

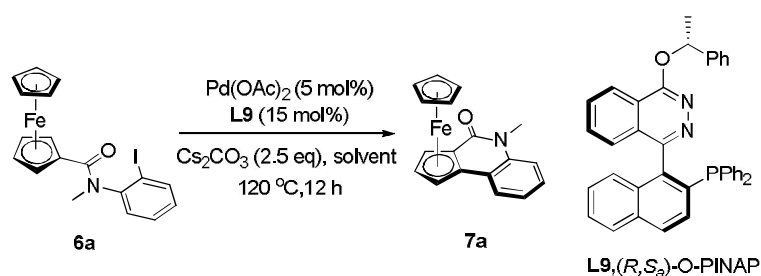
**Palladium-Catalyzed Intramolecular Cp-H Bond
Functionalization/Arylation: An Enantioselective Approach to
Planar Chiral Quinilinoferrocenes**

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Supporting Information

General Information. Proton nuclear magnetic resonance were recorded on 400MHz instruments internally referenced to tetramethylsilane (0.0 ppm) or residue of CDCl₃ (7.26 ppm) signals. All reactions were performed under an inert atmosphere of nitrogen in flame-dried glassware, unless otherwise stated. Solvents were distilled using standard techniques. Tetrahydrofuran was distilled over sodium under an atmosphere of nitrogen. Toluene, dichloromethane were distilled calcium hydride under an atmosphere of nitrogen.

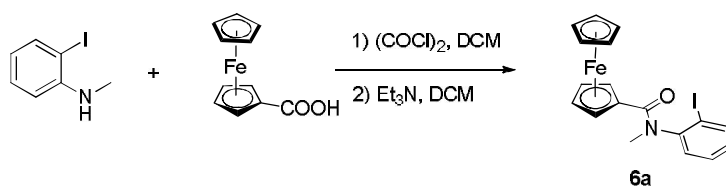
Table S1. Solvents Screening



solvent	yield	%ee
MeCN	NR	-
1,4-dioxane	30%	70
THF	NR	-
DMF	NR	-
toluene	98%	67
mesitylene ^a	43%	52
anisole	20%	50
benzene	98%	67
cyclohexane	20%	50

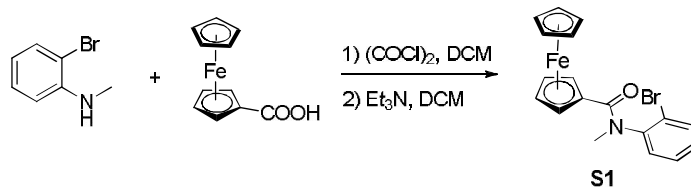
^a At 80 °C

Typical Procedures for the preparation of 6: (Typical Procedure A)



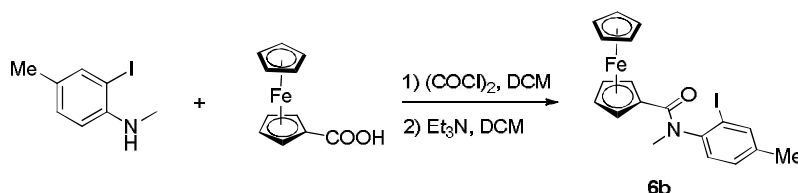
Oxalyl chloride (0.85 ml, 10.0 mmol) was added dropwise to a solution of ferrocenecarboxylic acid (1.15 g, 5.0 mmol) in CH_2Cl_2 (75 ml) at 25 °C under argon. The resulting solution was stirred at room temperature for 0.5 h and concentrated under reduced pressure. 2-Iodo-*N*-methylaniline (1.165 g, 5.0 mmol) and Et_3N (1.4 ml, 10.0 mmol) in CH_2Cl_2 (75 ml) was added to a solution of the above acid chloride in CH_2Cl_2 (125 ml) and stirred for 15 h. The reaction mixture was extracted with CH_2Cl_2 and the organic layer was washed with NaHCO_3 aq. and brine. The organic layer was dried over MgSO_4 , filtrated and concentrated under reduced pressure and the residue was purified by flash chromatography (hexanes/ EtOAc = 5:1) on silica gel to afford **6a** as a yellow solid (1.50 g, 66%). ¹H NMR (400 MHz, CDCl_3) δ 7.95 (dd, J = 8.0, 1.6 Hz, 1 H), 7.36 (td, J = 7.6, 1.2 Hz, 1 H), 7.18 (dd, J = 8.0, 1.6 Hz, 1 H), 7.07 (td, J = 7.6, 1.6 Hz, 1 H), 4.22 (bs, 1 H), 4.20 (s, 5 H), 4.16 (bs, 1 H), 4.08 (bs, 1 H), 3.80 (bs, 1 H), 3.30 (s, 3 H); ¹³C NMR (100 MHz, CDCl_3) δ 170.4, 147.1, 140.1, 129.74,

129.68, 129.5, 99.4, 75.4, 72.0, 70.3, 70.07, 70.03, 69.95, 37.7. HRMS (EI) calcd for $C_{18}H_{16}^{56}FeINO [M^+]$ 444.9626, found 444.9620.



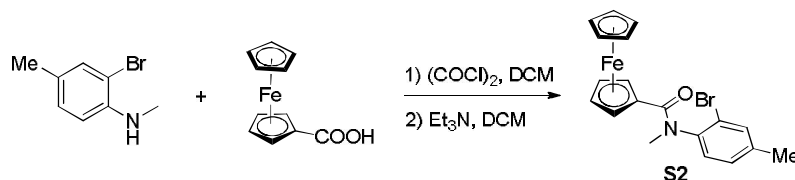
Compound **S1** was prepared following the Typical Procedure A.

The reaction of ferrocenecarboxylic acid (0.138 g, 0.60 mmol) and 2-bromo-*N*-methylaniline (0.111 g, 0.60 mmol) afforded **S1** (0.154 g, 65%). 1H NMR (400 MHz, $CDCl_3$) δ 7.70 (dd, $J = 8.0, 1.6$ Hz, 1 H), 7.33 (td, $J = 7.6, 1.6$ Hz, 1 H), 7.26 (td, $J = 7.6, 1.6$ Hz, 1 H), 7.19 (dd, $J = 8.0, 1.6$ Hz, 1 H), 4.29 (bs, 1 H), 4.19 (s, 5 H), 4.17 (bs, 1 H), 4.08 (bs, 1 H), 3.75 (bs, 1 H), 3.32 (s, 3 H); ^{13}C NMR (100 MHz, $CDCl_3$) δ 170.4, 143.6, 133.7, 130.5, 129.5, 128.7, 123.1, 75.3, 71.8, 70.0, 69.9, 37.3. HRMS (EI) calcd for $C_{18}H_{16}^{79}Br^{56}FeNO [M^+]$ 396.9765, found 396.9760.



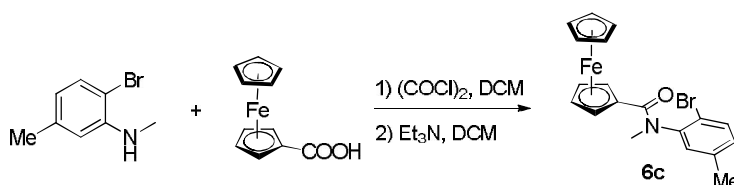
Compound **6b** was prepared following the Typical Procedure A.

The reaction of ferrocenecarboxylic acid (0.552 g, 2.4 mmol) and 2-iodo-*N*,4-dimethylaniline (0.593 g, 2.4 mmol) afforded **6b** (0.881 g, 80%). 1H NMR (400 MHz, $CDCl_3$) δ 7.77 (d, $J = 1.2$ Hz, 1 H), 7.15 (dd, $J = 8.0, 1.2$ Hz, 1 H), 7.04 (d, $J = 7.6$ Hz, 1 H), 4.21 (bs, 1 H), 4.20 (s, 5 H), 4.16 (bs, 1 H), 4.09 (bs, 1 H), 3.84 (bs, 1 H), 3.27 (s, 3 H), 2.36 (s, 3 H); ^{13}C NMR (100 MHz, $CDCl_3$) δ 170.4, 144.5, 140.4, 139.9, 130.4, 129.1, 99.2, 75.5, 71.9, 70.4, 70.03, 69.99, 69.92, 37.8, 20.6. HRMS (EI) calcd for $C_{19}H_{18}^{56}FeINO [M^+]$ 458.9782, found 458.9789.



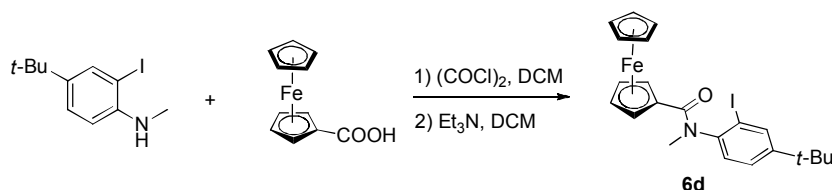
Compound **S2** was prepared following the Typical Procedure A.

The reaction of ferrocenecarboxylic acid (0.460 g, 2.0 mmol) and 2-bromo-*N*,4-dimethylaniline (0.494 g, 2.0 mmol) afforded **S2** (0.617 g, 75%). 1H NMR (400 MHz, $CDCl_3$) δ 7.52 (d, $J = 1.2$ Hz, 1 H), 7.11 (dd, $J = 8.0, 1.2$ Hz, 1 H), 7.06 (d, $J = 8.0$ Hz, 1 H), 4.28 (bs, 1 H), 4.18 (s, 5 H), 4.17 (bs, 1 H), 4.09 (bs, 1 H), 3.78 (bs, 1 H), 3.29 (s, 3 H), 2.38 (s, 3 H); ^{13}C NMR (100 MHz, $CDCl_3$) δ 170.4, 140.9, 139.9, 134.1, 129.9, 129.4, 122.7, 75.3, 71.8, 70.1, 70.0, 69.9, 37.4, 20.8. HRMS (EI) calcd for $C_{19}H_{18}^{79}Br^{56}FeNO [M^+]$ 410.9921, found 410.9914.



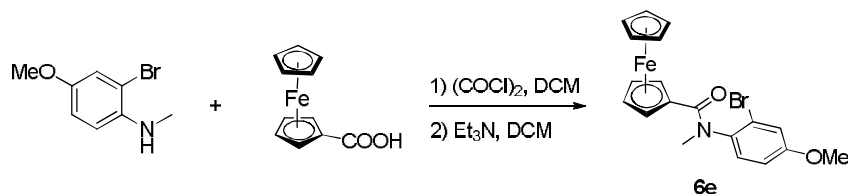
Compound **6c** was prepared following the Typical Procedure A.

The reaction of ferrocenecarboxylic acid (0.138 g, 0.60 mmol) and 2-bromo-*N*,5-dimethylaniline (0.119 g, 0.60 mmol) afforded **6c** (0.201 g, 82%). ¹H NMR (400 MHz, CDCl₃) δ 7.56 (d, *J* = 8.0 Hz, 1 H), 7.06 (d, *J* = 8.4 Hz, 1 H), 7.01 (s, 1 H), 4.33-4.03 (m, 8 H), 3.80 (bs, 1 H), 3.29 (s, 3 H), 2.30 (s, 3 H); ¹³C NMR (100 MHz, CDCl₃) δ 170.3, 143.3, 139.0, 133.3, 131.0, 130.4, 119.7, 75.4, 71.8, 70.2, 70.1, 69.9, 37.4, 20.8. HRMS (EI) calcd for C₁₉H₁₈⁷⁹Br⁵⁶FeNO [M⁺] 410.9921, found 410.9915.



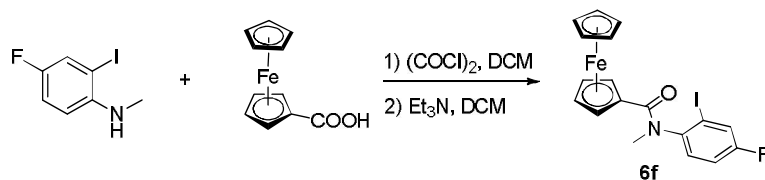
Compound **6d** was prepared following the Typical Procedure A.

The reaction of ferrocenecarboxylic acid (0.460 g, 2.0 mmol) and 4-(*tert*-butyl)-2-iodo-*N*-methylaniline (0.578 g, 2.0 mmol) afforded **6d** (0.731 g, 73%). ¹H NMR (400 MHz, CDCl₃) δ 7.91 (d, *J* = 2.0 Hz, 1 H), 7.35 (dd, *J* = 8.0, 2.0 Hz, 1 H), 7.07 (d, *J* = 8.0 Hz, 1 H), 4.29 (s, 1 H), 4.21 (s, 5 H), 4.17 (bs, 1 H), 4.08 (bs, 1 H), 3.75 (bs, 1 H), 3.28 (s, 3 H), 1.33 (s, 9 H); ¹³C NMR (100 MHz, CDCl₃) δ 170.5, 153.2, 144.4, 137.0, 128.9, 126.8, 99.2, 75.5, 72.0, 70.2, 70.1, 69.9, 37.8, 34.6, 31.2. HRMS (EI) calcd for C₂₂H₂₄⁵⁶FeINO [M⁺] 501.0252, found 501.0250.



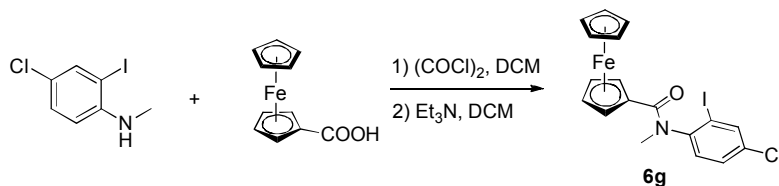
Compound **6e** was prepared following the Typical Procedure A.

The reaction of ferrocenecarboxylic acid (0.460 g, 2.0 mmol) and 2-bromo-4-methoxy-*N*-methylaniline (0.430 g, 2.0 mmol) afforded **6e** (0.599 g, 70%). ¹H NMR (400 MHz, CDCl₃) δ 7.22 (d, *J* = 2.8 Hz, 1 H), 7.10 (d, *J* = 8.8 Hz, 1 H), 6.85 (dd, *J* = 8.8, 2.8 Hz, 1 H), 4.26 (bs, 1 H), 4.21-4.15 (m, 6 H), 4.10 (bs, 1 H), 3.84 (s, 3 H), 3.81 (bs, 1 H), 3.28 (s, 3 H); ¹³C NMR (100 MHz, CDCl₃) δ 170.6, 159.5, 136.4, 130.7, 123.6, 118.5, 114.4, 75.4, 71.9, 70.2, 70.1, 70.0, 69.9, 55.8, 37.6. HRMS (EI) calcd for C₁₉H₁₈⁷⁹Br⁵⁶FeNO₂ [M⁺] 426.9870, found 426.9868.



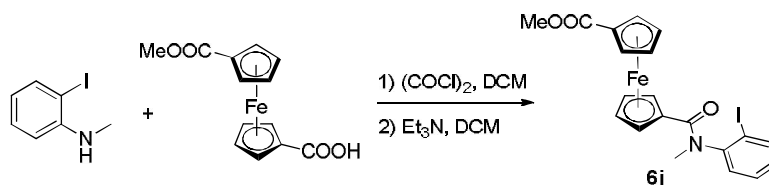
Compound **6f** was prepared following the Typical Procedure A.

The reaction of ferrocenecarboxylic acid (0.161 g, 0.70 mmol) and 4-fluoro-2-iodo-*N*-methylaniline (0.176 g, 0.70 mmol) afforded **6f** (0.253 g, 78%). ¹H NMR (400 MHz, CDCl₃) δ 7.66 (dd, *J* = 7.6, 2.8 Hz, 1 H), 7.17-7.04 (m, 2 H), 4.33-4.16 (m, 7 H), 4.12 (bs, 1 H), 3.79 (bs, 1 H), 3.27 (s, 3 H); ¹³C NMR (100 MHz, CDCl₃) δ 170.5, 162.3, 159.7, 143.6, 143.5, 130.3, 130.2, 126.9, 126.7, 116.8, 116.6, 99.3, 99.2, 75.3, 71.9, 70.2, 70.0, 37.8. HRMS (EI) calcd for C₁₈H₁₅F⁵⁶FeINO [M⁺] 462.9532, found 462.9524.



