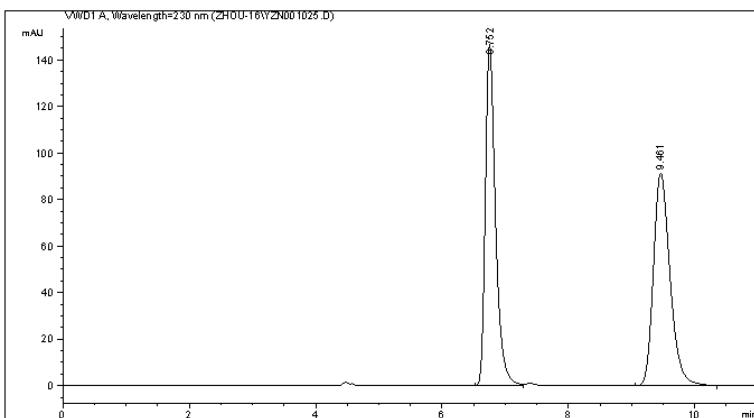
Peak RetTime	Type	Width	Area	Height	Area	
# [min]		[min]	[mAU]	*s [mAU]	%	
1	6.681	VB	0.2076	1879.85925	138.68797	94.9097
2	7.718	VB	0.2405	100.82166	6.30243	5.0903

Totals : 1980.68091 144.99041

```
=====
*** End of Report ***
```

Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001025.D
Sample Name: YJ-6-90A+-

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 4/15/2016 5:15:52 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 4/15/2016 5:09:35 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 7:52:28 PM
(modified after loading)
Sample Info : UJ-H, Hex/i-PrOH = 95/5, 0.7 mL/min, 30oC, 230 nm
```



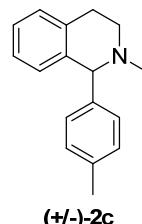
```
=====
Area Percent Report
```

Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

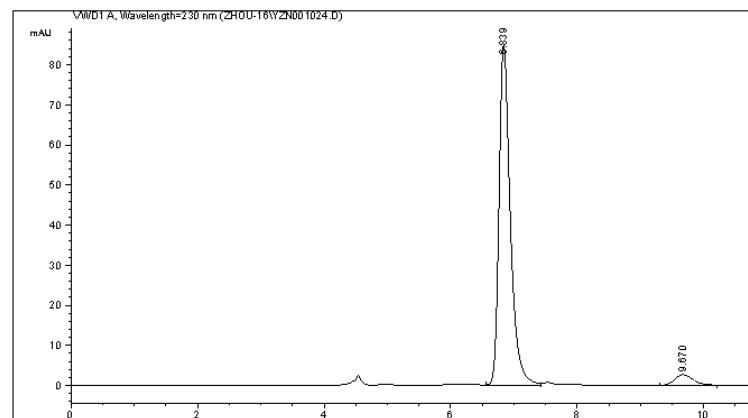
Peak RetTime	Type	Width	Area	Height	Area
# [min]		[min]	[mAU]	*s	[mAU]
1 6.752	BV	0.1700	1642.48145	146.59044	49.5974
2 9.461	BB	0.2800	1669.14624	91.40023	50.4026

Totals : 3311.62769 237.99067



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001024.D
Sample Name: YJ-6-90A

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 4/15/2016 4:56:53 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 4/15/2016 4:47:41 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 7:52:28 PM
(modified after loading)
Sample Info : UJ-H, Hex/i-PrOH = 95/5, 0.7 mL/min, 30oC, 230 nm
```



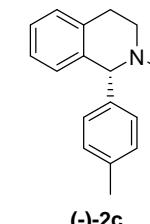
```
=====
Area Percent Report
```

Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

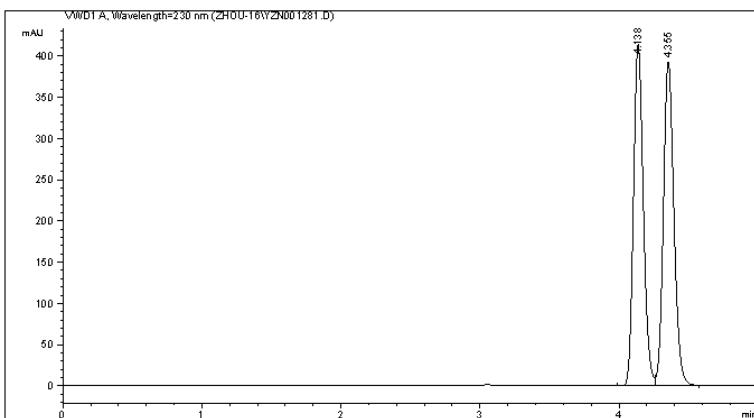
Peak RetTime	Type	Width	Area	Height	Area
# [min]		[min]	[mAU]	*s	[mAU]
1 6.839	BV	0.1866	1049.71289	84.79782	95.1694
2 9.670	BB	0.3106	53.28136	2.66300	4.8306

Totals : 1102.99425 87.46082



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001281.D
Sample Name: YJ-6-101A+

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/7/2016 11:17:37 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/7/2016 11:16:29 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 7:55:08 PM
(modified after loading)
Sample Info : UD-H, Hex/i-PrOH = 90/10, 1.0 mL/min, 30oC, 230 nm
```



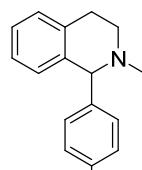
```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area
#		[min]	[min]	[mAU]	%
1	4.138	BV	0.0747	2034.47632	414.56564 49.5893
2	4.355	BV	0.0817	2068.17554	393.52551 50.4107

Totals : 4102.65186 808.09116



(+/-)-2d

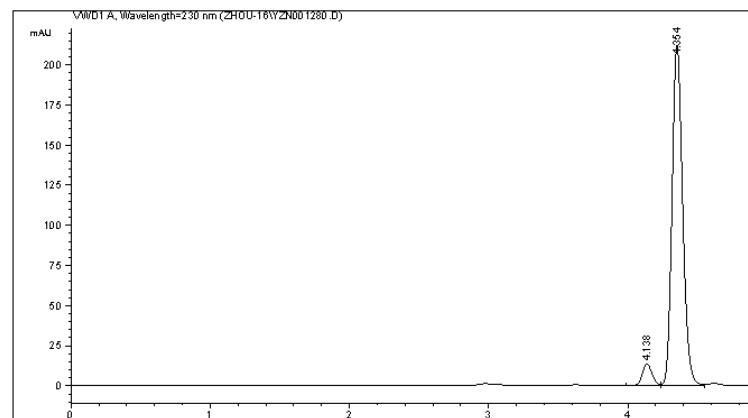
*** End of Report ***

Instrument 1 7/25/2016 7:55:31 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001280.D
Sample Name: YJ-6-101A

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/7/2016 11:10:12 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/7/2016 11:09:28 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 7:55:08 PM
(modified after loading)
Sample Info : UD-H, Hex/i-PrOH = 90/10, 1.0 mL/min, 30oC, 230 nm
```



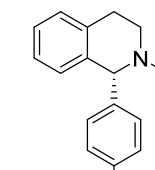
```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area
#		[min]	[min]	[mAU]	%
1	4.138	BV	0.0778	68.04399	13.48343 5.6933
2	4.354	BV	0.0808	1127.11353	212.41176 94.3067

Totals : 1195.15752 225.89518



(-)-2d

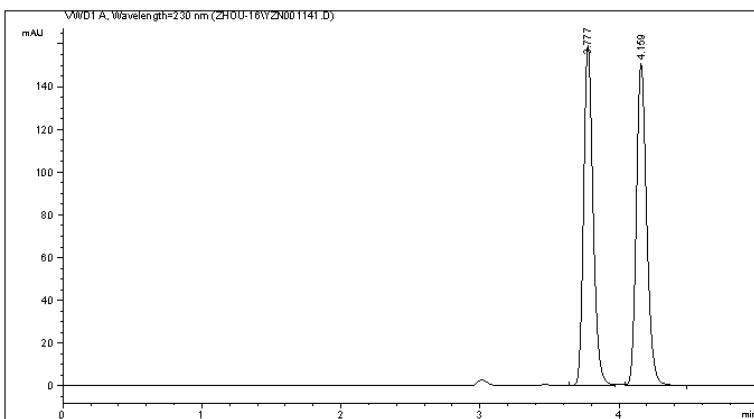
*** End of Report ***

Instrument 1 7/25/2016 7:55:52 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001141.D
Sample Name: YJ-6-92C+-

```
=====
Acq. Operator :                               Location : Vial 1
Acq. Instrument : Instrument 1             Location : Vial 1
Injection Date : 4/22/2016 10:09:32 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 4/22/2016 10:06:06 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 7:55:08 PM
(modified after loading)
Sample Info : UD-H, Hex/i-PrOH = 95/5, 1.0 mL/min, 30oC, 230 nm
```



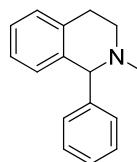
```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area
# [min]		[min]	[mAU]	*s	[mAU]
1 3.777	VV	0.0728	756.84216	159.51071	49.2602
2 4.159	VB	0.0793	779.57513	150.73552	50.7398

Totals : 1536.41730 310.24623



(+/-)-2e

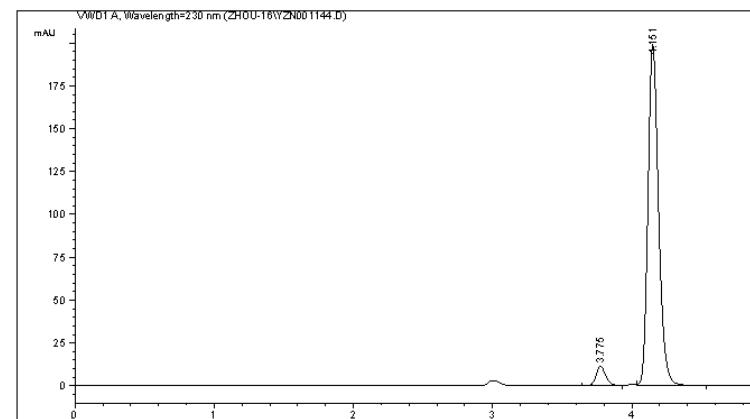
*** End of Report ***

Instrument 1 7/25/2016 7:57:33 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001144.D
Sample Name: YJ-6-92C