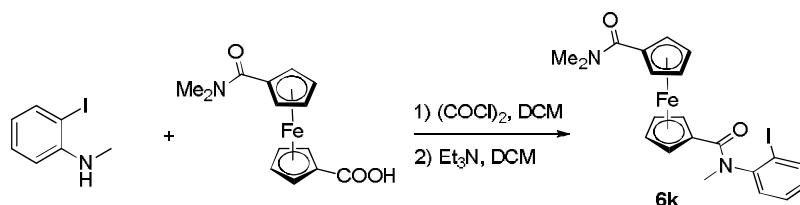
Compound **6g** was prepared following the Typical Procedure A.

The reaction of ferrocenecarboxylic acid (0.460 g, 2.0 mmol) and 2-iodo-4-chloro-*N*-methylaniline (0.534 g, 2.0 mmol) afforded **6g** (0.651 g, 68%). ¹H NMR (400 MHz, CDCl₃) δ 7.94 (d, *J* = 2.4 Hz, 1 H), 7.33 (dd, *J* = 8.4, 2.4 Hz, 1 H), 7.09 (d, *J* = 8.4 Hz, 1 H), 4.29 (bs, 1 H), 4.21 (s, 6 H), 4.13 (bs, 1 H), 3.82 (bs, 1 H), 3.27 (s, 3 H); ¹³C NMR (100 MHz, CDCl₃) δ 170.4, 145.9, 139.4, 134.2, 130.1, 129.8, 99.6, 75.2, 72.0, 70.2, 70.08, 70.05, 37.7. HRMS (EI) calcd for C₁₈H₁₅³⁵Cl⁵⁶FeINO [M⁺] 478.9236, found 478.9232.



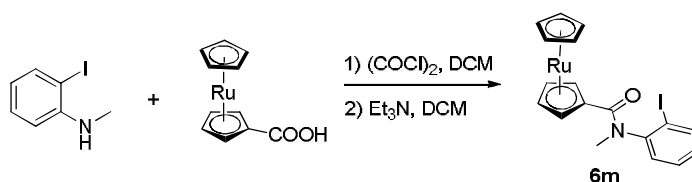
Compound **6j** was prepared following the Typical Procedure A.

The reaction of 1-methoxycarbonylferrocene-1-carboxylic acid (0.144 g, 0.50 mmol) and 2-iodo-*N*-methylaniline (0.117 g, 0.50 mmol) afforded **6j** (0.151 g, 60%). ¹H NMR (400 MHz, CDCl₃) δ 7.92 (dd, *J* = 7.6, 1.2 Hz, 1 H), 7.39 (td, *J* = 7.6, 1.2 Hz, 1 H), 7.28 (dd, *J* = 7.6, 1.2 Hz, 1 H), 7.07 (td, *J* = 8.0, 1.6 Hz, 1 H), 4.89-4.73 (m, 2 H), 4.49-4.39 (m, 2 H), 4.23-4.07 (m, 3 H), 3.98 (bs, 1 H), 3.80 (s, 3 H), 3.29 (s, 3 H); ¹³C NMR (100 MHz, CDCl₃) δ 171.3, 146.8, 140.1, 129.73, 129.68, 99.5, 73.61, 73.56, 72.44, 72.40, 72.0, 71.8, 71.7, 71.6, 71.2, 51.7, 37.6. HRMS (EI) calcd for C₂₀H₁₈⁵⁶FeINO₃ [M⁺] 502.9681, found 502.9673.



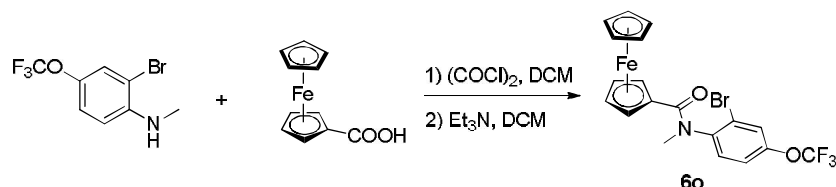
Compound **6k** was prepared following the Typical Procedure A.

The reaction of 1-carboxy-1'-[(dimethylamino) carbonyl] ferrocene (60.0 mg, 0.20 mmol) and 2-iodo-*N*-methylaniline (47.0 mg, 0.20 mmol) afforded **6k** (45.0 mg, 44%). ¹H NMR (400 MHz, CDCl₃) δ 7.91 (dd, *J* = 8.0, 1.2 Hz, 1 H), 7.37 (td, *J* = 7.2, 1.2 Hz, 3 H), 7.24 (dd, *J* = 8.0, 1.2 Hz, 1 H), 7.06 (td, *J* = 8.0, 1.6 Hz, 1 H), 4.66 (d, *J* = 12.4 Hz, 2 H), 4.38 (d, *J* = 10.8 Hz, 2 H), 4.28-4.08 (m, 3 H), 4.01 (bs, 1 H), 3.28 (s, 3 H), 3.23-2.91 (br, 6 H); ¹³C NMR (100 MHz, CDCl₃) δ 170.1, 169.8, 146.7, 140.0, 129.7, 129.6, 99.2, 79.2, 76.7, 72.9, 72.4, 72.3, 72.2, 72.1, 72.0, 71.9, 71.7, 39.0, 37.7, 36.6. HRMS (EI) calcd for C₂₁H₂₁⁵⁶FeIN₂O₂ [*M*⁺] 515.9997, found 515.9991.



Compound **6m** was prepared following the Typical Procedure A.

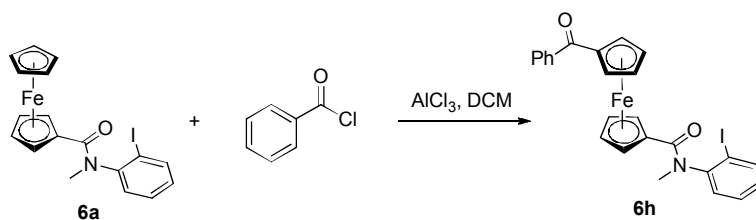
The reaction of ruthenocenecarboxylic acid (0.275 g, 1.0 mmol) and 2-iodo-*N*-methylaniline (0.233 g, 1.0 mmol) afforded **6m** (0.368 g, 75%). ¹H NMR (400 MHz, CDCl₃) δ 7.93 (d, *J* = 8.0 Hz, 1 H), 7.40 (td, *J* = 8.0, 1.6 Hz, 1 H), 7.24 (dd, *J* = 8.0, 1.2 Hz, 1 H), 7.09 (td, *J* = 8.0, 1.2 Hz, 1 H), 4.57 (s, 5 H), 4.47 (bs, 1 H), 4.43 (bs, 1 H), 4.32 (bs, 1 H), 4.28 (bs, 1 H), 3.24 (s, 3 H); ¹³C NMR (100 MHz, CDCl₃) δ 168.8, 146.8, 140.1, 129.71, 129.65, 129.58, 99.5, 79.7, 72.9, 72.4, 72.0, 71.9, 71.7, 37.7. HRMS (EI) calcd for C₁₈H₁₆¹⁰²RuINO [*M*⁺] 490.9320, found 490.9315.



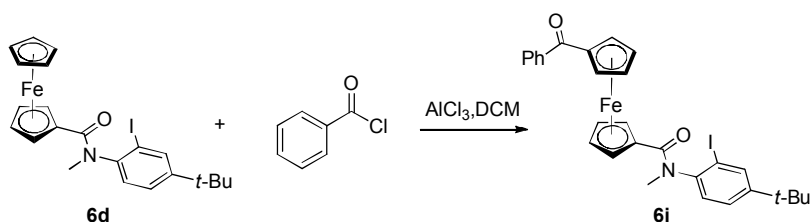
Compound **6o** was prepared following the Typical Procedure A.

The reaction of ferrocenecarboxylic acid (0.138 g, 0.60 mmol) and 2-bromo-*N*-methyl-4-(trifluoromethoxy)aniline (0.162 g, 0.60 mmol) afforded **6o** (87.0 mg, 30%). ¹H NMR (400 MHz, CDCl₃) δ 7.58 (d, *J* = 1.6 Hz, 1 H), 7.24-7.15 (m, 2 H), 4.35 (bs, 1 H), 4.26-4.16 (m, 6 H), 4.12 (bs, 1 H), 3.74 (bs, 1 H), 3.30 (s, 3 H); ¹³C NMR (100 MHz, CDCl₃) δ 170.5, 148.4, 142.5, 131.2, 126.1, 123.7, 121.5, 121.0, 118.9, 116.4, 75.2, 71.9, 70.3, 70.1, 69.9, 37.4. HRMS (EI) calcd for C₁₉H₁₅⁷⁹BrF₃⁵⁶FeNO₂ [*M*⁺] 480.9588, found 480.9591.

Typical Procedures for the preparation of 6h, 6i, 6l: (Typical Procedure B)

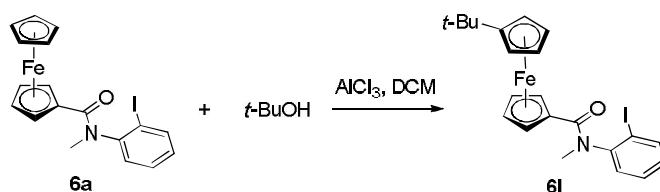


The benzoyl chloride (0.14 ml, 1.2 mmol, 1.2 equiv) was added to a mixture of **6a** (0.445 g, 1.0 mmol) and AlCl_3 (0.267 g, 2.0 mmol, 2.0 equiv) in CH_2Cl_2 (10 ml) at room temperature and stirred for additional 3 h. The reaction was quenched by the addition of crushed ice and extracted with CH_2Cl_2 twice. The combined organic phase was washed with water and dried over Na_2SO_4 . After filtration the solvent was removed by evaporation, and the resulting mixture was purified by column chromatography on silica gel (hexanes/ethyl acetate = 1:1) to afford the desired product **6h** (0.480 g, 85%). ^1H NMR (400 MHz, CDCl_3) δ 7.96-7.80 (m, 3 H), 7.56 (t, J = 7.2 Hz, 1 H), 7.46 (t, J = 8.0 Hz, 2 H), 7.33 (t, J = 7.2 Hz, 1 H), 7.17 (dd, J = 7.6, 1.2 Hz, 1 H), 7.06 (t, J = 7.6 Hz, 1 H), 4.93 (s, 2 H), 4.66 (s, 1 H), 4.65 (s, 1 H), 4.22-3.97 (m, 4 H), 3.27 (s, 3 H); ^{13}C NMR (100 MHz, CDCl_3) δ 198.5, 168.9, 146.6, 140.1, 139.3, 131.7, 129.75, 129.69, 128.2, 99.4, 79.0, 75.2, 73.2, 73.05, 72.98, 72.61, 72.56, 72.4, 72.0, 37.7. HRMS (EI) calcd for $\text{C}_{25}\text{H}_{20}^{56}\text{FeINO}_2$ [M^+] 548.9888, found 548.9895.



Compound **6i** was prepared following the **Typical Procedure B**.

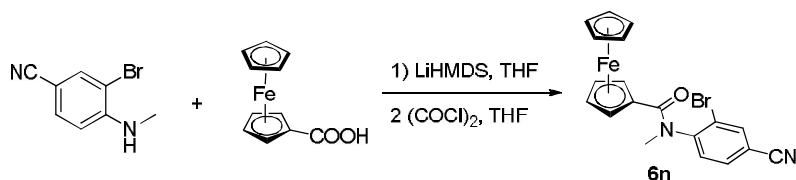
The reaction of benzoyl chloride (70 μl , 0.60 mmol, 1.2 equiv) and **6d** (0.251 g, 0.50 mmol) afforded **6i** (0.242 g, 80%). ^1H NMR (400 MHz, CDCl_3) δ 8.04-7.74 (m, 3 H), 7.64-7.31 (m, 4 H), 7.06 (bs, 1 H), 4.92 (bs, 2 H), 4.66 (bs, 2 H), 4.25-3.85 (m, 4 H), 3.32 (br, 3 H), 1.30 (s, 9 H). ^{13}C NMR (100 MHz, CDCl_3) δ 198.5, 169.1, 153.5, 143.8, 139.4, 136.9, 131.7, 128.9, 128.2, 126.9, 99.2, 78.9, 75.2, 73.2, 73.0, 72.7, 72.4, 72.1, 37.7, 34.6, 31.1. HRMS (EI) calcd for $\text{C}_{29}\text{H}_{28}^{56}\text{FeINO}_2$ [M^+] 605.0514, found 605.0511.



Compound **6l** was prepared following the **Typical Procedure B**.

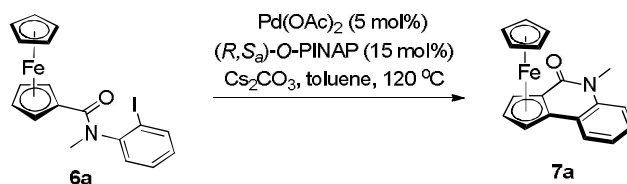
The reaction of *t*-BuOH (16.3 mg, 0.22 mmol, 1.1 equiv) and **6a** (89.0 mg, 0.20

mmol) afforded **6l** (40.1 mg, 40%). ¹H NMR (400 MHz, CDCl₃) δ 7.94 (d, *J* = 8.0 Hz, 1 H), 7.36 (t, *J* = 7.2 Hz, 1 H), 7.18 (d, *J* = 6.8 Hz, 1 H), 7.07 (t, *J* = 7.6 Hz, 1 H), 4.25-4.11 (m, 4 H), 4.10-4.02 (m, 3 H), 3.87 (bs, 1 H), 3.29 (s, 3 H), 1.17 (s, 9 H); ¹³C NMR (100 MHz, CDCl₃) δ 170.6, 147.2, 140.1, 129.7, 129.6, 129.4, 103.2, 99.4, 75.5, 71.8, 70.6, 70.5, 70.04, 69.98, 66.9, 66.4, 37.7, 31.4, 30.2. HRMS (EI) calcd for C₂₂H₂₄⁵⁶FeINO₂ [M⁺] 501.0252, found 501.0250.



The LiHMDS (1.0 M in THF) (10.0 ml, 10.0 mmol) was added to 3-bromo-4-(methylamino)benzonitrile (0.422 g, 2.0 mmol) in THF (20 ml) at -10 °C. After stirring for 20 min at -10 °C the reaction mixture was allowed to warm to room temperature and stirred for additional 2 h.

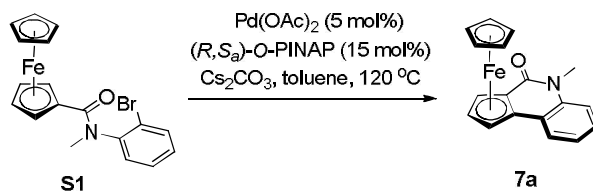
Oxalyl chloride (0.43 ml, 5.0 mmol) was added dropwise to a solution of ferrocenecarboxylic acid (0.576 g, 2.5 mmol) in CH₂Cl₂ (20 ml) at 25 °C under argon. The resulting solution was stirred at room temperature for 0.5 h and concentrated under reduced pressure. The residue was diluted with THF (12 ml). This solution was added to the above amide solution at 0 °C in 25 min then the reaction mixture was allowed to warm to room temperature and stirred for another 20 min. After quenched with water, the residue was extracted with ethyl acetate twice. The combined organic layer was washed with brine, dried over Na₂SO₄, filtrated and concentrated, and purified by column chromatography (hexanes/EtOAc = 4:1) on silica gel to afford **6n** (0.558 g, 66%). ¹H NMR (400 MHz, CDCl₃) δ 7.99 (d, *J* = 1.6 Hz, 1 H), 7.59 (dd, *J* = 8.0, 1.6 Hz, 1H), 7.26 (d, *J* = 8.0 Hz, 1H), 4.41 (bs, 1 H), 4.32-3.95 (m, 7 H), 3.73 (bs, 1 H), 3.31 (s, 3 H); ¹³C NMR (100 MHz, CDCl₃) δ 170.4, 148.3, 137.3, 132.3, 131.3, 123.7, 116.6, 113.1, 75.1, 71.9, 70.3, 70.2, 69.8, 37.2. HRMS (EI) calcd for C₁₉H₁₅⁷⁹Br⁵⁶FeN₂O [M⁺] 421.9717, found 421.9709.



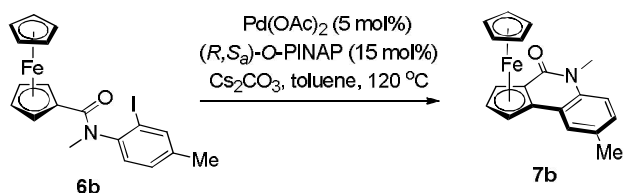
Typical procedure for the palladium-catalyzed asymmetric C-H functionalization/cyclization reaction: (**Typical Procedure C**).

A mixture of **6a** (44.5 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)₂ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S_a*)-O-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs₂CO₃ (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) was stirred at 120 °C for 12 h (the reaction was monitored by TLC). After cooling to room temperature the solvent was removed by evaporation and the residue was purified by column chromatography on silica gel (hexanes/ethyl acetate = 5:1) to afford **7a** (31.0 mg, 98%, 67% ee). [α]_D²³

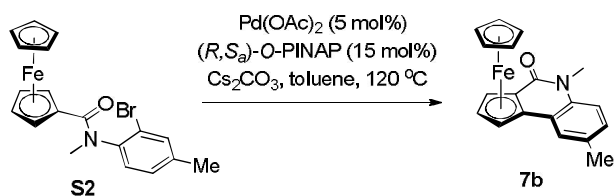
+19.6 (c 1.0, CHCl₃). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 8:92, flow: 1.0 ml/min, λ = 254 nm. **¹H NMR** (400 MHz, CDCl₃) δ 7.72 (d, J = 7.6 Hz, 1 H), 7.39 (t, J = 7.6 Hz, 1 H), 7.27 (d, J = 7.6 Hz, 1 H), 7.18 (t, J = 7.6 Hz, 1 H), 5.22 (s, 1 H), 5.15 (s, 1 H), 4.49 (s, 1 H), 3.94 (s, 5 H), 3.70 (s, 3 H); **¹³C NMR** (100 MHz, CDCl₃) δ 167.8, 138.2, 127.2, 123.5, 123.0, 122.3, 114.9, 83.4, 71.5, 71.3, 70.4, 67.2, 63.6, 29.3. HRMS (EI) calcd for C₁₈H₁₅⁵⁶FeNO [M⁺] 317.0503, found 317.0497.



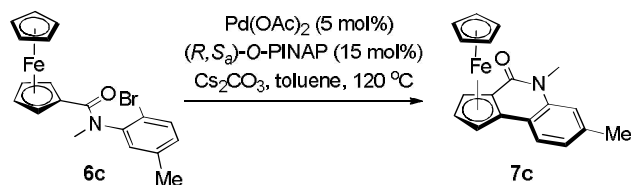
The reaction of **S1** (39.8 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)₂ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S_a*)-O-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs₂CO₃ (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7a** (29.8 mg, 94%, 49% ee).



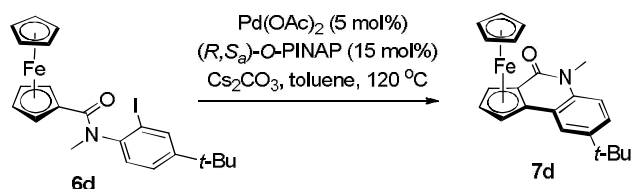
The reaction of **6b** (45.9 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)₂ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S_a*)-O-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs₂CO₃ (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7b** (R_f = 0.30, PE:EA = 5:1) (29.8 mg, 90%, 67% ee). [α]_D²³ +185.0 (c 1.0, CHCl₃). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 6:94, flow: 1.0 ml/min, λ = 254 nm. **¹H NMR** (400 MHz, CDCl₃) δ 7.51 (s, 1 H), 7.23-7.11 (m, 2 H), 5.20 (s, 1 H), 5.12 (s, 1 H), 4.47 (s, 1 H), 3.94 (s, 5 H), 3.67 (s, 3 H), 2.44 (s, 3 H); **¹³C NMR** (100 MHz, CDCl₃) δ 167.7, 136.1, 131.7, 128.1, 123.7, 122.8, 114.9, 83.5, 71.6, 71.2, 70.4, 67.1, 63.5, 29.3, 20.7. HRMS (EI) calcd for C₁₉H₁₇⁵⁶FeNO [M⁺] 331.0660, found 331.0656.



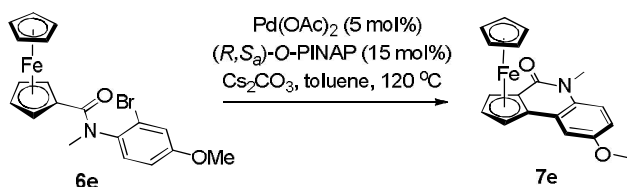
The reaction of **S2** (41.1 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)₂ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S_a*)-O-PINAP (8.4 mg, 0.015 mmol, 15 mol %) and Cs₂CO₃ (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7b** (31.4 mg, 95%, 57% ee).



The reaction of **6c** (41.1 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)_2 (1.1 mg, 0.005 mmol, 5 mol%), $(R,S_a)\text{-O-PINAP}$ (8.4 mg, 0.015 mmol, 15 mol%) and Cs_2CO_3 (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7c** (R_f = 0.20, PE:EA = 5:1) (32.1 mg, 97%, 51% ee). $[\alpha]^{23}_D$ +6.4 (c 1.0, CHCl_3). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 6:94, flow: 1.0 ml/min, λ = 254 nm. **$^1\text{H NMR}$** (400 MHz, CDCl_3) δ 7.60 (d, J = 7.6 Hz, 1 H), 7.08 (s, 1 H), 7.00 (d, J = 7.6 Hz, 1 H), 5.19 (dd, J = 2.4, 1.2 Hz, 1 H), 5.11 (dd, J = 2.4, 1.2 Hz, 1 H), 4.46 (t, J = 2.4 Hz, 1 H), 3.93 (s, 5 H), 3.68 (s, 3 H), 2.46 (s, 3 H); **$^{13}\text{C NMR}$** (100 MHz, CDCl_3) δ 168.1, 138.2, 137.2, 123.3, 123.2, 120.0, 115.6, 83.8, 71.15, 71.06, 70.3, 66.9, 63.3, 29.3, 22.0. HRMS (EI) calcd for $\text{C}_{19}\text{H}_{17}^{56}\text{FeNO}$ [M^+] 331.0660, found 331.0655.

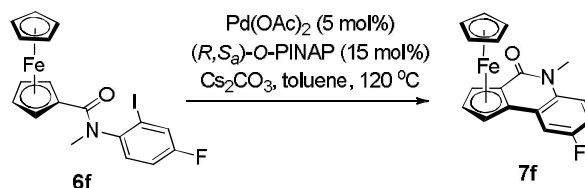


The reaction of **6d** (50.1 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)_2 (1.1 mg, 0.005 mmol, 5 mol%), $(R,S_a)\text{-O-PINAP}$ (8.4 mg, 0.015 mmol, 15 mol%) and Cs_2CO_3 (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7d** (R_f = 0.30, PE:EA = 5:1) (32.1 mg, 86%, 60% ee). $[\alpha]^{23}_D$ +293.8 (c 1.0, CHCl_3). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 0.5:99.5, flow: 1.0 ml/min, λ = 254 nm. **$^1\text{H NMR}$** (400 MHz, CDCl_3) δ 7.70 (d, J = 2.4 Hz, 1 H), 7.42 (dd, J = 8.8, 2.4 Hz, 1 H), 7.21 (d, J = 8.8 Hz, 1 H), 5.24-5.18 (m, 1 H), 5.18-5.13 (m, 1 H), 4.48 (t, J = 2.8 Hz, 1 H), 3.94 (s, 5 H), 3.68 (s, 3 H), 1.41 (s, 9 H); **$^{13}\text{C NMR}$** (100 MHz, CDCl_3) δ 167.8, 145.2, 136.0, 124.5, 122.4, 119.9, 114.7, 83.8, 71.6, 71.1, 70.4, 67.1, 63.4, 34.3, 31.5, 29.2. HRMS (EI) calcd for $\text{C}_{22}\text{H}_{23}^{56}\text{FeNO}$ [M^+] 373.1129, found 373.1126.

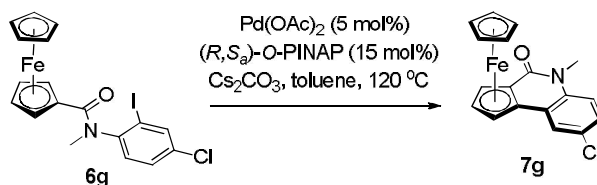


The reaction of **6e** (42.8 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)_2 (1.1 mg, 0.005 mmol, 5 mol%), $(R,S_a)\text{-O-PINAP}$ (8.4 mg, 0.015 mmol, 15 mol%) and Cs_2CO_3 (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7e** (R_f = 0.30, PE:EA = 3:1) (31.2 mg, 90%, 64% ee). $[\alpha]^{23}_D$ +198.9 (c 1.0, CHCl_3). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 6:94, flow: 1.0 ml/min, λ = 254 nm. **$^1\text{H NMR}$** (400 MHz, CDCl_3) δ 7.22 (d, J = 2.8 Hz, 1 H), 7.20 (d, J = 9.2 Hz, 1 H),

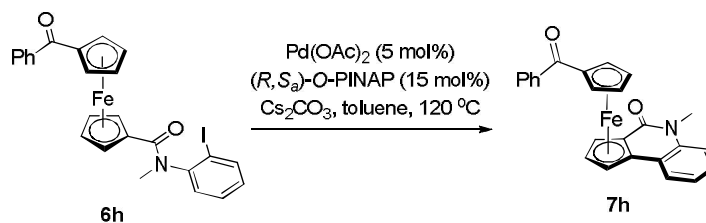
6.96 (dd, $J = 9.2, 2.8$ Hz, 1 H), 5.21 (dd, $J = 2.4, 1.2$ Hz, 1 H), 5.10 (dd, $J = 2.4, 1.2$ Hz, 1 H), 4.48 (t, $J = 2.4$ Hz, 1 H), 3.95 (s, 5 H), 3.91 (s, 3 H), 3.67 (s, 3 H); ^{13}C NMR (100 MHz, CDCl_3) δ 167.3, 154.9, 132.3, 124.2, 116.0, 112.8, 108.2, 83.0, 71.8, 71.3, 70.4, 67.3, 63.7, 55.6, 29.4. HRMS (EI) calcd for $\text{C}_{19}\text{H}_{17}^{56}\text{FeNO}_2$ [M^+] 347.0609, found 347.0610.



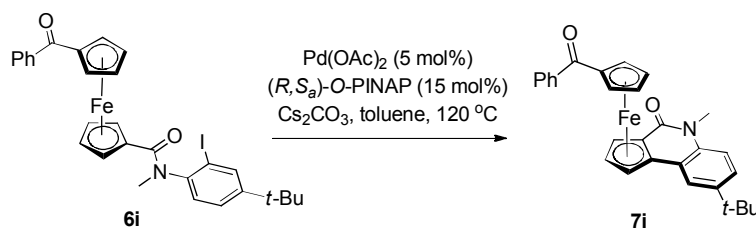
The reaction of **6f** (46.3 mg, 0.10 mmol, 1.0 equiv), $\text{Pd}(\text{OAc})_2$ (1.1 mg, 0.005 mmol, 5 mol%), (R,S_a)-*O*-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs_2CO_3 (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7f** ($R_f = 0.30$, PE:EA = 5:1) (28.5 mg, 85%, 45% ee). $[\alpha]_D^{23}$ -67.0 (c 1.0, CHCl_3). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 8:92, flow: 1.0 ml/min, $\lambda = 254$ nm. ^1H NMR (400 MHz, CDCl_3) δ 7.39 (dd, $J = 8.8, 3.2$ Hz, 1 H), 7.21 (dd, $J = 8.8, 4.8$ Hz, 1 H), 7.09 (td, $J = 8.8, 2.8$ Hz, 1 H), 5.23 (d, $J = 1.2$ Hz, 1 H), 5.10 (d, $J = 1.6$ Hz, 1 H), 4.52 (t, $J = 2.8$ Hz, 1 H), 3.97 (s, 5 H), 3.68 (s, 3 H); ^{13}C NMR (100 MHz, CDCl_3) δ 167.4, 159.6, 157.2, 134.5, 124.94, 124.86, 116.3, 116.2, 114.0, 113.8, 109.7, 109.5, 82.2, 71.65, 71.62, 70.5, 67.6, 64.0, 29.5. HRMS (EI) calcd for $\text{C}_{18}\text{H}_{14}\text{F}^{56}\text{FeNO}$ [M^+] 335.0409, found 335.0402.



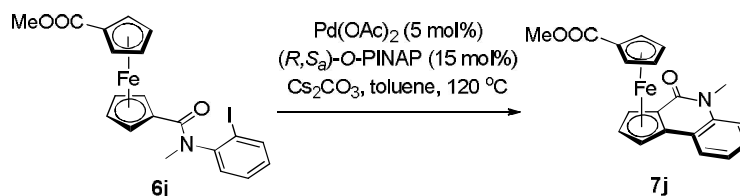
The reaction of **6g** (49.4 mg, 0.10 mmol, 1.0 equiv), $\text{Pd}(\text{OAc})_2$ (1.1 mg, 0.005 mmol, 5 mol%), (R,S_a)-*O*-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs_2CO_3 (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7g** ($R_f = 0.30$, PE:EA = 3:1) (33.0 mg, 94%, 28% ee). $[\alpha]_D^{23}$ +93.0 (c 1.0, CHCl_3). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 6:94, flow: 1.0 ml/min, $\lambda = 254$ nm. ^1H NMR (400 MHz, CDCl_3) δ 7.66 (d, $J = 2.4$ Hz, 1 H), 7.33 (dd, $J = 8.8, 2.4$ Hz, 1 H), 7.18 (d, $J = 8.8$ Hz, 1 H), 5.23 (s, 1 H), 5.11 (s, 1 H), 4.52 (s, 1 H), 3.97 (s, 5 H), 3.67 (s, 3 H); ^{13}C NMR (100 MHz, CDCl_3) δ 167.5, 136.8, 127.7, 126.9, 124.9, 122.9, 116.2, 82.0, 71.7, 71.5, 70.6, 67.6, 63.9, 29.4. HRMS (EI) calcd for $\text{C}_{18}\text{H}_{14}^{35}\text{Cl}^{56}\text{FeNO}$ [M^+] 351.0113, found 351.0109.



The reaction of **6h** (54.9 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)₂ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S*_a)-*O*-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs₂CO₃ (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7h** (R_f = 0.20, PE:EA = 1:1) (21.1 mg, 50%, 61% ee). [α]_D²³ -465.0 (c 1.0, CHCl₃). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 8:92, flow: 1.0 ml/min, λ = 254 nm. ¹H NMR (400 MHz, CDCl₃) δ 7.62-7.48 (m, 4 H), 7.46-7.33 (m, 3 H), 7.22 (d, *J* = 8.4 Hz, 1 H), 7.15 (t, *J* = 7.6 Hz, 1 H), 5.23 (s, 1 H), 5.12 (s, 1 H), 4.85 (s, 1 H), 4.55 (s, 1 H), 4.47 (s, 1 H), 4.45 (s, 1 H), 4.31 (s, 1 H), 3.58 (s, 3 H); ¹³C NMR (100 MHz, CDCl₃) δ 196.7, 166.5, 138.9, 138.4, 131.7, 128.3, 128.00, 127.96, 124.3, 122.4, 120.5, 115.2, 85.7, 79.4, 74.6, 74.0, 73.6, 72.9, 72.5, 72.2, 68.0, 65.5, 29.3. HRMS (EI) calcd for C₂₅H₁₉⁵⁶FeNO₂ [M⁺] 421.0765, found 421.0767.