```
=====
Acq. Operator :                               Location : Vial 1
Acq. Instrument : Instrument 1             Location : Vial 1
Injection Date : 4/22/2016 10:30:03 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 4/22/2016 10:28:41 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 7:55:08 PM
(modified after loading)
Sample Info : UD-H, Hex/i-PrOH = 95/5, 1.0 mL/min, 30oC, 230 nm
```



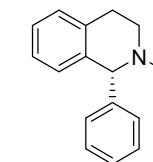
```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area
# [min]		[min]	[mAU]	*s	[mAU]
1 3.775	BV	0.0744	55.18276	11.30744	4.9818
2 4.151	VB	0.0806	1052.50842	198.99071	95.0182

Totals : 1107.69118 210.29815



(-)-2e

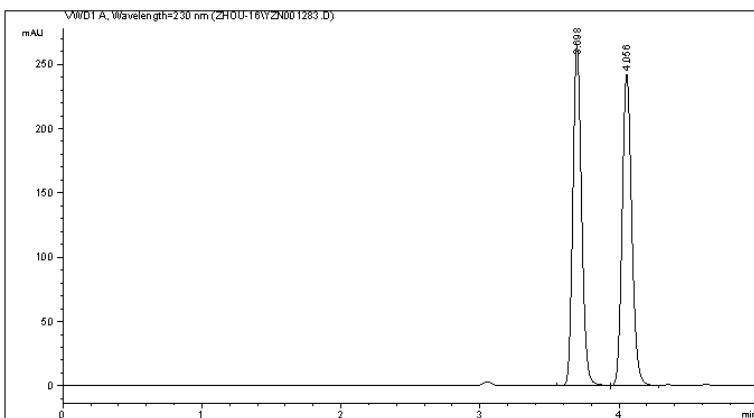
*** End of Report ***

Instrument 1 7/25/2016 7:57:53 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001283.D
Sample Name: YJ-6-101B+-

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/7/2016 11:29:50 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/7/2016 11:29:16 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 7:55:08 PM
(modified after loading)
Sample Info : UD-H, Hex/i-PrOH = 90/10, 1.0 mL/min, 30oC, 230 nm
```



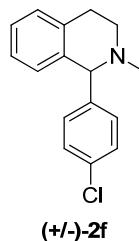
```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

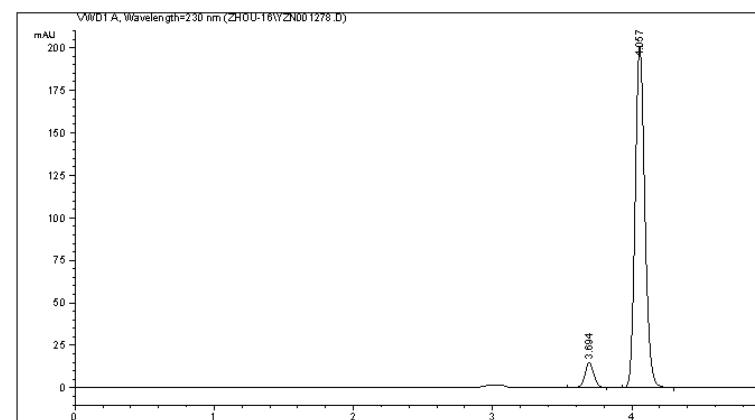
Peak RetTime	Type	Width	Area	Height	Area		
#		[min]	[mAU]	*s	[mAU]	1	%
1	3.698	0.0670	1157.58936	265.09619	49.8513		
2	4.056	0.0735	1164.49731	242.44028	50.1487		

Totals : 2322.08667 507.53647



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001278.D
Sample Name: YJ-6-101B

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/7/2016 10:57:17 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/7/2016 10:56:56 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 7:55:08 PM
(modified after loading)
Sample Info : UD-H, Hex/i-PrOH = 90/10, 1.0 mL/min, 30oC, 230 nm
```



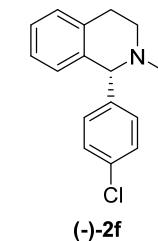
```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

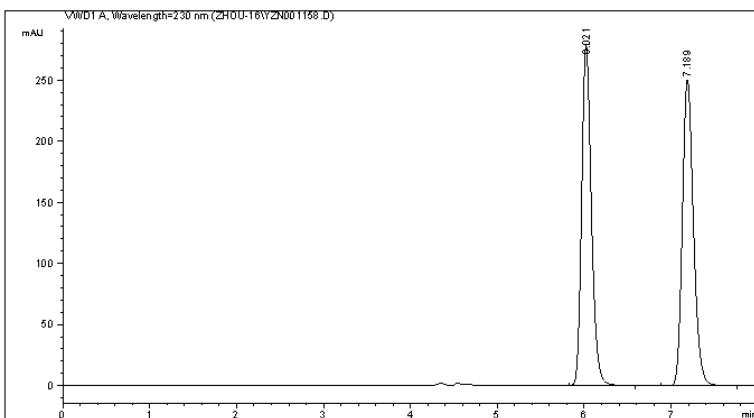
Peak RetTime	Type	Width	Area	Height	Area		
#		[min]	[mAU]	*s	[mAU]	1	%
1	3.694	0.0686	66.68333	14.78734	6.4298		
2	4.057	0.0754	970.41309	200.56610	93.5702		

Totals : 1037.09642 215.35345



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001158.D
Sample Name: YJ-6-93D+-

```
=====
Acq. Operator :                               Location : Vial 1
Acq. Instrument : Instrument 1             Location : Vial 1
Injection Date : 4/23/2016 3:59:44 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 4/23/2016 3:57:44 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 8:01:48 PM
(modified after loading)
Sample Info : UD-H, Hex/i-PrOH = 98/2, 0.7 mL/min, 30oC, 230 nm
```

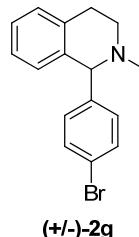


```
=====
Area Percent Report
```

Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

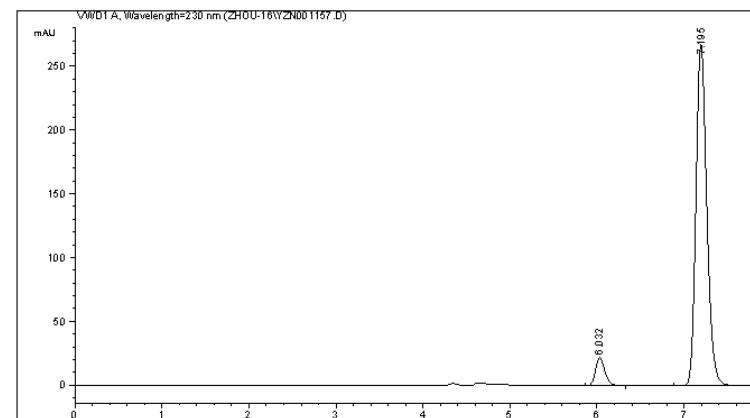
Signal 1: VWD1 A, Wavelength=230 nm

#	RetTime	Type	Width	Area	Height	Area
	[min]		[min]	[mAU]	*s	[mAU]
1	6.021	VB	0.1200	2174.69897	279.14999	49.4372
2	7.189	BB	0.1370	2224.21753	250.62343	50.5628
Totals :				4398.91650	529.77342	



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001157.D
Sample Name: YJ-6-93D

```
=====
Acq. Operator :                               Location : Vial 1
Acq. Instrument : Instrument 1             Location : Vial 1
Injection Date : 4/23/2016 3:49:47 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 4/23/2016 3:48:08 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 8:01:48 PM
(modified after loading)
Sample Info : UD-H, Hex/i-PrOH = 98/2, 0.7 mL/min, 30oC, 230 nm
```

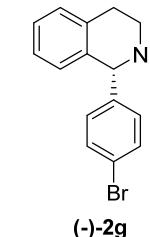


```
=====
Area Percent Report
```

Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

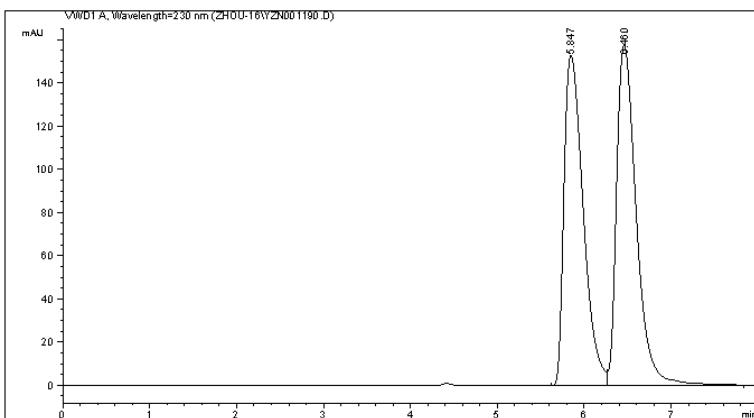
Signal 1: VWD1 A, Wavelength=230 nm

#	RetTime	Type	Width	Area	Height	Area
	[min]		[min]	[mAU]	*s	[mAU]
1	6.032	VB	0.1147	164.21527	22.00098	6.6115
2	7.195	BB	0.1332	2319.56982	267.58517	93.3885
Totals :				2483.78510	289.58615	