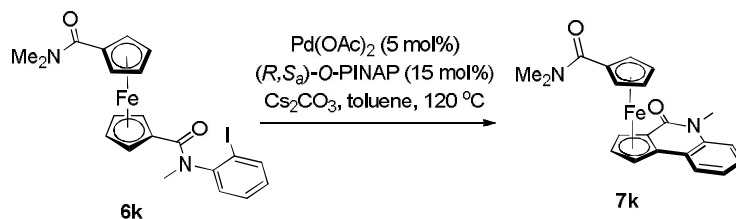


The reaction of **6i** (60.5 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)₂ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S*_a)-*O*-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs₂CO₃ (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7i** (R_f = 0.20, PE:EA = 2:1) (35.8 mg, 75%, 63% ee). [α]_D²³ -242.1 (c 1.0, CHCl₃). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 6:94, flow: 1.0 ml/min, λ = 254 nm. ¹H NMR (400 MHz, CDCl₃) δ 7.61-7.52 (m, 3 H), 7.52-7.42 (m, 2 H), 7.37 (t, *J* = 7.2 Hz, 2 H), 7.16 (d, *J* = 8.4 Hz, 1 H), 5.22 (s, 1 H), 5.16 (s, 1 H), 4.86 (s, 1 H), 4.56 (s, 1 H), 4.46 (s, 1 H), 4.43 (s, 1 H), 4.30 (s, 1 H), 3.57 (s, 3 H), 1.37 (s, 9 H); ¹³C NMR (100 MHz, CDCl₃) δ 196.7, 166.5, 145.3, 139.0, 136.2, 131.7, 128.3, 128.0, 125.4, 120.7, 119.9, 115.0, 86.3, 79.5, 74.5, 73.9, 73.6, 72.8, 72.6, 72.3, 68.0, 65.2, 34.2, 31.4, 29.2. HRMS (EI) calcd for C₂₉H₂₇⁵⁶FeNO₂ [M⁺] 477.1391, found 477.1389.

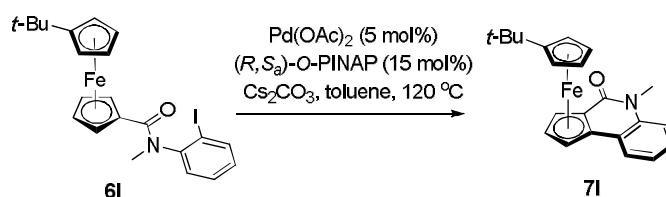


The reaction of **6j** (50.3 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)₂ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S*_a)-*O*-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs₂CO₃ (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7j** (R_f = 0.40, PE:EA = 2:1) (30.0 mg, 80%, 55% ee). [α]_D²³ -284.5 (c 1.0, CHCl₃). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 6:94, flow: 1.0 ml/min, λ = 254 nm. ¹H NMR (400 MHz, CDCl₃) δ 7.65 (dd, *J* = 7.6, 1.6 Hz, 1 H), 7.45 (td, *J* = 7.6, 1.6 Hz, 1 H), 7.29 (d, *J* = 8.4 Hz, 1 H), 7.20 (td, *J* = 7.6, 1.2 Hz, 1 H), 5.24 (dd, *J* = 2.4, 1.2 Hz, 1 H), 5.14 (dd, *J* = 2.4, 1.2 Hz, 1 H), 4.65 (dt, *J* = 2.4, 1.2 Hz, 1 H), 4.49 (t, *J* = 2.4 Hz, 1 H), 4.44-4.35 (m, 2 H), 4.31 (td, *J* = 2.4, 1.2 Hz, 1 H), 3.68 (s, 3 H), 3.34

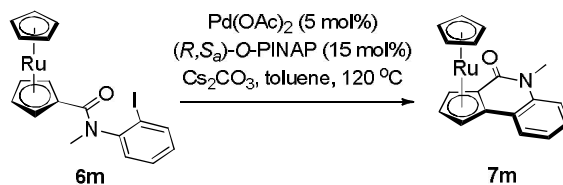
(s, 3 H); ^{13}C NMR (100 MHz, CDCl_3) δ 168.8, 166.6, 138.6, 127.9, 124.2, 122.2, 120.2, 115.1, 85.4, 74.1, 72.9, 72.8, 72.2, 71.9, 71.2, 67.8, 64.5, 51.6, 29.4. HRMS (EI) calcd for $\text{C}_{20}\text{H}_{17}^{56}\text{FeNO}_3$ [M^+] 375.0558, found 375.0550.



The reaction of **6k** (51.6 mg, 0.10 mmol, 1.0 equiv), $\text{Pd}(\text{OAc})_2$ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S_a*)-*O*-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs_2CO_3 (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7k** (R_f = 0.20, ethyl acetate) (29.1 mg, 75%, 53% ee). $[\alpha]^{23}_{\text{D}}$ +31.3 (c 1.0, CHCl_3). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 6:94, flow: 1.0 ml/min, λ = 254 nm. ^1H NMR (400 MHz, CDCl_3) δ 7.73 (dd, J = 7.6, 1.6 Hz, 1 H), 7.42 (td, J = 7.2, 1.6 Hz, 1 H), 7.27 (d, J = 7.6 Hz, 1 H), 7.19 (td, J = 7.6, 0.8 Hz, 1 H), 5.23 (dd, J = 2.4, 1.2 Hz, 1 H), 5.21 (dd, J = 2.8, 1.2 Hz, 1 H), 4.61 (dt, J = 2.4, 1.2 Hz, 1 H), 4.58 (t, J = 2.4 Hz, 1 H), 4.23 (dt, J = 2.4, 1.2 Hz, 1 H), 4.16 (td, J = 2.4, 1.2 Hz, 1 H), 4.02 (td, J = 2.4, 1.2 Hz, 1 H), 3.69 (s, 3 H), 2.88 (br, 6 H); ^{13}C NMR (100 MHz, CDCl_3) δ 168.3, 167.2, 138.4, 127.6, 124.2, 122.2, 121.5, 114.9, 84.7, 79.8, 73.5, 73.1, 72.3, 72.0, 71.9, 71.3, 67.9, 66.5, 38.8, 36.6, 29.4. HRMS (EI) calcd for $\text{C}_{21}\text{H}_{20}^{56}\text{FeN}_2\text{O}_2$ [M^+] 388.0874, found 388.0871.

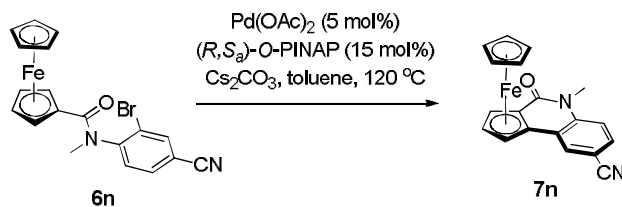


The reaction of **6l** (50.1 mg, 0.10 mmol, 1.0 equiv), $\text{Pd}(\text{OAc})_2$ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S_a*)-*O*-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs_2CO_3 (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7l** (R_f = 0.30, PE:EA = 2:1) (19.4 mg, 52%, 60% ee). $[\alpha]^{23}_{\text{D}}$ -26.0 (c 1.0, CHCl_3). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 8:92, flow: 1.0 ml/min, λ = 254 nm. ^1H NMR (400 MHz, CDCl_3) δ 7.71 (dd, J = 7.6, 1.2 Hz, 1 H), 7.38 (t, J = 7.6, 1 H), 7.26 (d, J = 7.6 Hz, 1 H), 7.17 (t, J = 7.2 Hz, 1 H), 5.20 (d, J = 1.2 Hz, 1 H), 5.12 (s, 1 H), 4.48 (t, J = 2.4 Hz, 1 H), 3.98 (s, 1 H), 3.77 (s, 1 H), 3.74 (s, 1 H), 3.69 (s, 3 H), 3.61 (s, 1 H), 1.18 (s, 9 H); ^{13}C NMR (100 MHz, CDCl_3) δ 167.9, 138.2, 127.1, 123.6, 123.0, 122.2, 114.9, 103.3, 83.3, 71.9, 71.3, 70.8, 70.5, 67.5, 67.1, 66.5, 64.1, 31.3, 30.2, 29.3. HRMS (EI) calcd for $\text{C}_{22}\text{H}_{23}^{56}\text{FeNO}$ [M^+] 373.1129, found 373.1126.

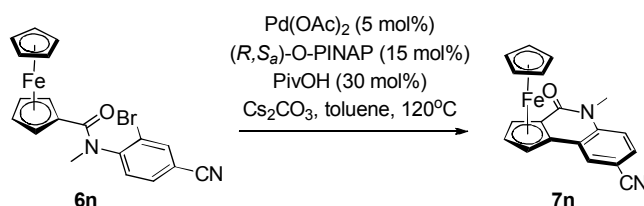


The reaction of **6m** (49.0 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)₂ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S_a*)-O-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs₂CO₃ (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7m** (R_f = 0.20, PE:EA = 5:1) (35.5 mg, 98%, 65% ee). [α]_D²³ -22.9 (c 1.0, CHCl₃). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 1:99, flow: 1.0 ml/min, λ = 254 nm. ¹H NMR (400 MHz, CDCl₃) δ 7.64 (dd, *J* = 7.6, 1.6 Hz, 1 H), 7.34 (td, *J* = 7.6, 1.6 Hz, 1 H), 7.22 (d, *J* = 8.0 Hz, 1 H), 7.12 (td, *J* = 7.6, 0.8 Hz, 1 H), 5.51 (dd, *J* = 2.4, 0.8 Hz, 2 H), 4.84 (t, *J* = 2.4 Hz, 1 H), 4.38 (s, 5 H), 3.62 (s, 3 H); ¹³C NMR (100 MHz, CDCl₃) δ 166.4, 137.8, 127.1, 123.2, 122.1, 121.8, 114.8, 86.6, 76.2, 72.8, 72.1, 69.3, 66.3, 29.3. HRMS (EI) calcd for C₁₈H₁₅¹⁰²RuNO [M⁺] 363.0197, found 363.0195.

Scheme 3. The effects by the addition of PivOH

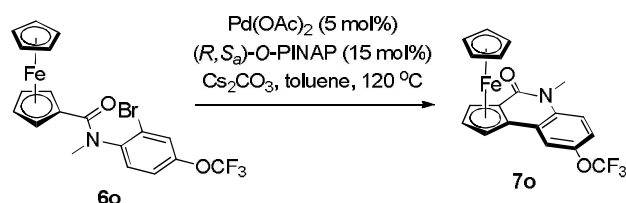


The reaction of **6n** (42.3 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)₂ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S_a*)-O-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs₂CO₃ (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7n** (R_f = 0.20, PE:EA = 2:1) (14.4 mg, 42%, 88% ee). [α]_D²³ +197.8 (c 1.0, CHCl₃). HPLC conditions: Chiralcel OJ-H, isopropanol/hexane = 15:85, flow: 1.0 ml/min, λ = 254 nm. ¹H NMR (400 MHz, CDCl₃) δ 7.97 (d, *J* = 2.0 Hz, 1 H), 7.65 (dd, *J* = 8.8, 1.6 Hz, 1 H), 7.31 (d, *J* = 8.8 Hz, 1 H), 5.26 (d, *J* = 1.6 Hz, 1 H), 5.18 (d, *J* = 1.6 Hz, 1 H), 4.59 (t, *J* = 2.4 Hz, 1 H), 3.99 (s, 5 H), 3.70 (s, 3 H); ¹³C NMR (100 MHz, CDCl₃) δ 167.7, 141.3, 130.6, 127.1, 124.4, 118.9, 115.4, 105.5, 81.2, 72.3, 71.1, 70.7, 67.9, 64.2, 29.5. HRMS (EI) calcd for C₁₉H₁₄⁵⁶FeN₂O [M⁺] 342.0456, found 342.0453.

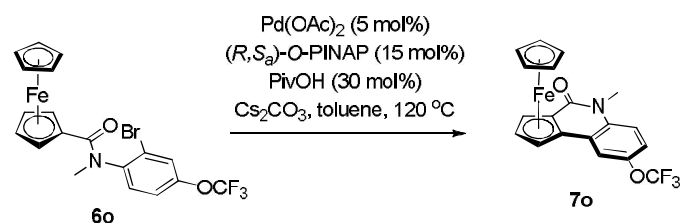


The reaction of **6n** (42.3 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)₂ (1.1 mg, 0.005

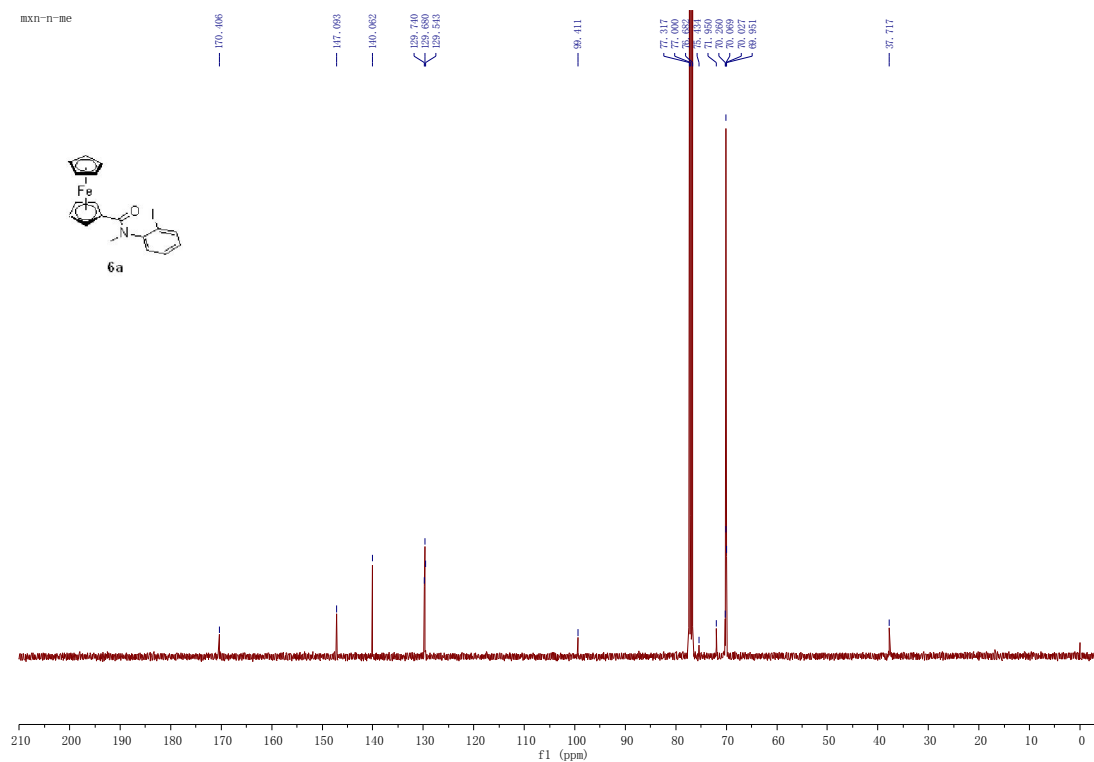
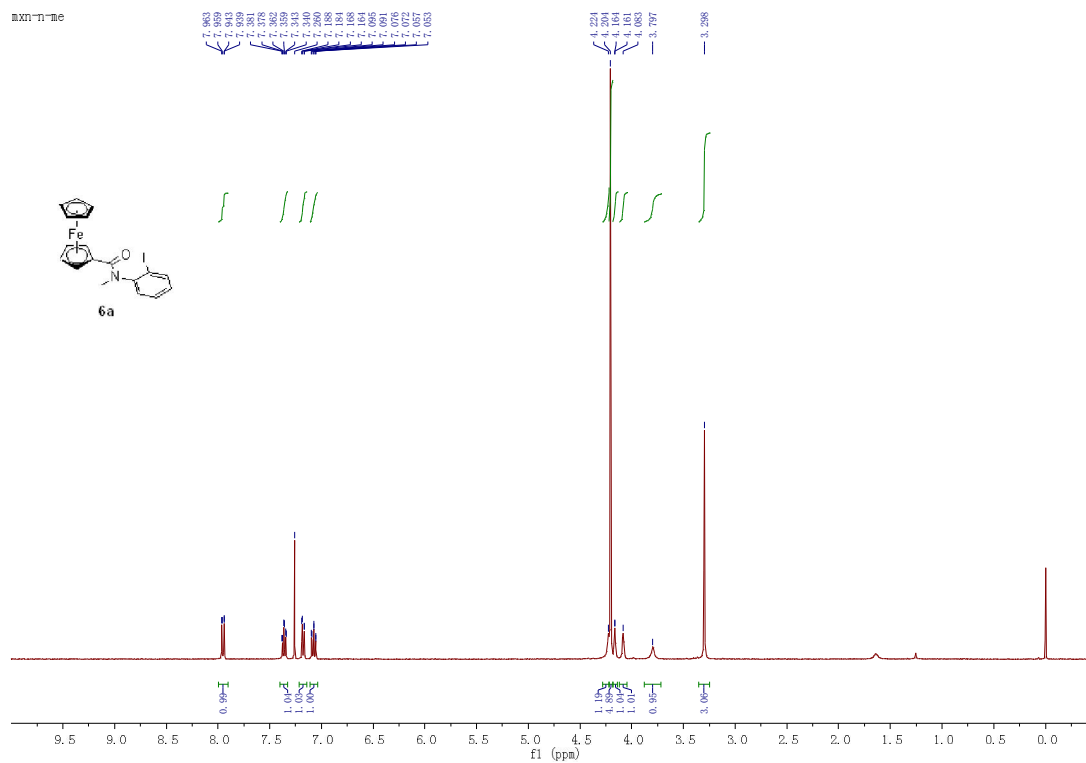
mmol, 5 mol%), (*R,S_a*)-*O*-PINAP (8.4 mg, 0.015 mmol, 15 mol%), PivOH (3.1 mg, 0.030 mmol, 30 mol %) and Cs₂CO₃ (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7n** (R_f = 0.20, PE:EA = 2:1) (33.5 mg, 98%, 4% ee).

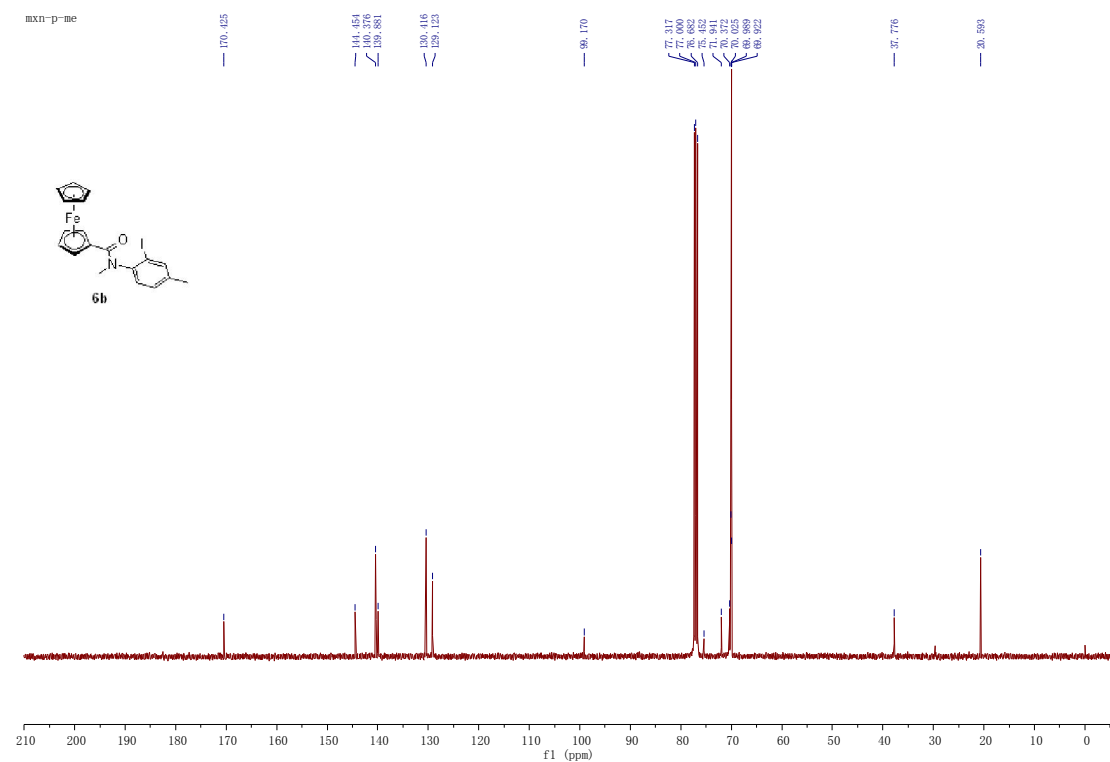
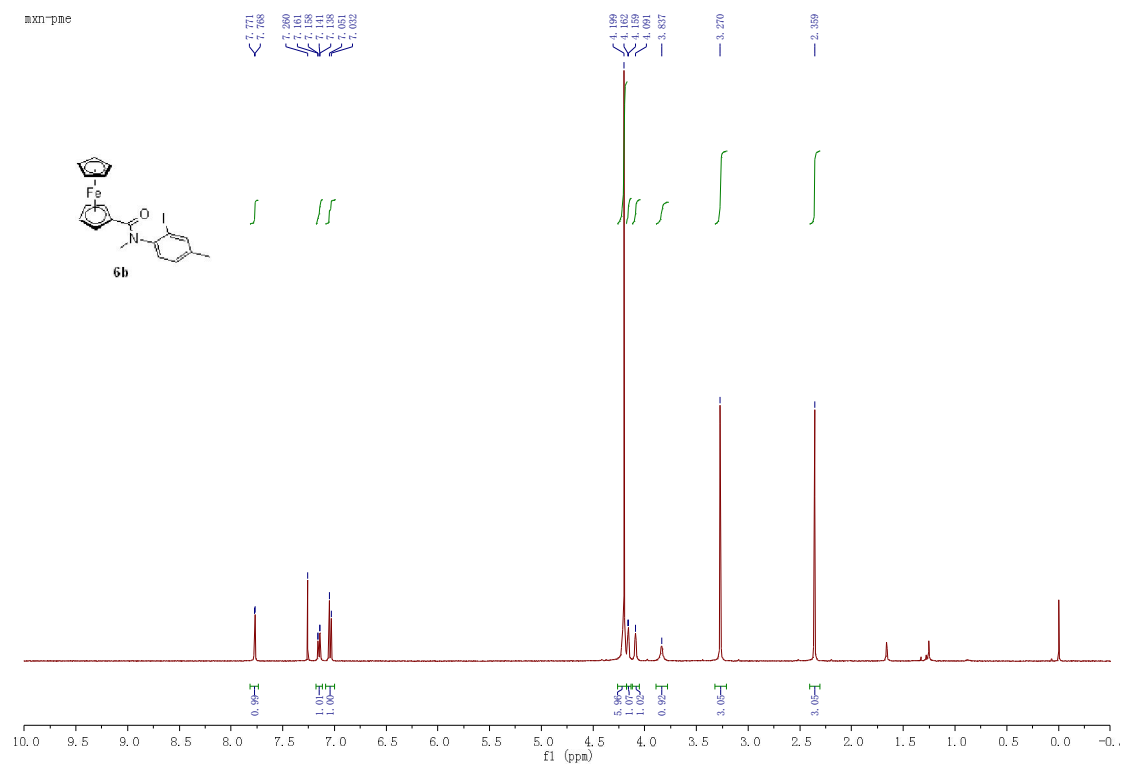


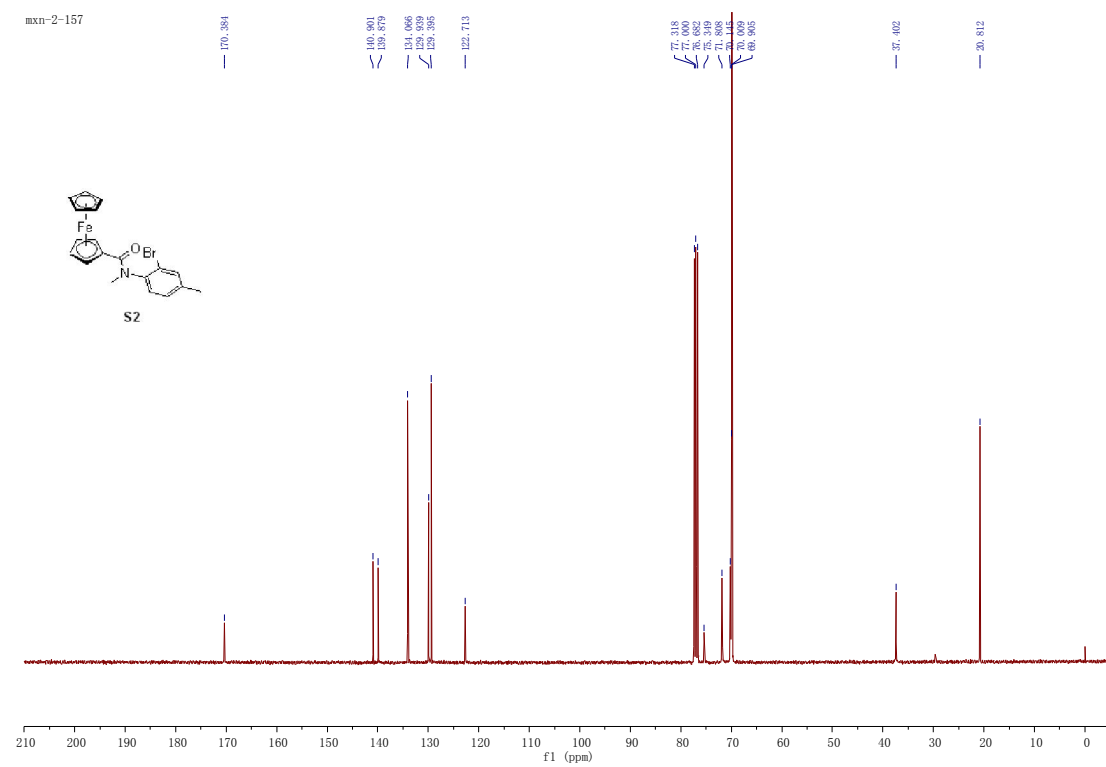
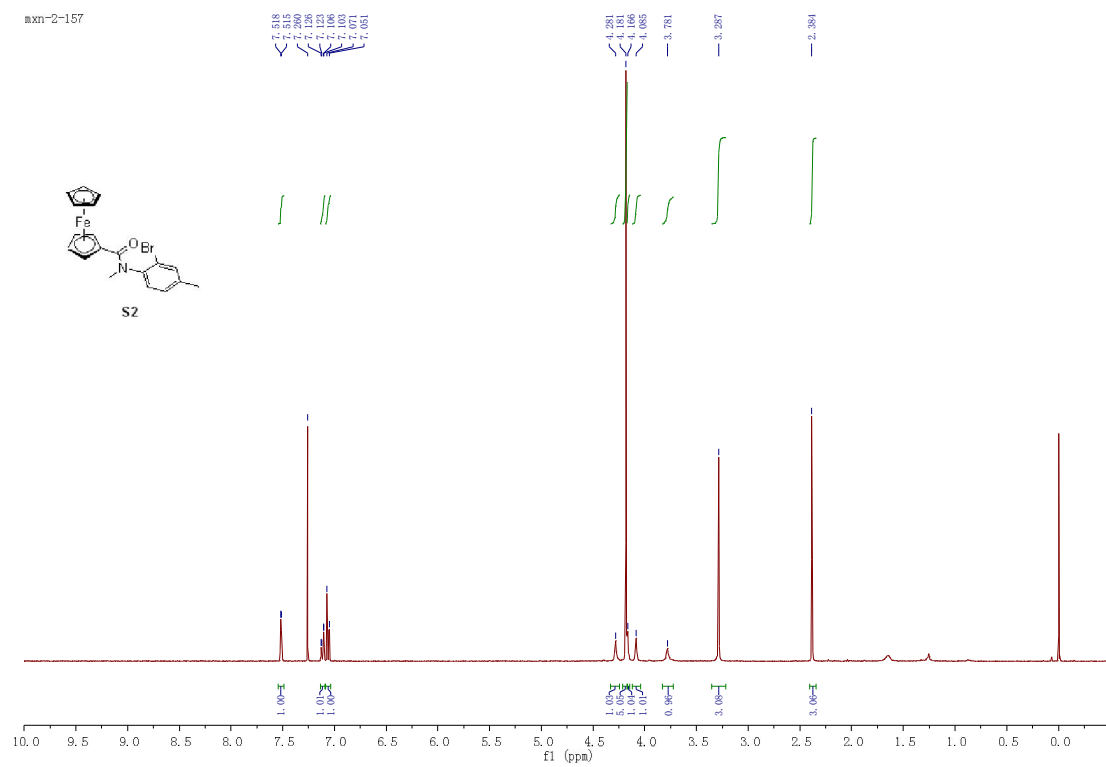
The reaction of **6o** (48.1 mg, 0.10 mmol, 1.0 equiv), Pd(OAc)₂ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S_a*)-*O*-PINAP (8.4 mg, 0.015 mmol, 15 mol%) and Cs₂CO₃ (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7o** (R_f = 0.30, PE:EA = 3:1) (13.2 mg, 33%, 48% ee). [α]_D²³ -11.0 (c 1.0, CHCl₃). HPLC conditions: Chiralcel AD-H, isopropanol/hexane = 6:94, flow: 1.0 ml/min, λ = 254 nm. ¹H NMR (400 MHz, CDCl₃) δ 7.54 (s, 1 H), 7.26-7.22 (m, 2 H), 5.24 (dd, *J* = 2.4, 1.2 Hz, 1 H), 5.13 (dd, *J* = 2.4, 1.2 Hz, 1 H), 4.54 (t, *J* = 2.4 Hz, 1 H), 3.97 (s, 5 H), 3.69 (s, 3 H); ¹³C NMR (100 MHz, CDCl₃) δ 167.5, 144.1, 136.8, 124.8, 124.4, 121.9, 119.6, 119.3, 116.0, 115.8, 81.9, 71.8, 71.5, 70.6, 67.7, 64.0, 29.5. HRMS (EI) calcd for C₁₉H₁₄F₃⁵⁶FeNO₂ [*M*⁺] 401.0326, found 401.0319.