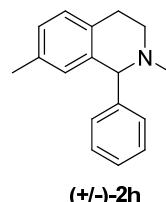
Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001190.D
Sample Name: YJ-6-97A+-

```
=====
Acq. Operator :                               Location : Vial 1
Acq. Instrument : Instrument 1             Location : Vial 1
Injection Date : 4/26/2016 6:19:45 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 4/26/2016 6:19:05 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 8:05:43 PM
(modified after loading)
Sample Info : 0G, Hex/i-PrOH = 99.6/0.4, 0.7 mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```



Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area
# [min]		[min]	[mAU]	*s	[mAU]
1 5.847	BV	0.2412	2342.20557	153.01801	48.4621
2 6.460	BV	0.2439	2490.86523	157.74657	51.5379

Totals : 4833.07080 310.76457

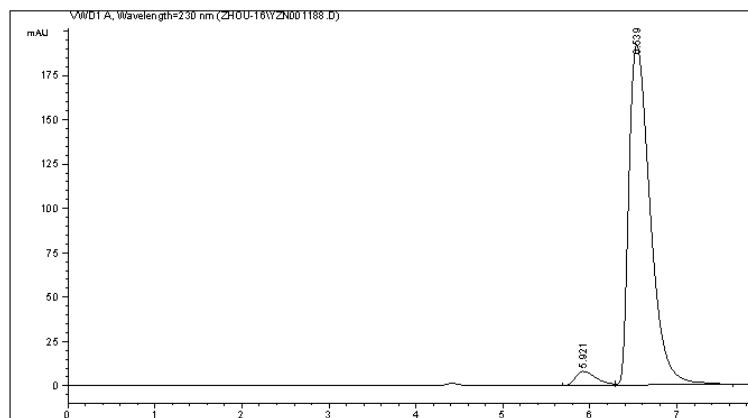
*** End of Report ***

Instrument 1 7/25/2016 8:05:46 PM

Page 1 of 1

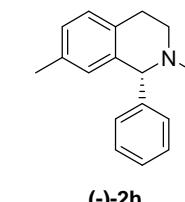
Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001188.D
Sample Name: YJ-6-97A

```
=====
Acq. Operator :                               Location : Vial 1
Acq. Instrument : Instrument 1             Location : Vial 1
Injection Date : 4/26/2016 5:59:34 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 4/26/2016 5:58:43 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 8:05:43 PM
(modified after loading)
Sample Info : 0G, Hex/i-PrOH = 99.6/0.4, 0.7 mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```



Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area
# [min]		[min]	[mAU]	*s	[mAU]
1 5.921	BV	0.2533	126.44871	7.79561	3.7467
2 6.539	BV	0.2647	3248.46802	191.68631	96.2533

Totals : 3374.91673 199.48192

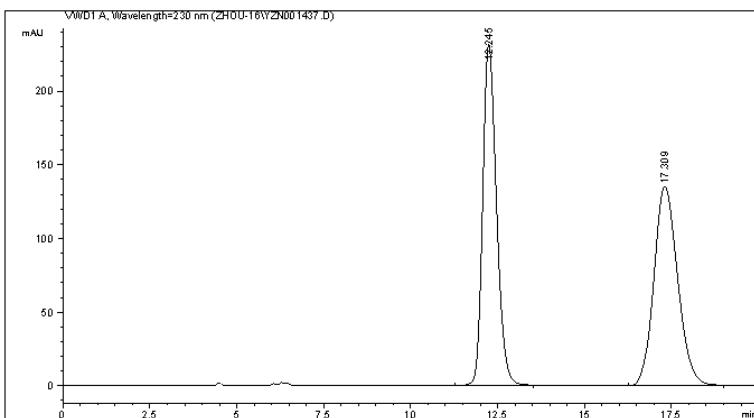
*** End of Report ***

Instrument 1 7/25/2016 8:06:31 PM

Page 1 of 1

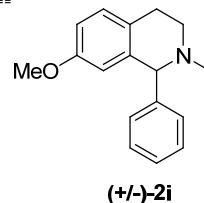
Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001437.D
Sample Name: YJ-7-13A+-

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/23/2016 10:41:44 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/23/2016 10:39:32 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 9:44:36 PM
(modified after loading)
Sample Info : UJ-H, Hex/i-PrOH = 90/10, 0.7 mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```



Peak RetTime	Type	Width	Area	Height	Area	
#	[min]	[min]	[mAU]	*s	[mAU]	%
1	12.245	BB	0.4236	6374.08203	231.14511	49.8491
2	17.309	BB	0.7351	6412.67432	135.25278	50.1509

Totals : 1.27868e4 366.39789

*** End of Report ***

Instrument 1 7/25/2016 9:44:38 PM

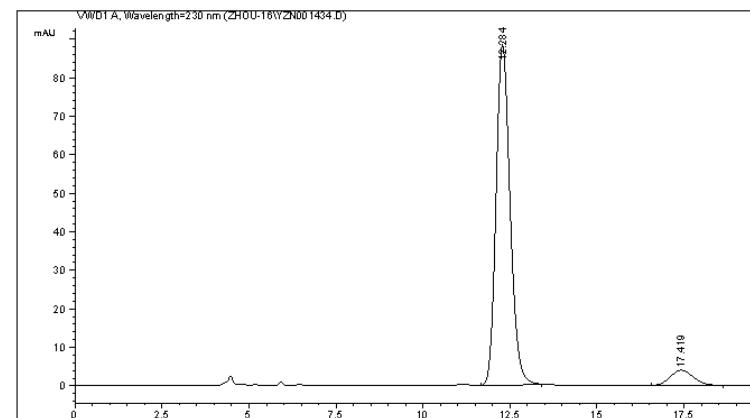
Page 1 of 1

Instrument 1 7/25/2016 8:14:42 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001434.D
Sample Name: YJ-7-13A

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/23/2016 9:41:12 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/23/2016 9:40:02 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 8:14:51 PM
(modified after loading)
Sample Info : UJ-H, Hex/i-PrOH = 90/10, 0.7 mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
```

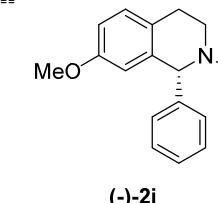
```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area	
#	[min]	[min]	[mAU]	*s	[mAU]	%
1	12.284	BB	0.4330	2485.96509	88.37807	92.9638
2	17.419	BB	0.7181	188.15681	3.98641	7.0362

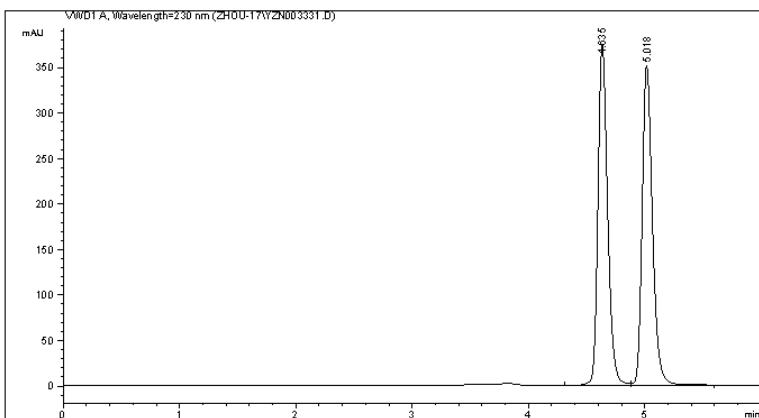
Totals : 2674.12190 92.36447

*** End of Report ***

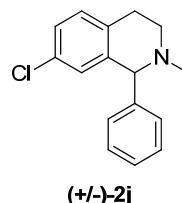


Data File C:\CHEM32\1\DATA\ZHOU-17\YZN003331.D
Sample Name: YJ-8-67A+-

```
Acq. Operator : 0                                         Location : Vial 1
Acq. Instrument : Instrument 1
Injection Date : 2/18/2017 10:13:45 PM
Acq. Method : C:\CHRM32\1\METHODS\DEF LC.M
Last changed : 2/18/2017 10:12:41 PM by 0
                                         (modified after loading)
Analysis Method : C:\CHRM32\1\METHODS\DEF LC.M
Last changed : 2/19/2017 7:15:55 PM by 0
                                         (modified after loading)
Sample Info : AD-H, Hexane-i-PrOH = 95/5, 0.8mL/min, 30oC, 230 nm
```



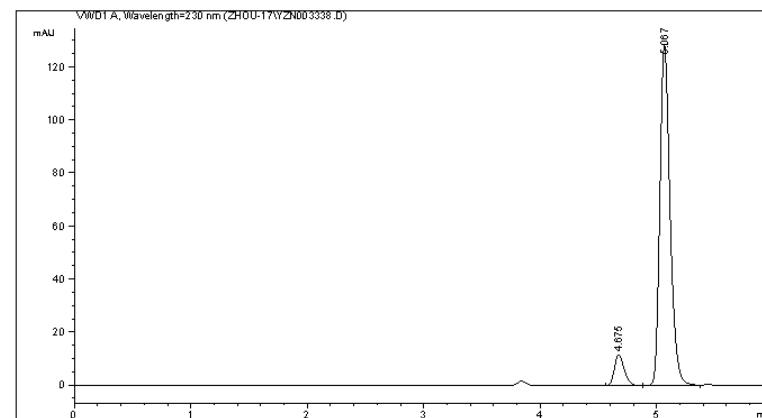
Signal 1: WVUD 1 A, Wavelength=230 nm						
Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]
1	4.635	BV	0.0893	2161.72510	373.91660	50.0947
2	5.018	VB	0.0933	2153.55273	351.52271	49.9053
Totals :				4315.27283	725.43930	