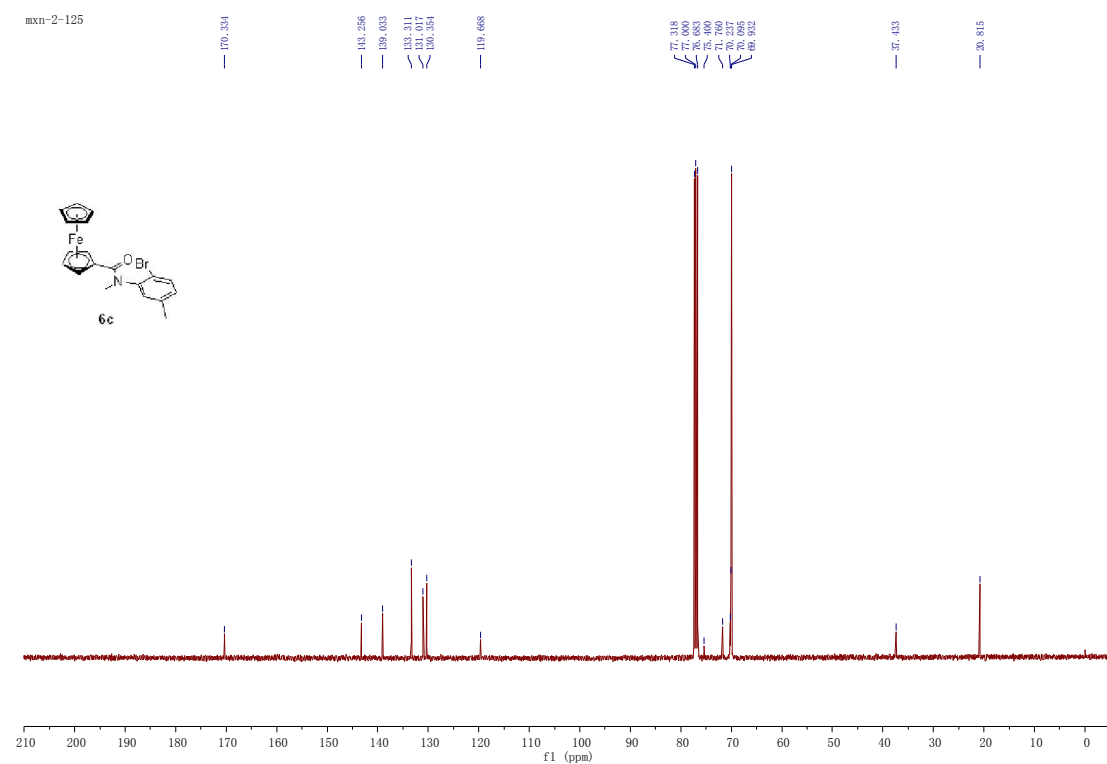
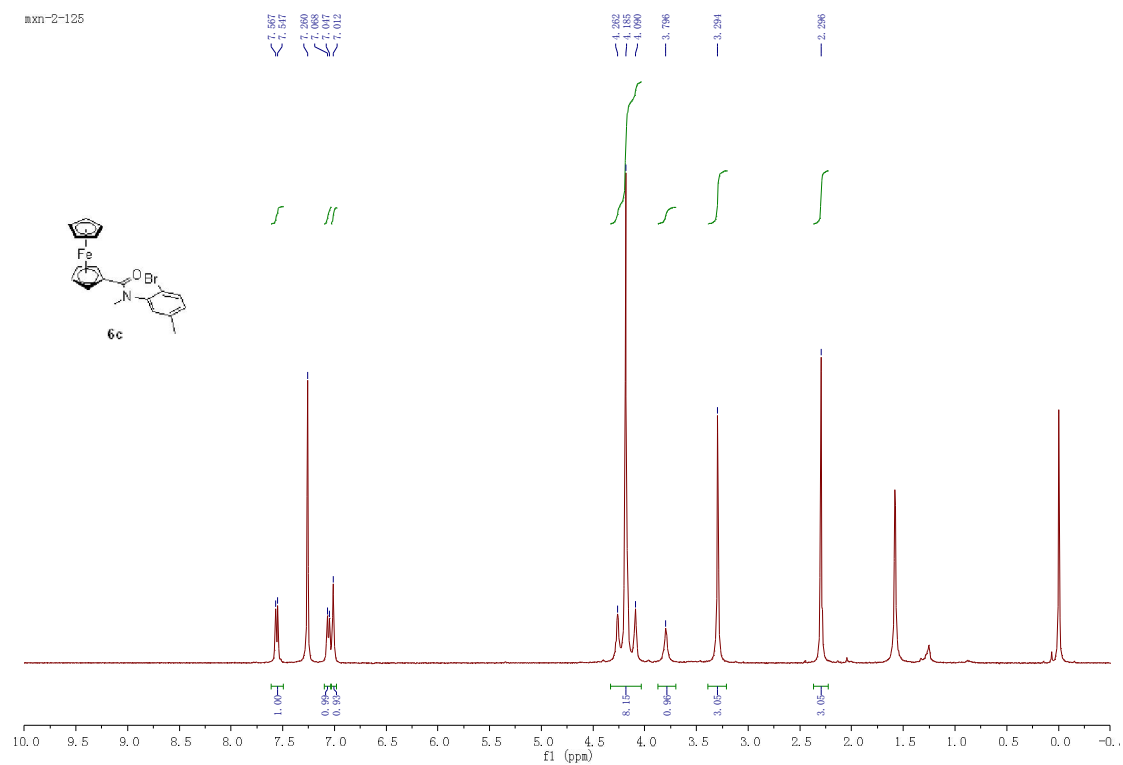


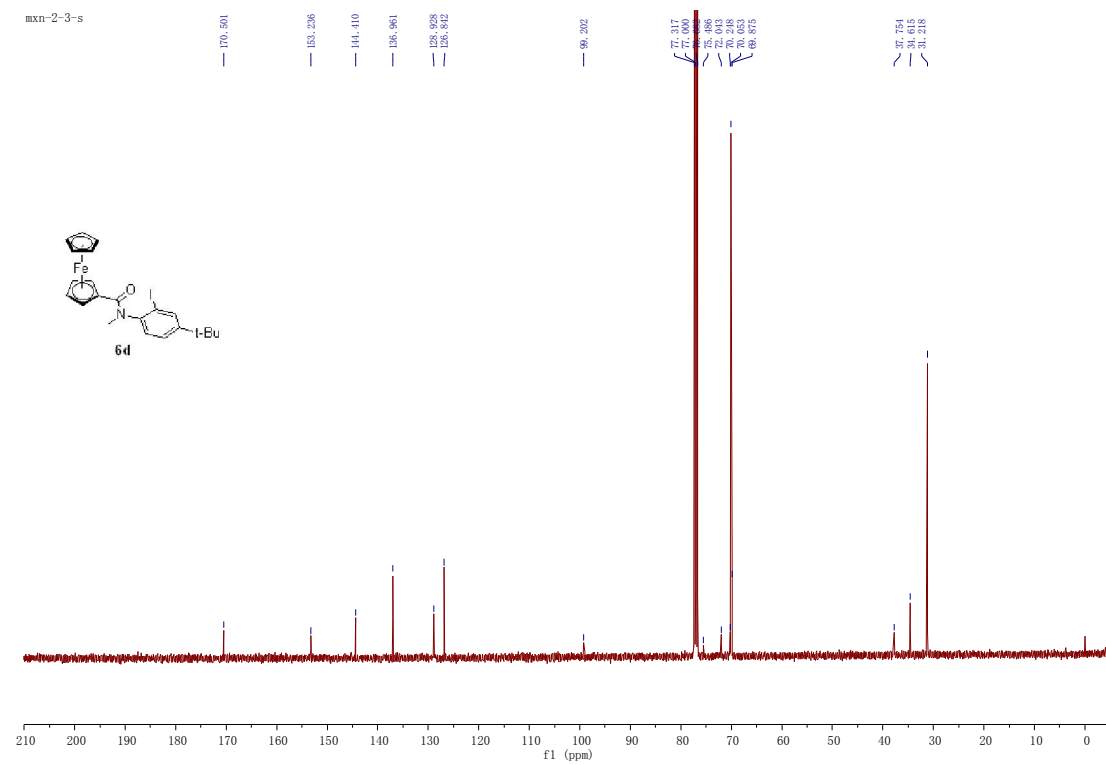
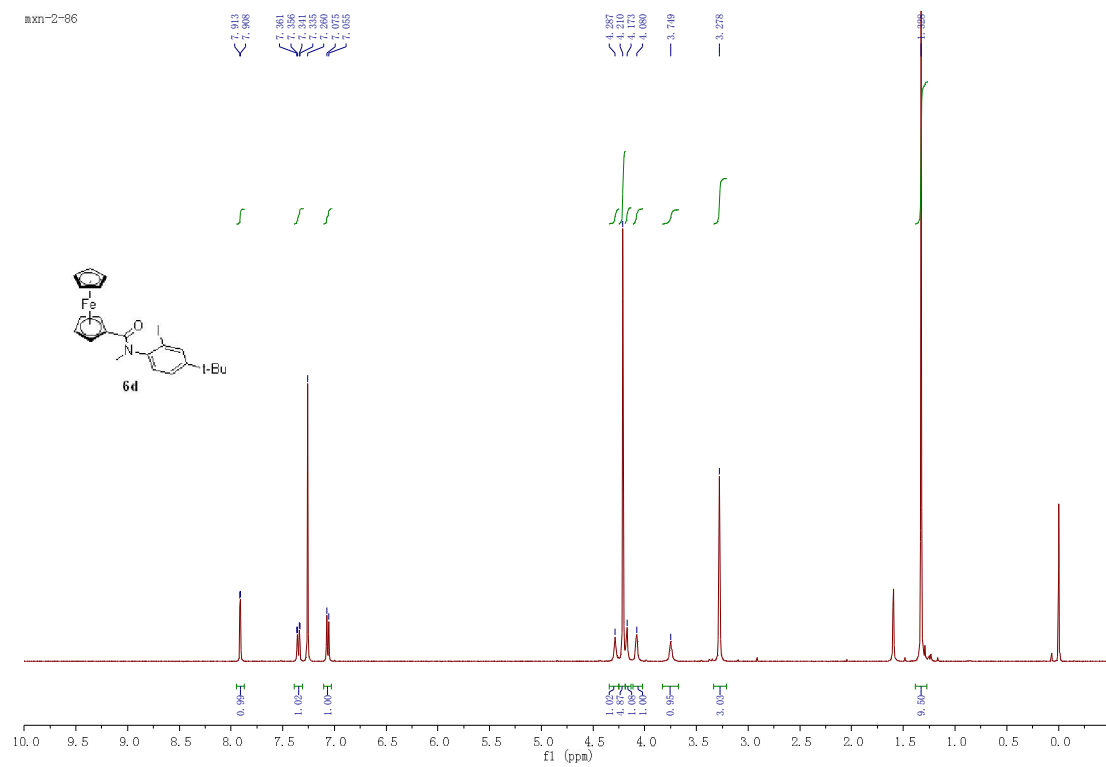
The reaction of **6o** (48.1mg, 0.10 mmol, 1.0 equiv), Pd(OAc)₂ (1.1 mg, 0.005 mmol, 5 mol%), (*R,S_a*)-*O*-PINAP (8.4 mg, 0.015 mmol, 15 mol%), PivOH (3.1 mg, 0.030 mmol, 30 mol %) and Cs₂CO₃ (81.5 mg, 0.25 mmol, 2.5 equiv) in toluene (2.0 ml) at 120 °C for 12 h afforded **7o** (R_f = 0.30, PE:EA = 3:1) (16.8 mg, 42%, 16% ee).



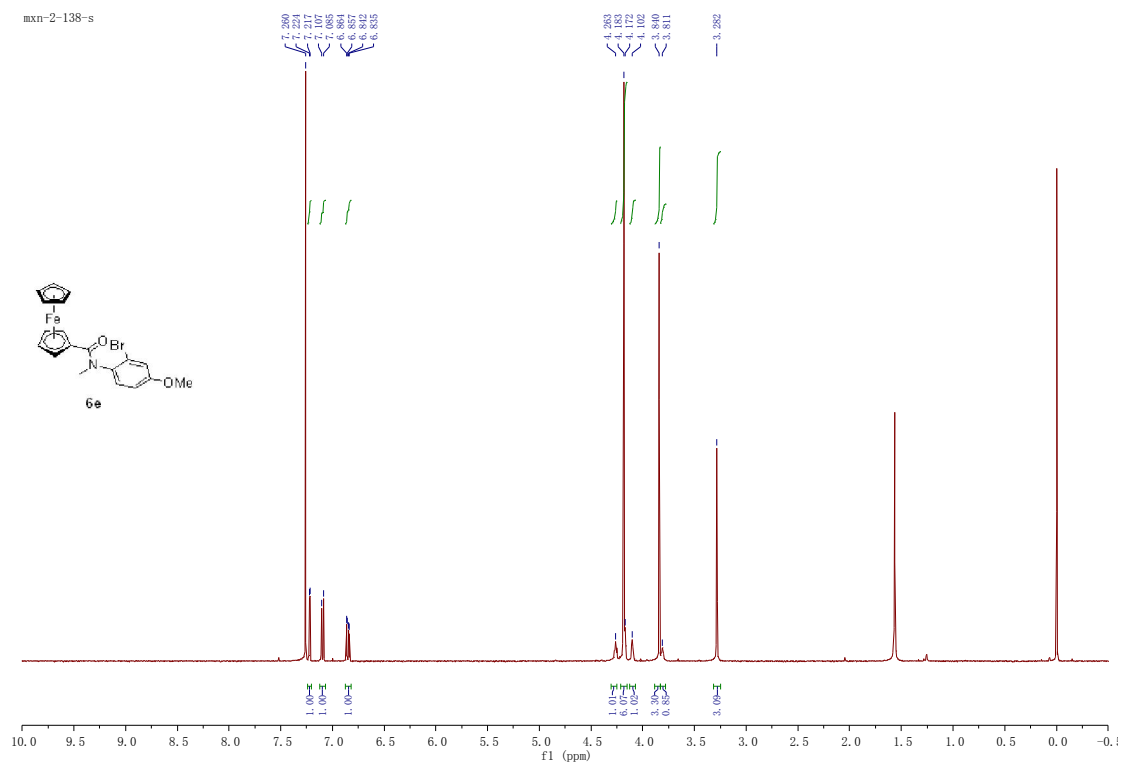




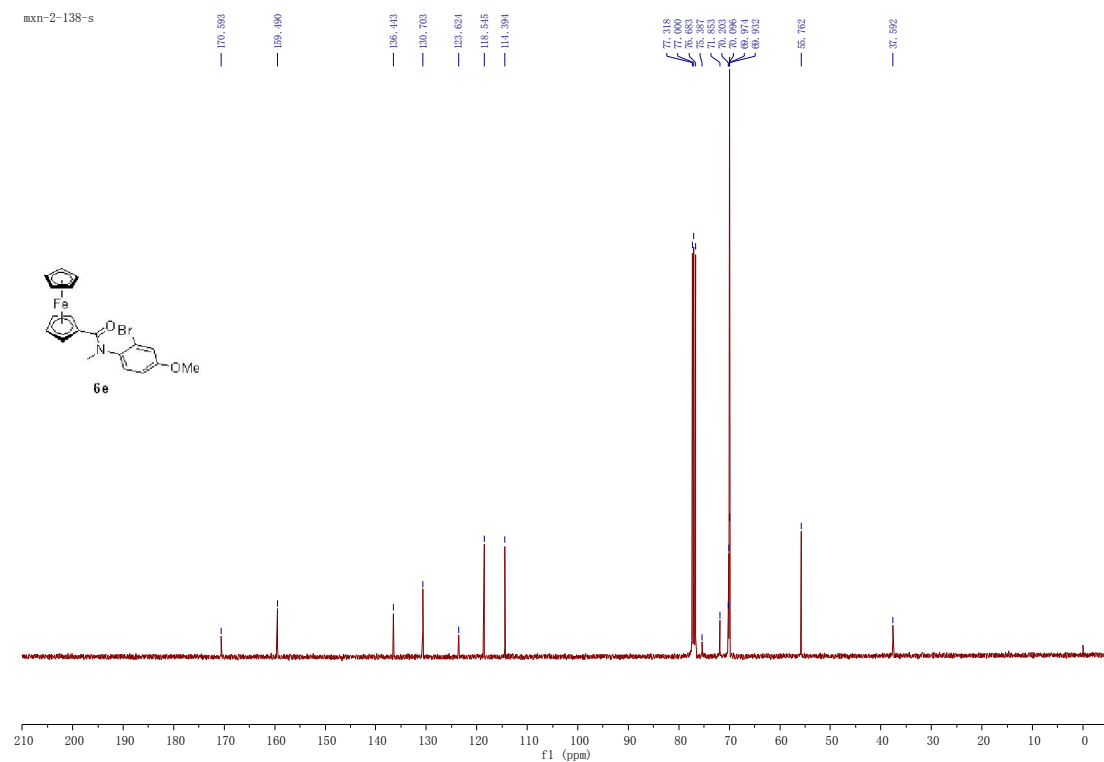


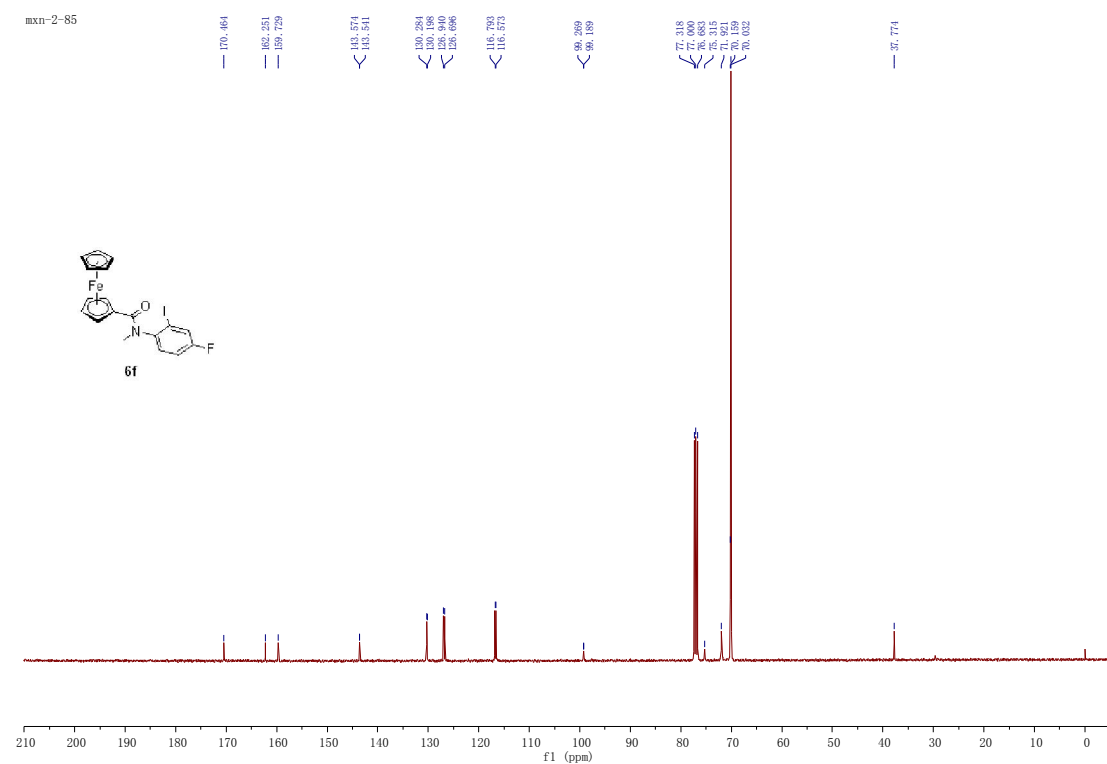
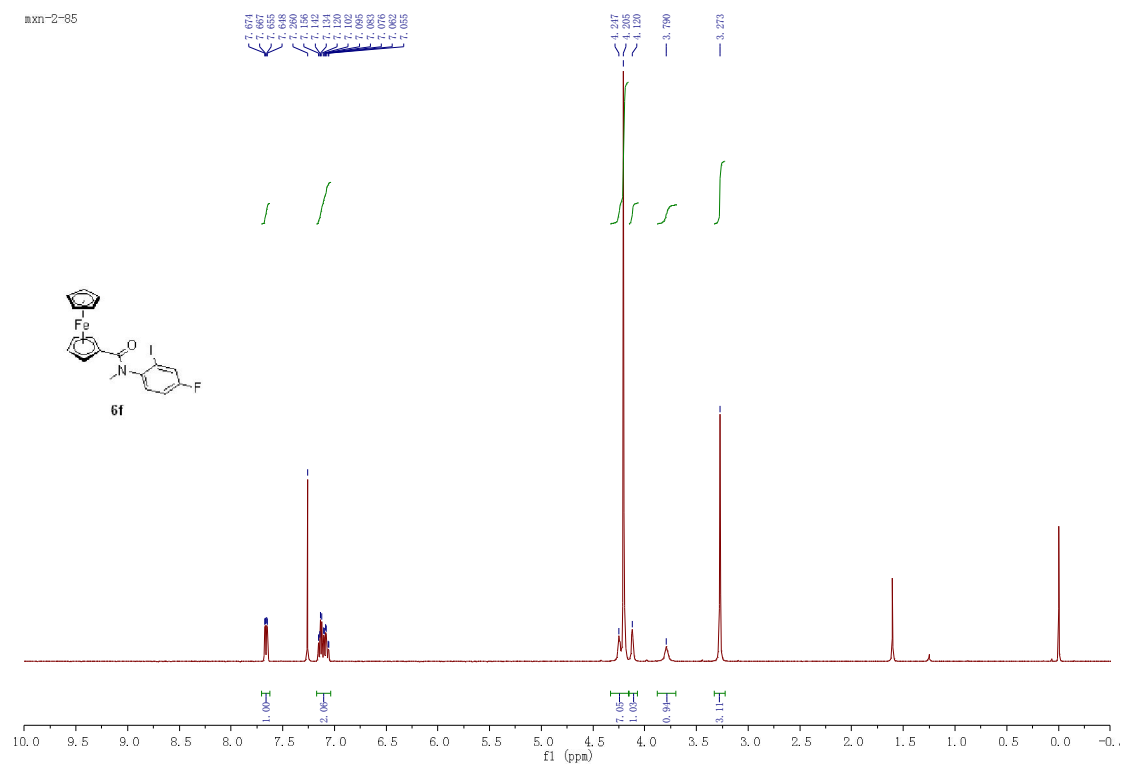


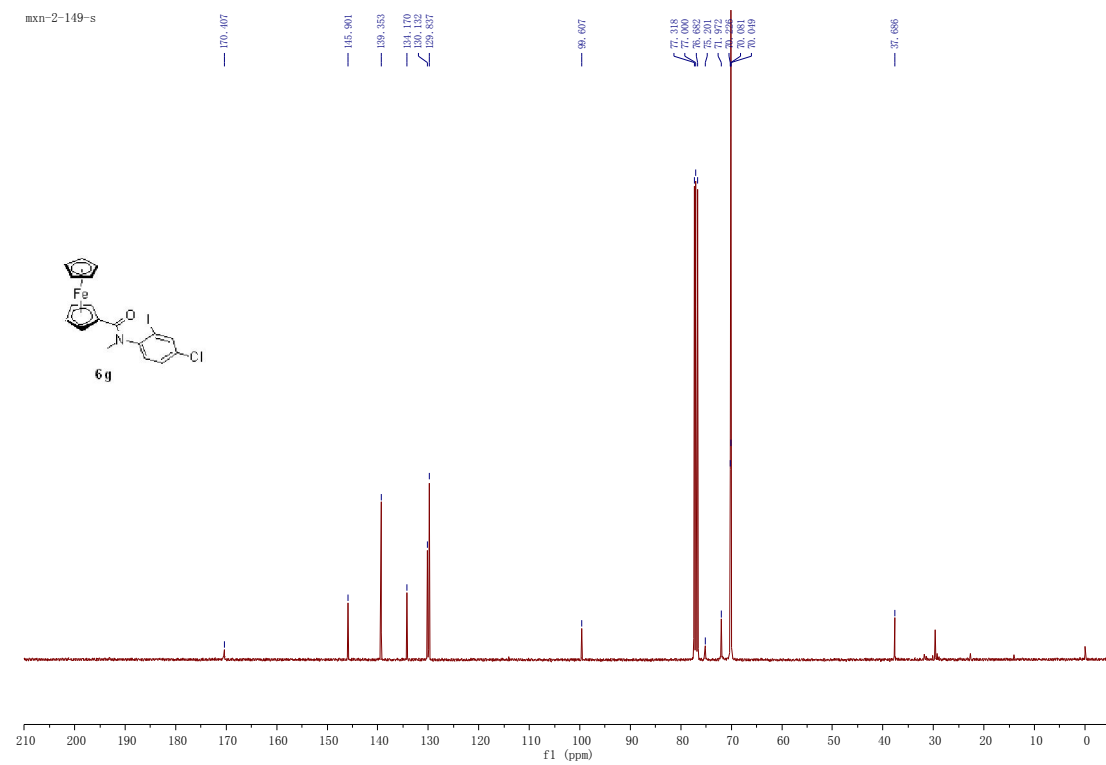
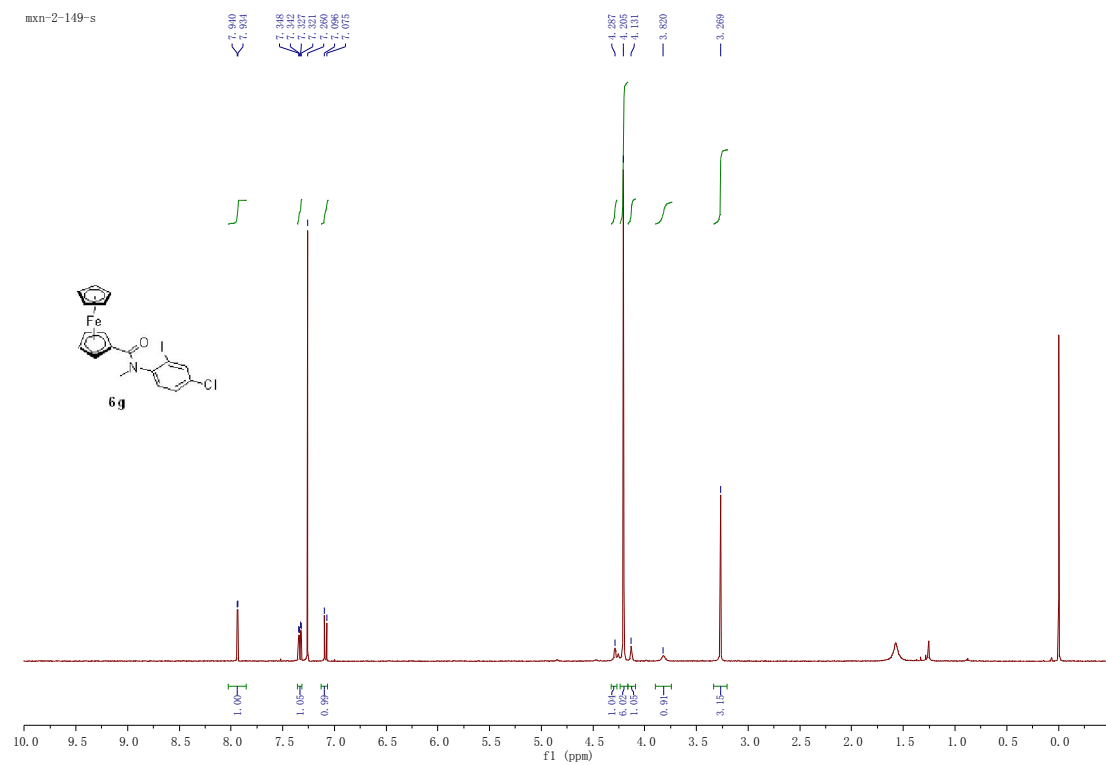
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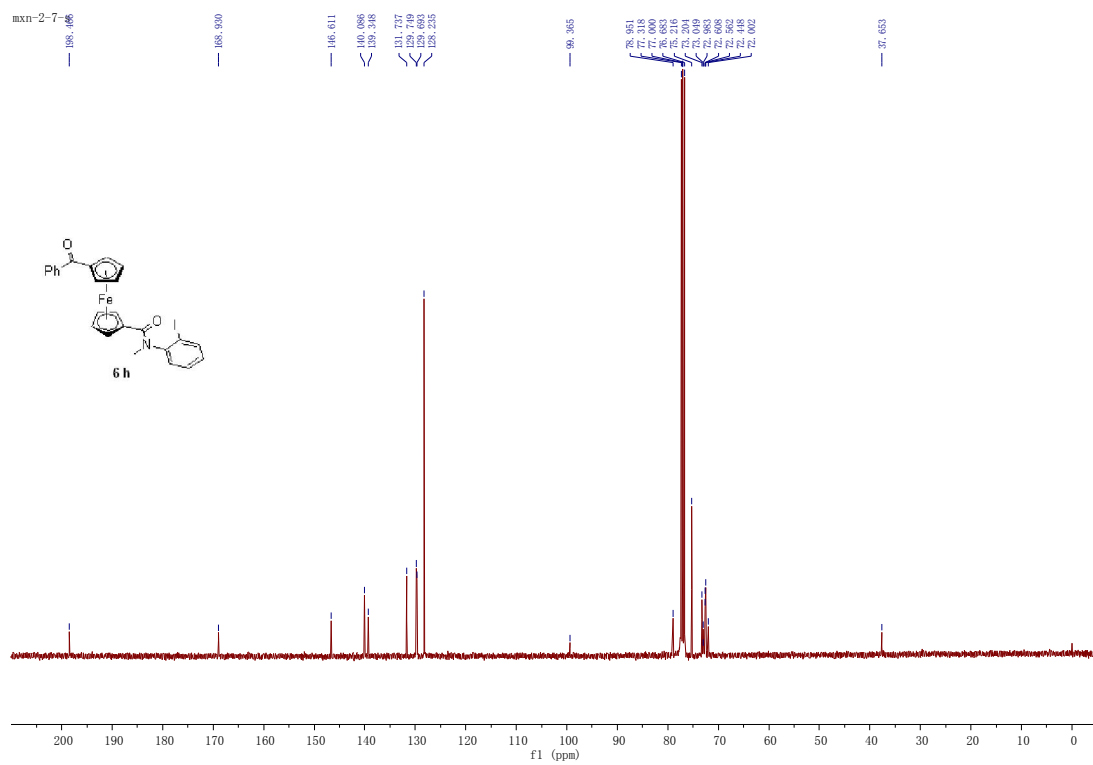
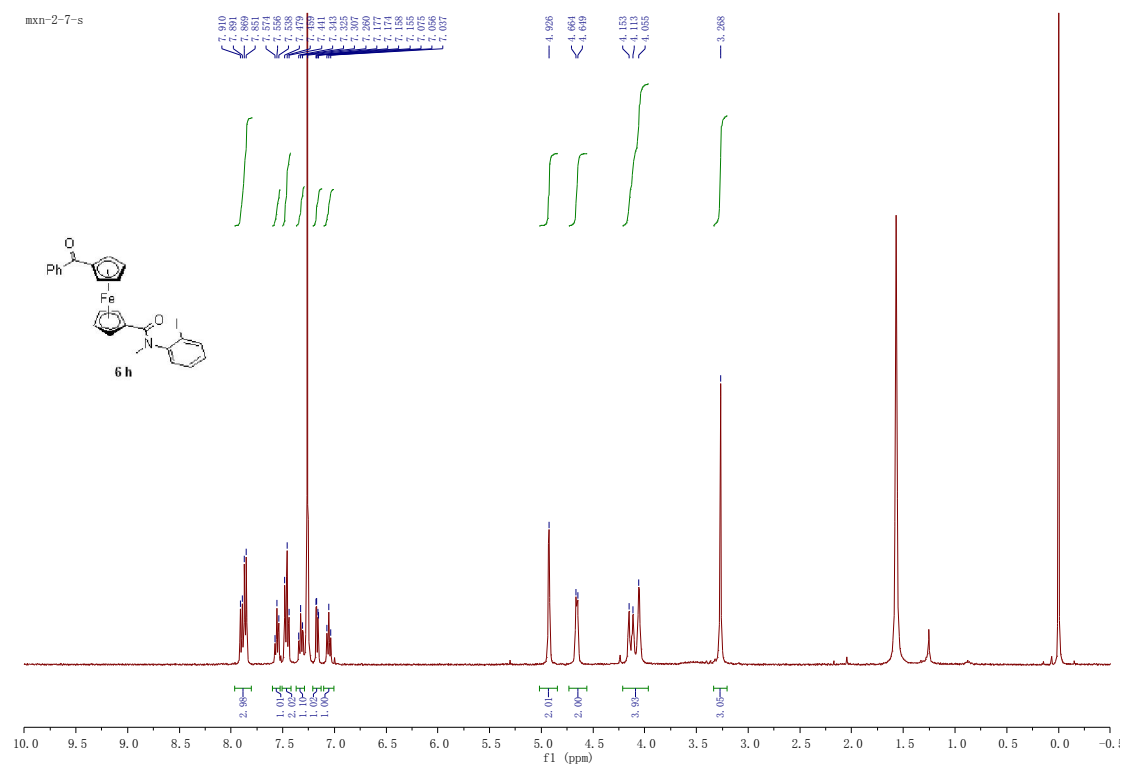