=====
*** End of Report ***

Data File C:\CHEM32\1\DATA\ZHOU-17\YZN003338.I
Sample Name: YJ-8-67A

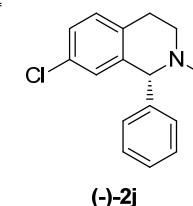
```
=====
Acq. Operator   : 0                                         Location : Vial 1
Acq. Instrument: Instrument 1
Injection Date : 2/19/2017 6:51:42 PM
Acq. Method    : C:\CHEM32\1\METHODS\DEF LC.M
Last changed   : 2/19/2017 6:45:41 PM by 0
                           (modified after loading)
Analysis Method: C:\CHEM32\1\METHODS\DEF LC.M
Last changed   : 2/19/2017 7:15:55 PM by 0
                           (modified after loading)
Sample Info    : AD-H, Hexane-i-PrOH = 95/5, 0.8mL/min, 300C, 230 m
```



=====
Area Percent Report
=====

Sorted by : signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDS

Signal 1: WMD1 A, Wavelength=230 nm						
Peak #	RetTime [min]	Type	Width [min]	Area [mAU]	Height [mAU]	Area %
1	4.675	BV	0.0866	68.86459	11.77551	8.0668
2	5.067	VE	0.0930	784.62659	128.58343	91.9314



Totals : 853.49118 140.3589

=====
*** End of Report ***

Instrument 1 2/19/2017 7:16:24 PM 0

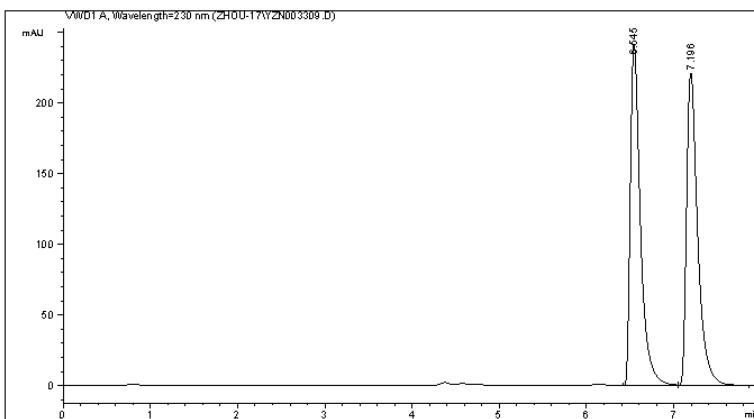
Page 1 of 1

Instrument 1 2/19/2017 7:16:00 PM

Page 1 of

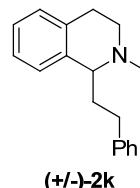
Data File C:\CHEM32\1\DATA\ZHOU-17\YZN003309.D
Sample Name: YJ-8-64+-

```
=====
Acq. Operator : 0
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date : 2/15/2017 9:48:37 PM
Acq. Method   : C:\CHEM32\1\METHODS\DEF LC.M
Last changed   : 2/15/2017 9:47:28 PM by 0
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed   : 2/19/2017 7:22:10 PM by 0
(modified after loading)
Sample Info    : UD-H, Hexane/i-PrOH = 98/2, 0.7 mL/min, 300°C, 230 nm
```



```
=====
Area Percent Report
```

```
Sorted By      : Signal
Multiplier:   : 1.0000
Dilution:     : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```



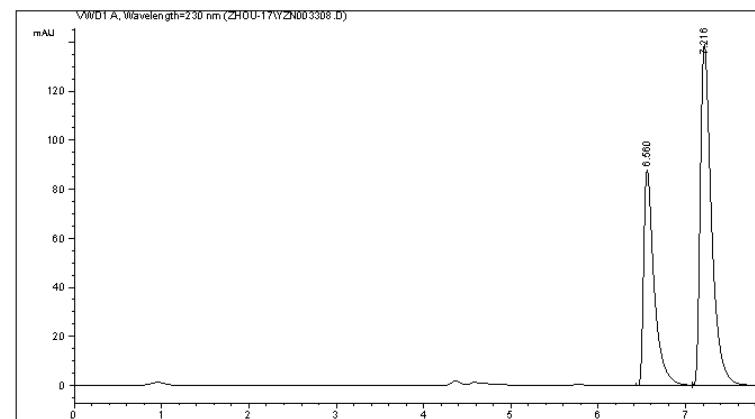
=====
*** End of Report ***

Instrument 1 2/19/2017 7:22:42 PM 0

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-17\YZN003308.D
Sample Name: YJ-8-64

```
=====
Acq. Operator : 0
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date : 2/15/2017 9:38:40 PM
Acq. Method   : C:\CHEM32\1\METHODS\DEF LC.M
Last changed   : 2/15/2017 9:37:50 PM by 0
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed   : 2/19/2017 7:22:10 PM by 0
(modified after loading)
Sample Info    : UD-H, Hexane/i-PrOH = 98/2, 0.7 mL/min, 300°C, 230 nm
```



```
=====
Area Percent Report
```

```
Sorted By      : Signal
Multiplier:   : 1.0000
Dilution:     : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime Type Width Area Height Area

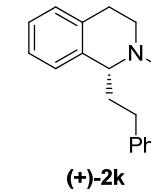
#	[min]	[min]	mAU	*s	[mAU]	%
1	6.560	BV	0.1235	734.09674	88.00777	37.1689
2	7.216	VB	0.1348	1240.93079	138.96794	62.8311

Totals :

1975.02753	226.97572
------------	-----------

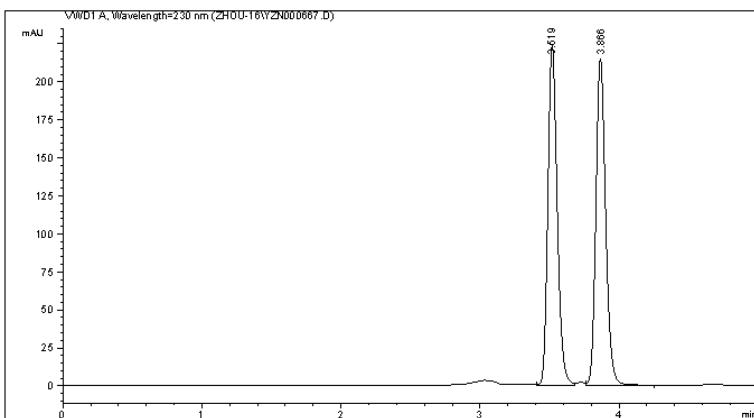
=====
*** End of Report ***

Instrument 1 2/19/2017 7:22:13 PM 0



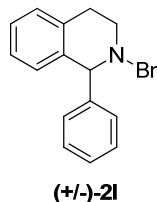
Data File C:\CHEM32\1\DATA\ZHOU-16\YZN000667.D
Sample Name: YJ-6-67B+-

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 3/19/2016 3:22:12 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/19/2016 3:05:26 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 9:44:36 PM
(modified after loading)
Sample Info : AD-H, H2-PrOH = 90/10, 1.0 mL/min, 30 oC, 230 nm
```



```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```



Peak RetTime	Type	Width	Area	Height	Area	
#	[min]	[min]	[mAU]	*s	[mAU]	%
1	3.519	0.0718	1043.47729	224.11391	49.7838	
2	3.866	0.0744	1052.54089	215.67365	50.2162	

Totals : 2096.01819 439.78755

=====
*** End of Report ***

Instrument 1 7/25/2016 9:45:14 PM

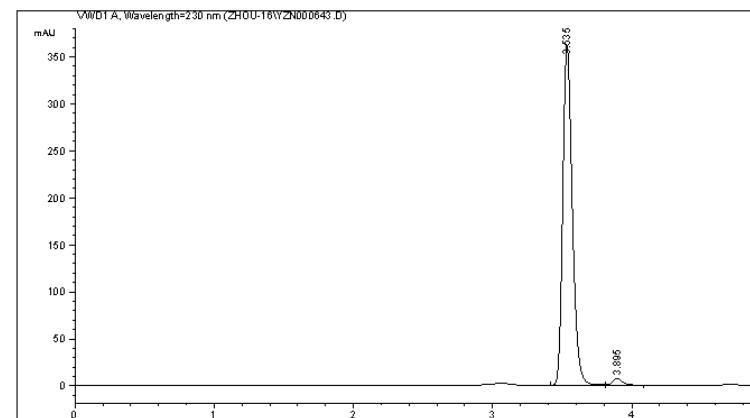
Page 1 of 1

Instrument 1 7/25/2016 8:08:37 PM

Page 1 of 1

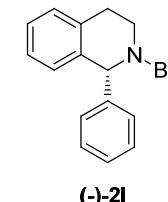
Data File C:\CHEM32\1\DATA\ZHOU-16\YZN000643.D
Sample Name: YJ-6-67B

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 3/18/2016 4:37:37 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/18/2016 4:28:33 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 8:08:34 PM
(modified after loading)
Sample Info : AD-H, H2-PrOH = 90/10, 1.0 mL/min, 30 oC, 230 nm
```



```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```



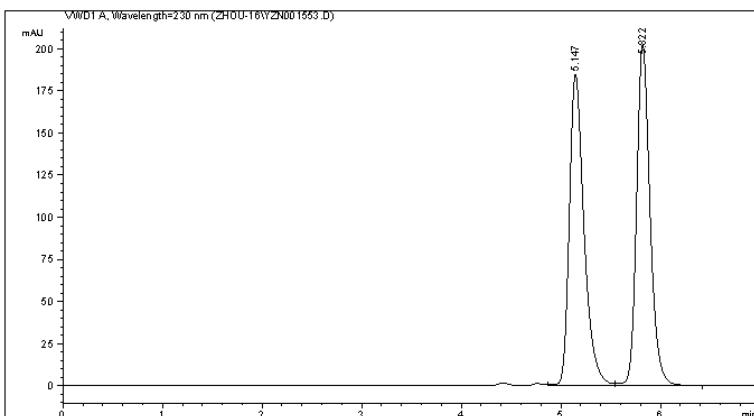
Peak RetTime	Type	Width	Area	Height	Area	
#	[min]	[min]	[mAU]	*s	[mAU]	%
1	3.535	0.0732	1733.31335	362.79758	97.7554	
2	3.895	0.0809	39.79905	7.32114	2.2446	

Totals : 1773.11240 370.11872

=====
*** End of Report ***

Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001553.D
Sample Name: YJ-7-4A++

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 6/8/2016 4:00:54 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/8/2016 3:32:26 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 9:43:18 PM
(modified after loading)
Sample Info : AD-H, Hex/i-PrOH = 98/2, 0.7 mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
```

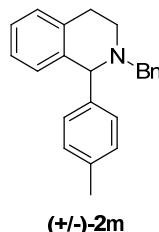
```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area
#		[min]	[mAU]	*s	[mAU]
1	5.147	0.1549	1899.97949	184.64914	49.9071
2	5.822	0.1434	1907.05432	202.39923	50.0929

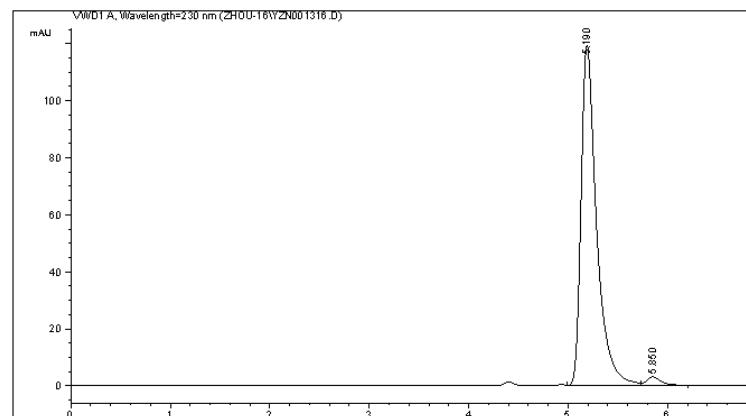
Totals : 3807.03381 387.24837

*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001316.D
Sample Name: YJ-7-4A

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/10/2016 3:54:07 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/10/2016 3:51:28 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 9:43:18 PM
(modified after loading)
Sample Info : AD-H, Hex/i-PrOH = 98/00, 0.7 mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
```

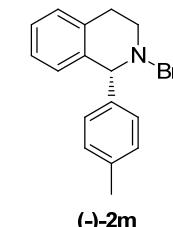
```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area
#		[min]	[mAU]	*s	[mAU]
1	5.190	0.1597	1293.68225	119.53790	97.5849
2	5.850	0.1508	32.01711	3.06859	2.4151

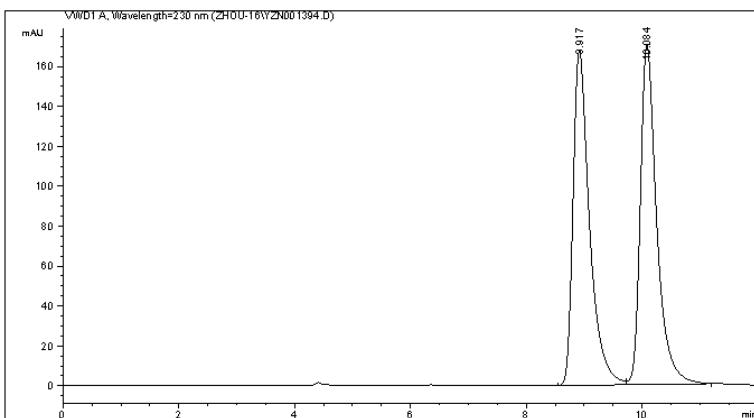
Totals : 1325.69936 122.60649

*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001394.D
Sample Name: YJ-7-4B++

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/17/2016 10:06:35 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/17/2016 10:04:29 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 8:12:06 PM
(modified after loading)
Sample Info : AD-H, Hex/i-PrOH = 99.2/0.8, 0.7 mL/min, 30oC, 230 nm
```



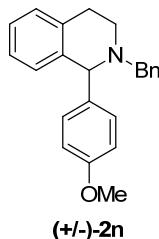
```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

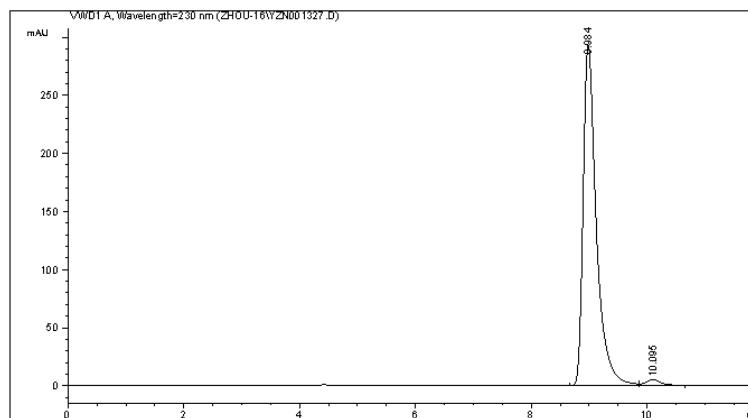
#	RetTime	Type	Width	Area	Height	Area	
	[min]		[min]	[mAU]	*s	[mAU]	%
1	8.917	BV	0.2942	3309.72656	167.81749	49.8220	
2	10.084	BV	0.2926	3333.36963	170.19594	50.1780	

Totals : 6643.09619 338.01343



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001327.D
Sample Name: YJ-7-4B

```
=====
Acq. Operator : 
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/10/2016 6:36:22 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/10/2016 6:35:37 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 8:12:06 PM
(modified after loading)
Sample Info : AD-H, Hex/i-PrOH = 99.2/0.8, 0.7 mL/min, 30oC, 230 nm
```



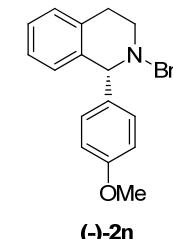
```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

#	RetTime	Type	Width	Area	Height	Area	
	[min]		[min]	[mAU]	*s	[mAU]	%
1	8.984	BV	0.2371	4647.17334	293.51337	97.9049	
2	10.095	BV	0.2849	99.44447	5.15190	2.0951	

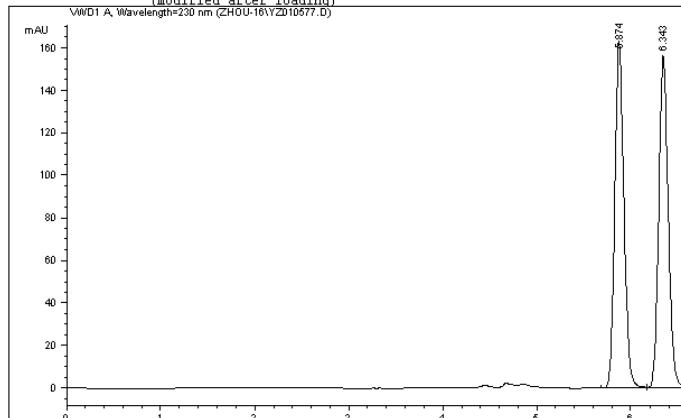
Totals : 4746.61781 298.66526



Data File C:\HPCHEM\1\DATA\ZHOU-16\Y2010577.D

AD-H, H/i-PrOH = 98/2, 0.7 mL/min, 30 oC, 230 nm

=====
Injection Date : 16/05/2016 11:09:02
Sample Name : YJ-7-B+-
Acq. Operator :
Acq. Method : C:\HPCHEM\1\METHODS\DEF LC.M
Last changed : 16/05/2016 10:22:30 by
(modified after loading)
Analysis Method : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed : 25/07/2016 19:39:08 by
(modified after loading)

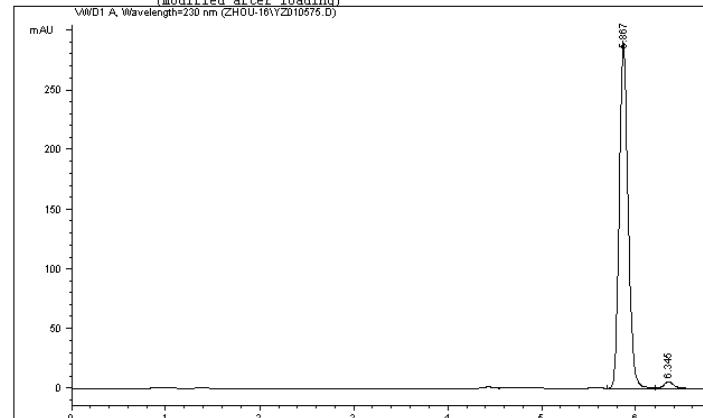


Sample Name: YJ-7-B+-

Data File C:\HPCHEM\1\DATA\ZHOU-16\Y2010575.D

AD-H, H/i-PrOH = 98/2, 0.7 mL/min, 30 oC, 230 nm

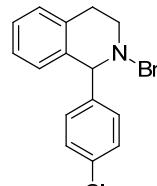
=====
Injection Date : 16/05/2016 10:51:53
Sample Name : YJ-7-B
Acq. Operator :
Acq. Method : C:\HPCHEM\1\METHODS\DEF LC.M
Last changed : 16/05/2016 10:22:30 by
(modified after loading)
Analysis Method : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed : 25/07/2016 19:38:12 by
(modified after loading)



Sample Name: YJ-7-B

=====
Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Sample Amount : 1.00000 [ng/ul] (not used in calc.)



Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area	
#	[min]	[min]	mAU	*s	[mAU]	%
1	5.874	VV	0.1036	1102.66138	163.18604	50.1811
2	6.343	VB	0.1092	1094.70215	156.58330	49.8189

Totals : 2197.36353 319.76933

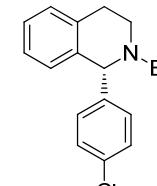
Results obtained with enhanced integrator!

*** End of Report ***

(±)-2o

=====
Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Sample Amount : 1.00000 [ng/ul] (not used in calc.)



Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area	
#	[min]	[min]	mAU	*s	[mAU]	%
1	5.867	VV	0.0996	1897.85339	290.34201	97.7297
2	6.345	VV	0.1163	44.08792	5.71054	2.2703

Totals : 1941.94132 296.05255

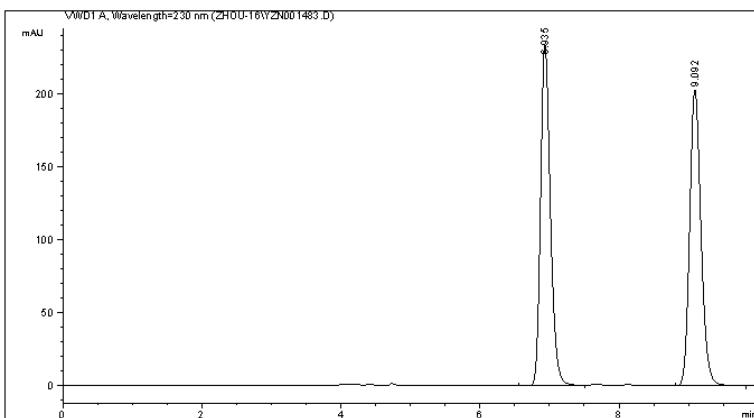
Results obtained with enhanced integrator!

*** End of Report ***

(-)-2o

Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001483.D
Sample Name: YJ-7-15B+-

```
=====
Acq. Operator :                               Location : Vial 1
Acq. Instrument : Instrument 1             Location : Vial 1
Injection Date : 5/28/2016 7:11:28 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/28/2016 7:10:43 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 6:15:59 PM
(modified after loading)
Sample Info : AD-H, Hex/i-PrOH = 98/2, 0.7 mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
```

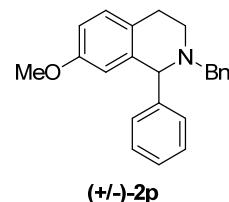
```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area
# [min]		[min]	[mAU]	*s	[mAU]
1 6.935	BB	0.1513	2301.68066	233.74991	50.0196
2 9.092	BB	0.1744	2299.87524	202.99535	49.9804

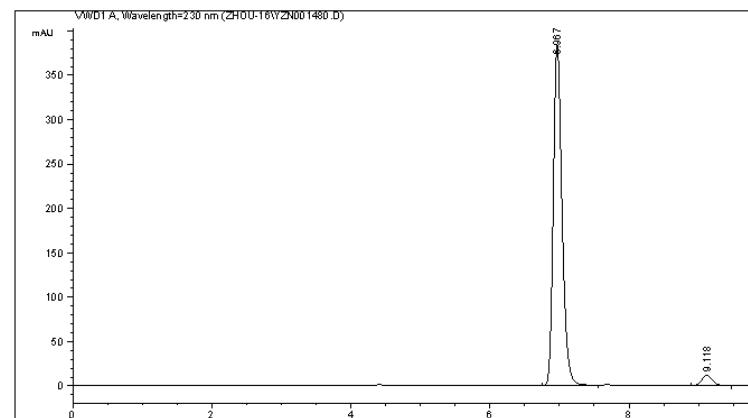
Totals : 4601.55591 436.74525

*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001480.D
Sample Name: YJ-7-15B

```
=====
Acq. Operator :                               Location : Vial 1
Acq. Instrument : Instrument 1             Location : Vial 1
Injection Date : 5/28/2016 6:38:18 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/28/2016 6:37:35 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 6:15:59 PM
(modified after loading)
Sample Info : AD-H, Hex/i-PrOH = 98/2, 0.7 mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
```

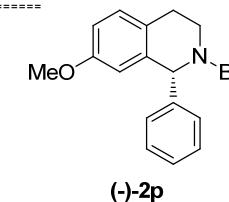
```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area
# [min]		[min]	[mAU]	*s	[mAU]
1 6.967	VB	0.1345	3374.26196	384.25894	96.3851
2 9.118	BB	0.1654	126.55072	11.84902	3.6149

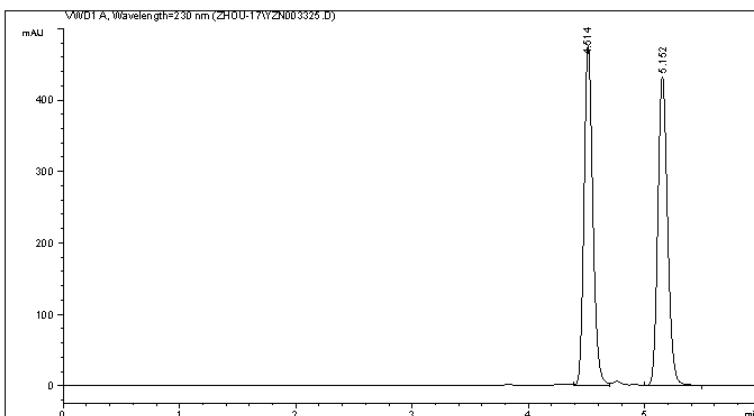
Totals : 3500.81268 396.10796

*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-17\YZN003325.D
Sample Name: YJ-8-67B+-

```
=====
Acq. Operator : 0
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date : 2/18/2017 4:16:47 PM
Acq. Method   : C:\CHEM32\1\METHODS\DEF LC.M
Last changed   : 2/18/2017 4:15:19 PM by 0
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed   : 2/19/2017 7:15:55 PM by 0
(modified after loading)
Sample Info     : AD-H, Hexane/i-PrOH = 95/5, 0.8mL/min, 30oC, 230 nm
```



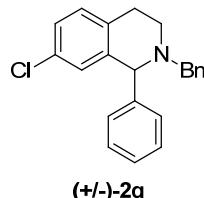
```
=====
Area Percent Report
=====
```

Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area	
# [min]		[min]	[mAU]	*s	[mAU]	%
1 4.514	VV	0.0805	2516.94360	477.26559	49.9573	
2 5.152	VV	0.0897	2521.24121	433.69730	50.0427	

Totals : 5038.18481 910.96289



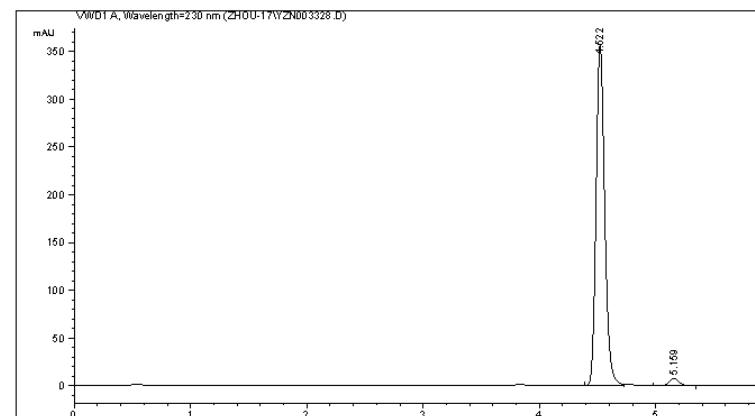
=====
*** End of Report ***
=====

Instrument 1 2/19/2017 7:19:14 PM 0

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-17\YZN003328.D
Sample Name: YJ-8-67B

```
=====
Acq. Operator : 0
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date : 2/18/2017 5:06:43 PM
Acq. Method   : C:\CHEM32\1\METHODS\DEF LC.M
Last changed   : 2/18/2017 5:05:14 PM by 0
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed   : 2/19/2017 7:15:55 PM by 0
(modified after loading)
Sample Info     : AD-H, Hexane/i-PrOH = 95/5, 0.8mL/min, 30oC, 230 nm
```



```
=====
Area Percent Report
=====
```

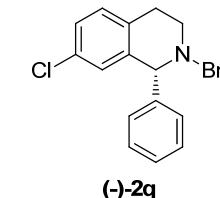
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area	
# [min]		[min]	[mAU]	*s	[mAU]	%
1 4.522	VV	0.0809	1855.14490	357.57715	97.3386	
2 5.159	VV	0.0962	50.72193	7.95742	2.6614	

Totals : 1905.86683 365.53456

=====
*** End of Report ***
=====

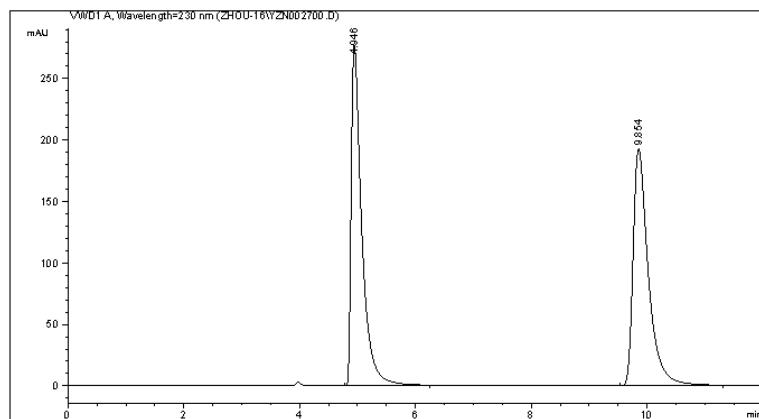


Instrument 1 2/19/2017 7:20:33 PM 0

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-16\YZN002700.D
Sample Name: YJ-8-10D+-

```
=====
Acq. Operator : 0
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 10/31/2016 3:17:26 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 10/31/2016 3:16:19 PM by 0
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 12/17/2016 11:12:34 AM by 0
(modified after loading)
Sample Info : IC, Hexane/i-PrOH = 70/30, 0.8 mL/min, 30oC, 230 nm
```



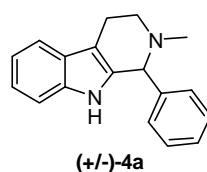
```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

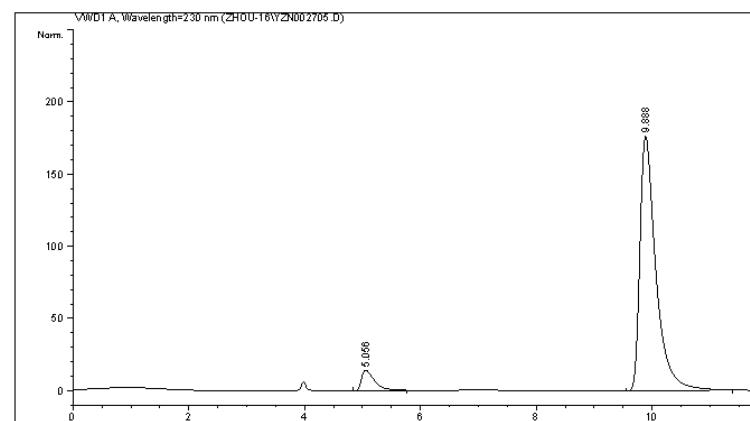
Peak RetTime	Type	Width	Area	Height	Area	
# [min]		[min]	[mAU]	*s	[mAU]	1 %
1 4.946	VB	0.1857	3514.16748	277.30234	49.9325	
2 9.854	BB	0.2715	3523.67017	192.78810	50.0675	

Totals : 7037.83765 470.09044



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN002705.D
Sample Name: YJ-8-10D

```
=====
Acq. Operator : 0
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 10/31/2016 8:55:15 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 10/31/2016 8:51:06 PM by 0
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 12/17/2016 11:14:05 AM by 0
(modified after loading)
Sample Info : IC, Hexane/i-PrOH = 70/30, 0.8 mL/min, 30oC, 230 nm
```



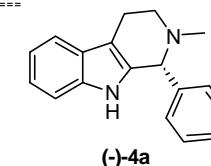
```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area	
# [min]		[min]	[mAU]	*s	[mAU]	1 %
1 5.056	BB	0.2261	218.57561	14.09365	6.0728	
2 5.888	BB	0.2849	3380.66309	176.29660	93.9272	

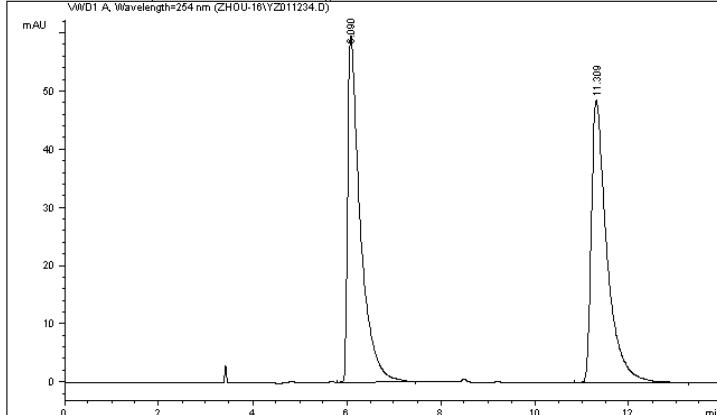
Totals : 3599.23869 190.39025



Data File C:\HPCHEM\1\DATA\ZHOU-16\Y2011234.D

IC, Hexane/iPrOH = 70/30, 0.7 mL/min, 30 oC, 254 nm

=====
Injection Date : 21/09/2016 09:49:04
Sample Name : YJ-7-81+- Location : Vial 1
Acc. Operator :
Acq. Method : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed : 21/09/2016 09:39:09 by
(modified after loading)
Analysis Method : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed : 17/12/2016 10:57:23 by
(modified after loading)

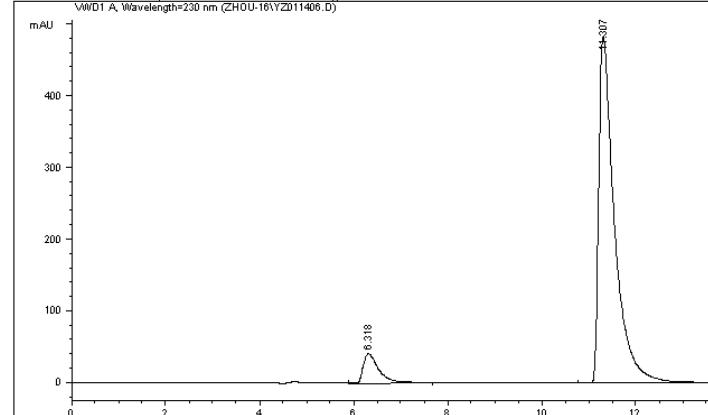


Sample Name: YJ-7-81+-

Data File C:\HPCHEM\1\DATA\ZHOU-16\Y2011406.D

IC, Hexane/iPrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm

=====
Injection Date : 01/11/2016 09:10:29
Sample Name : YJ-8-9A Location : Vial 1
Acc. Operator :
Acq. Method : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed : 01/11/2016 08:49:50 by
(modified after loading)
Analysis Method : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed : 17/12/2016 10:55:34 by
(modified after loading)



Sample Name: YJ-8-9A

=====
Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000

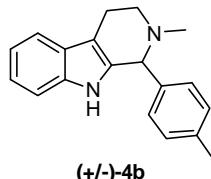
Signal 1: VWD1 A, Wavelength=254 nm

#	RetTime	Type	Width	Area	Height	Area
	[min]		[min]	[mAU]	*s	[mAU]
1	6.090	VB	0.2667	1115.77148	59.53540	49.8372
2	11.309	BB	0.3364	1123.06030	48.62826	50.1628

Totals : 2238.83179 108.16366

Results obtained with enhanced integrator!

*** End of Report ***



(+/-)-4b

=====
Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000

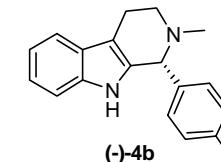
Signal 1: VWD1 A, Wavelength=230 nm

#	RetTime	Type	Width	Area	Height	Area
	[min]		[min]	[mAU]	*s	[mAU]
1	6.318	VB	0.3679	1008.58759	41.09137	8.2139
2	11.307	VB	0.3390	1.12705e4	483.37457	91.7861

Totals : 1.22791e4 524.46594

Results obtained with enhanced integrator!

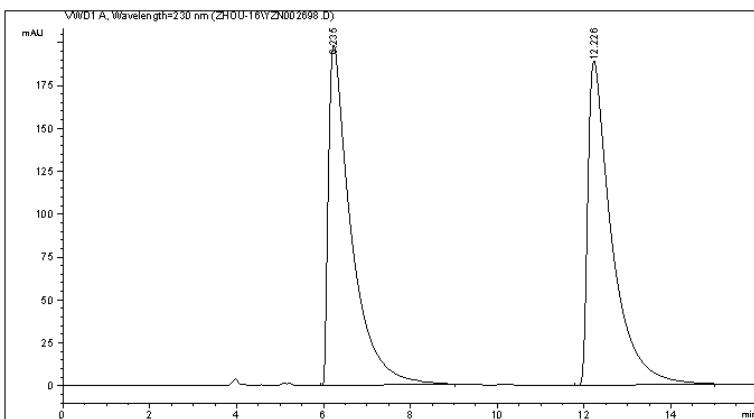
*** End of Report ***



(-)-4b

Data File C:\CHEM32\1\DATA\ZHOU-16\YZN002698.D
Sample Name: YJ-8-10C+-

```
=====
Acq. Operator : 0
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 10/31/2016 2:40:44 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 10/31/2016 2:27:18 PM by 0
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 12/17/2016 11:10:28 AM by 0
(modified after loading)
Sample Info : IC, Hexane/i-PrOH = 70/30, 0.8 mL/min, 30oC, 230 nm
```



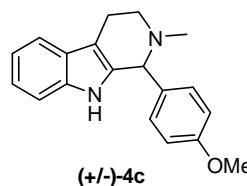
```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

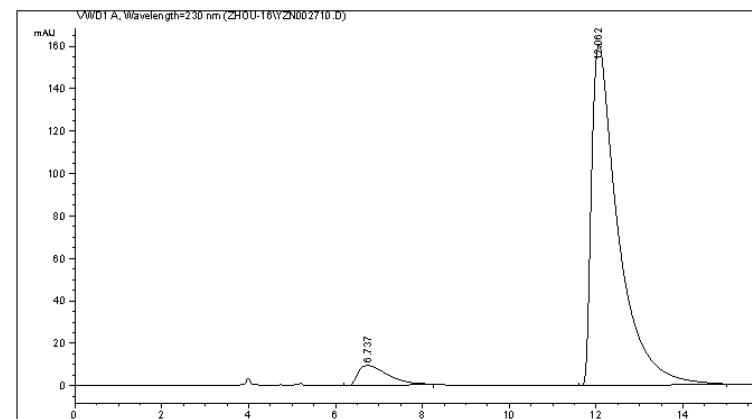
Peak RetTime	Type	Width	Area	Height	Area	
# [min]		[min]	[mAU]	*s	[mAU]	%
1 6.235	VB	0.5162	7172.03760	198.48344	49.9516	
2 12.226	BB	0.5453	7185.93018	189.08002	50.0484	

Totals : 1.43580e4 387.56346



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN002710.D
Sample Name: YJ-8-10C-

```
=====
Acq. Operator : 0
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 10/31/2016 10:24:55 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 10/31/2016 10:16:44 PM by 0
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 12/17/2016 11:11:38 AM by 0
(modified after loading)
Sample Info : IC, Hexane/i-PrOH = 70/30, 0.8 mL/min, 30oC, 230 nm
```



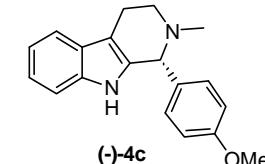
```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

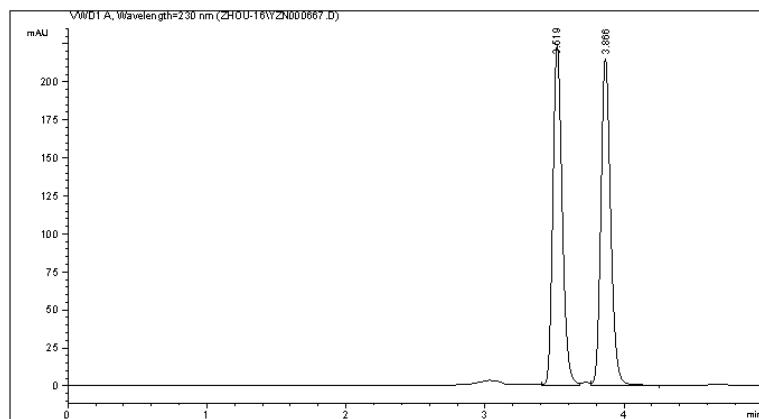
Peak RetTime	Type	Width	Area	Height	Area	
# [min]		[min]	[mAU]	*s	[mAU]	%
1 6.737	VB	0.7350	452.88751	9.38207	6.2161	
2 12.062	BB	0.6189	6832.87402	160.52557	93.7839	

Totals : 7285.76154 169.90764



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN000667.D
Sample Name: YJ-6-67E+-

```
=====
Acq. Operator : Location : Vial 1
Acq. Instrument : Instrument 1
Injection Date : 3/19/2016 3:22:12 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/19/2016 3:05:26 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/25/2016 9:44:36 PM
(modified after loading)
Sample Info : AD-H, H/i-PrOH = 90/10, 1.0 mL/min, 30 oC, 230 nm
```



```
=====
Area Percent Report
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area
# [min]		[min]	[mAU]	*s	[mAU]
1 3.519	VV	0.0718	1043.47729	224.11391	49.7838
2 3.866	VB	0.0744	1052.54089	215.67365	50.2162

Totals : 2096.01819 439.78755

*** End of Report ***



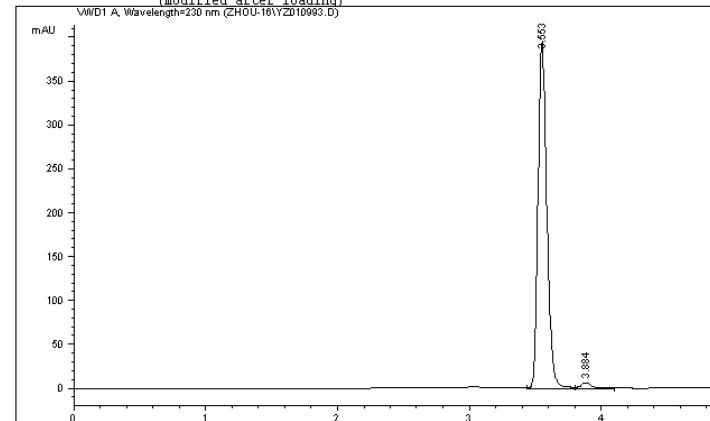
(±)-2l

Data File C:\HPCHEM\1\DATA\ZHOU-16\YZ010993.D

Sample Name: YJ-7-62

```
AD-H, H/i-PrOH = 90/10, 1.0 mL/min, 30 oC, 230 nm
```

```
=====
Injection Date : 23/08/2016 21:01:10
Sample Name : YJ-7-62
Location : Vial 1
Acq. Operator :
Acq. Method : C:\HPCHEM\1\METHODS\DEF_LC1.M
Last changed : 23/08/2016 20:42:07 by
(modified after loading)
Analysis Method : C:\HPCHEM\1\METHODS\DEF_LC1.M
Last changed : 27/08/2016 09:12:59 by
(modified after loading)
```



```
=====
Area Percent Report
```

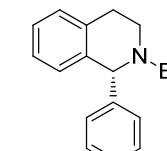
```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak RetTime	Type	Width	Area	Height	Area
# [min]		[min]	[mAU]	*s	[mAU]
1 3.553	VV	0.0721	1843.46289	393.63498	97.8740
2 3.884	VV	0.0861	40.04324	6.81000	2.1260

Totals : 1883.50613 400.44498

Results obtained with enhanced integrator!



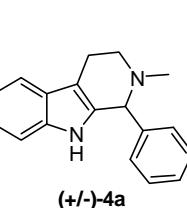
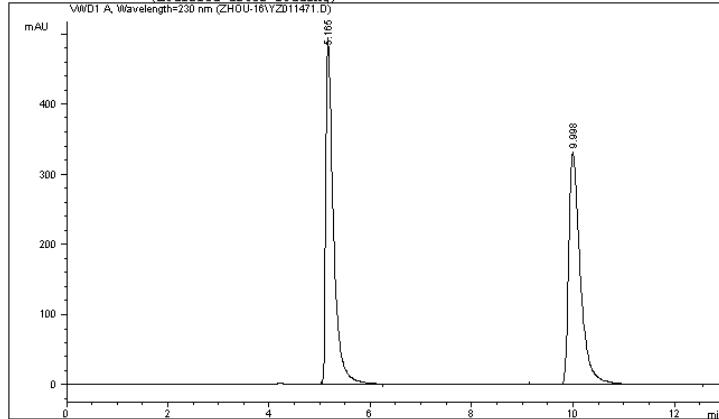
(-)-2l

*** End of Report ***

Data File C:\HPCHEM\1\DATA\ZHOUE-16\Y2011471.D

IC, Hexane/iPrOH = 70/30, 0.8 mL/min, 30 oC, 230 nm

=====
Injection Date : 19/11/2016 10:44:16
Sample Name : YJ-8-25+
Acq. Operator :
Acq. Method : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed : 19/11/2016 10:41:46 bv
(modified after loading)
Analysis Method : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed : 17/12/2016 10:59:15 bv
(modified after loading)



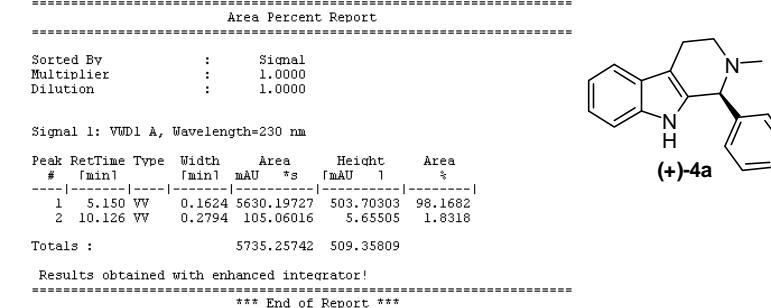
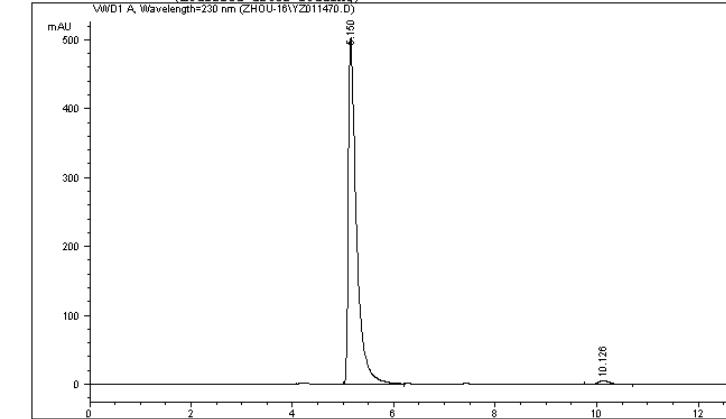
Instrument 1 17/12/2016 10:59:17

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Data File C:\HPCHEM\1\DATA\ZHOUE-16\Y2011470.D

IC, Hexane/iPrOH = 70/30, 0.8 mL/min, 30 oC, 230 nm

=====
Injection Date : 19/11/2016 10:26:50
Sample Name : YJ-8-25
Acq. Operator :
Acq. Method : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed : 19/11/2016 10:07:11 bv
(modified after loading)
Analysis Method : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed : 17/12/2016 11:00:49 bv
(modified after loading)



Instrument 1 17/12/2016 11:00:52

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