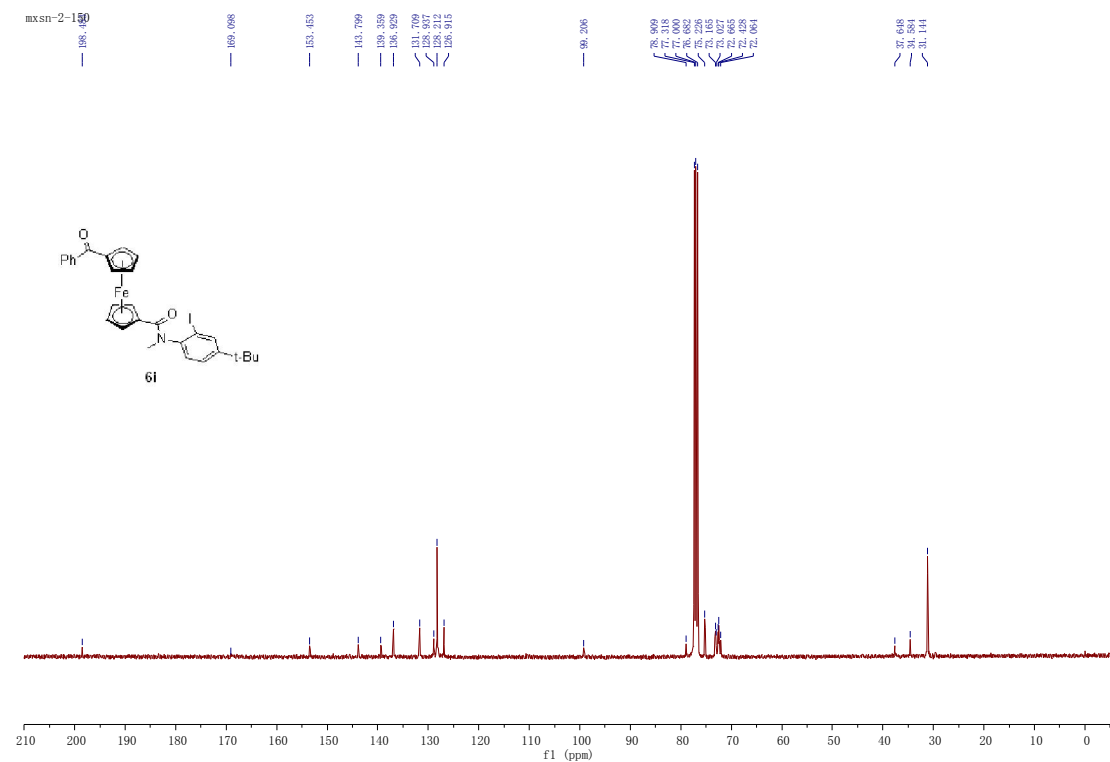
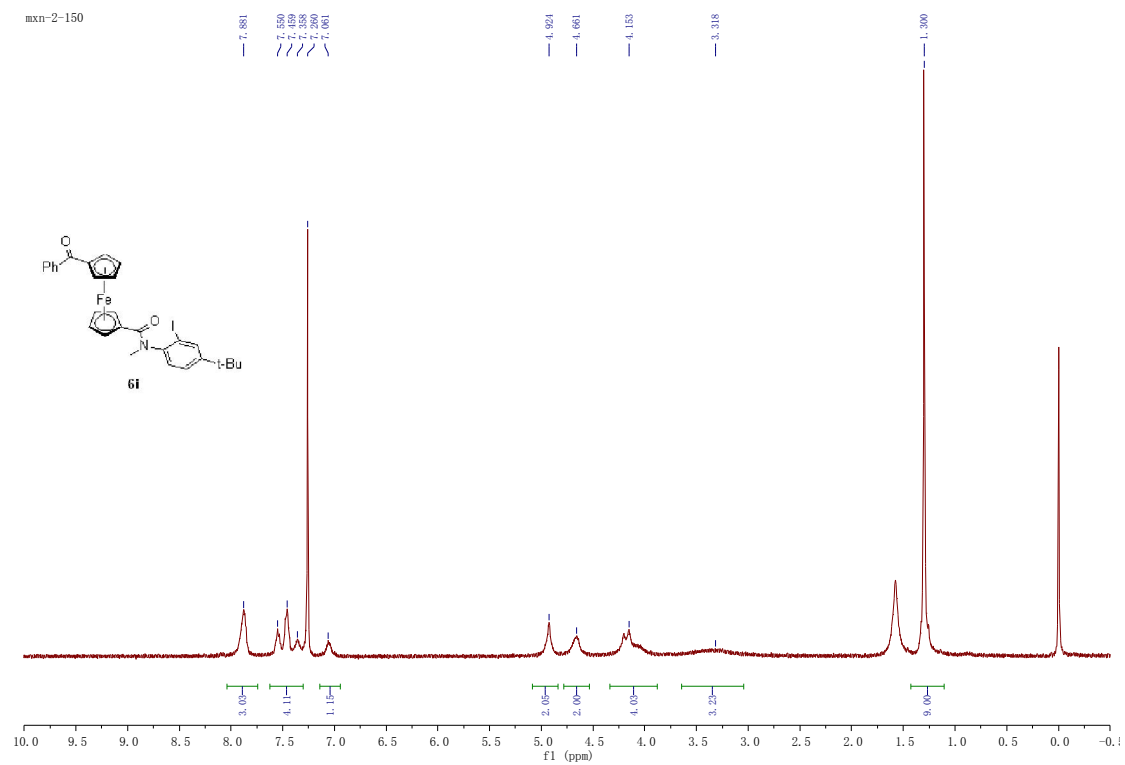
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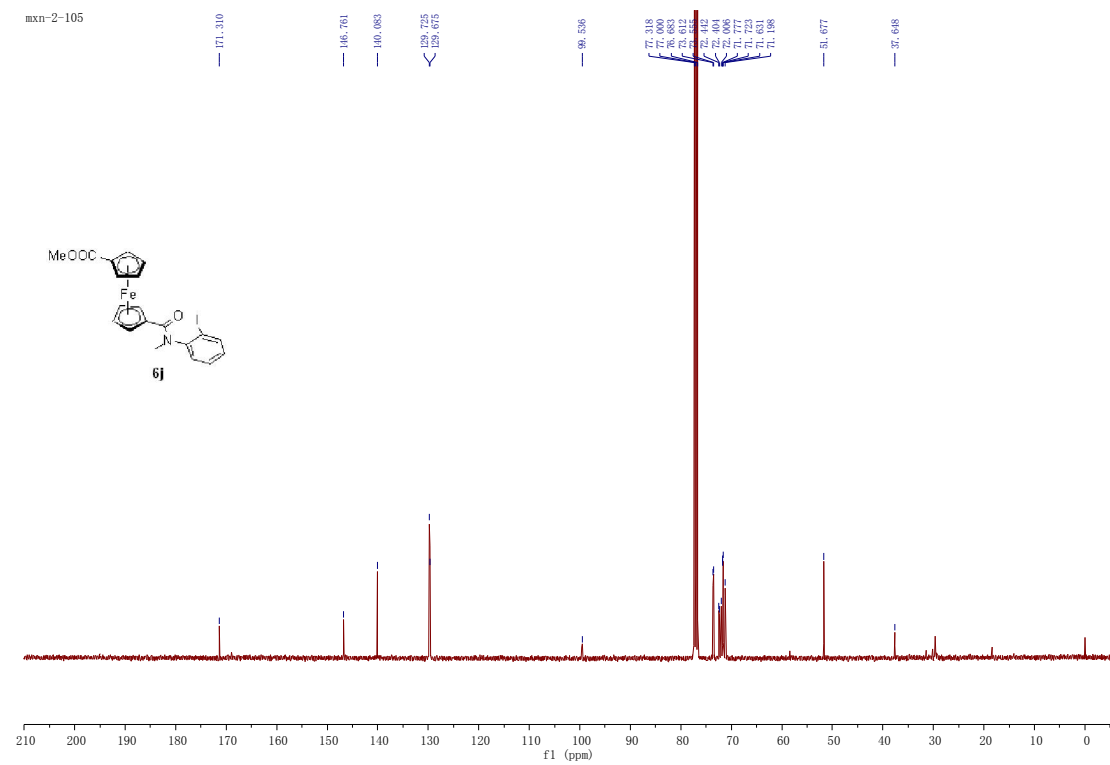
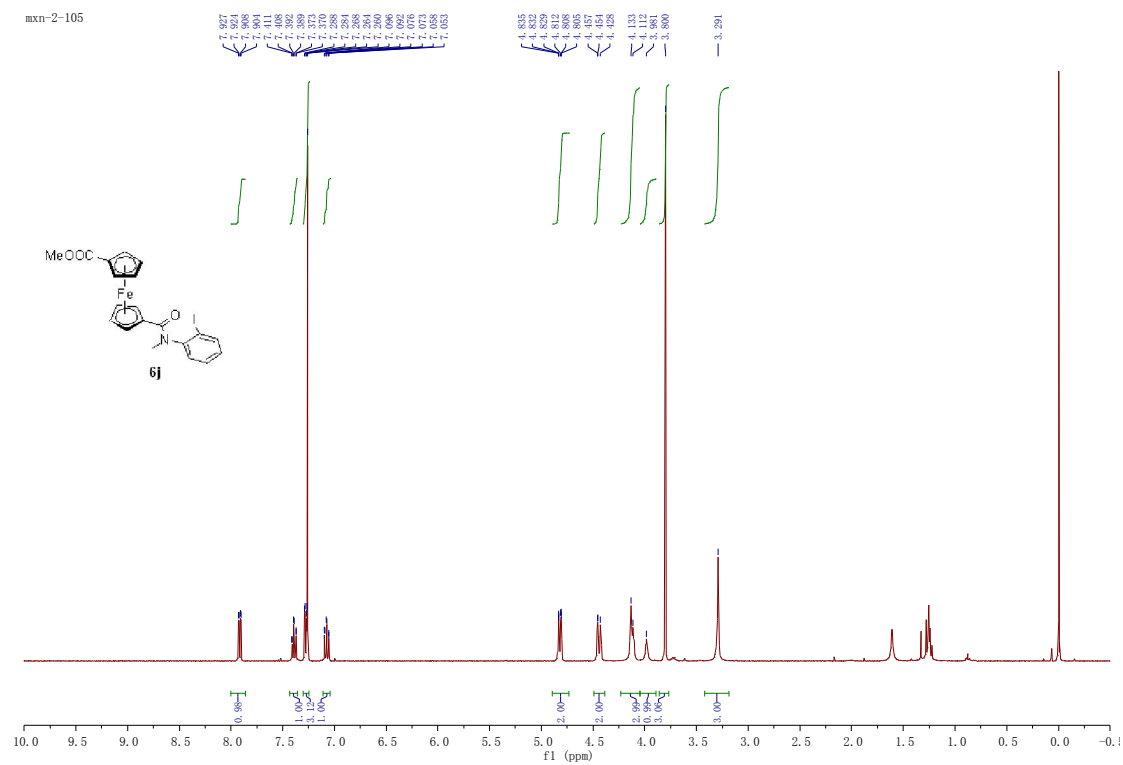




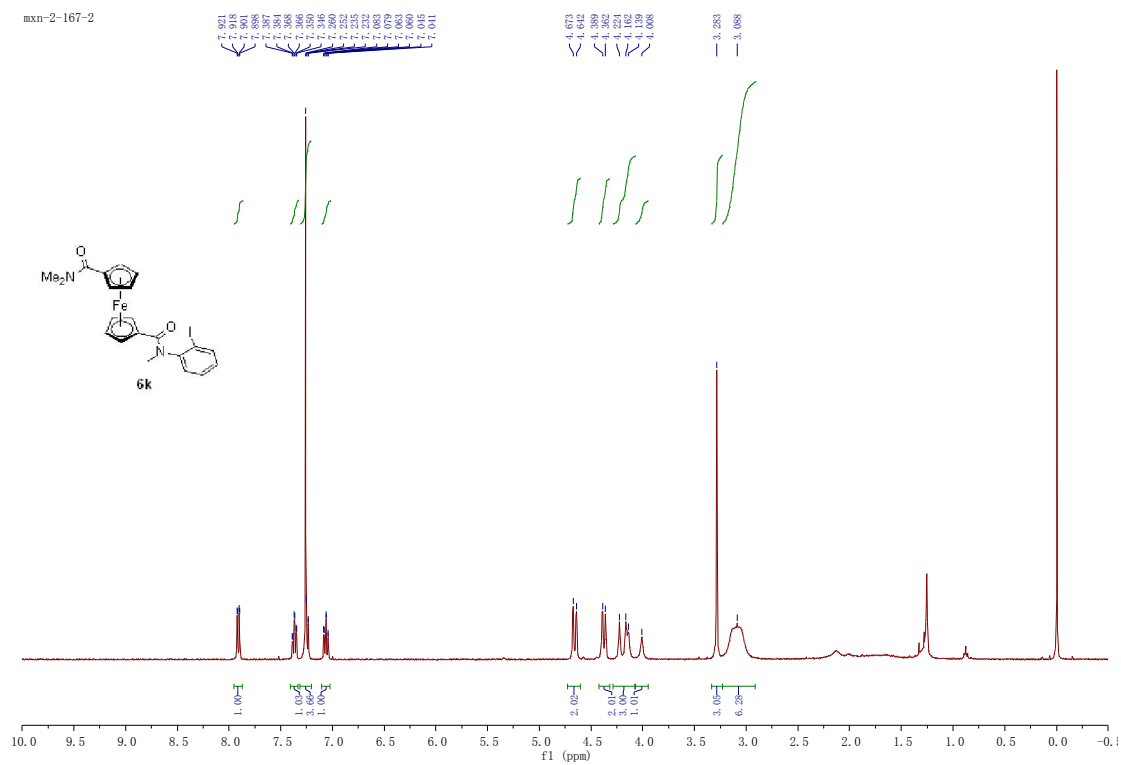




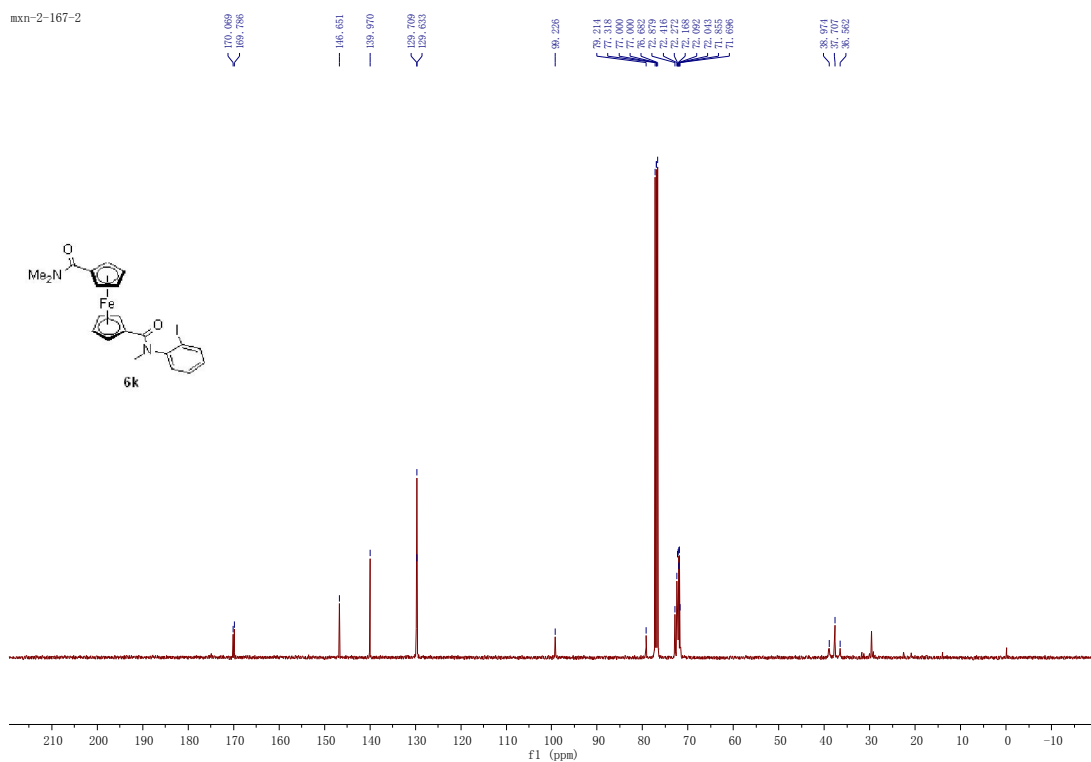


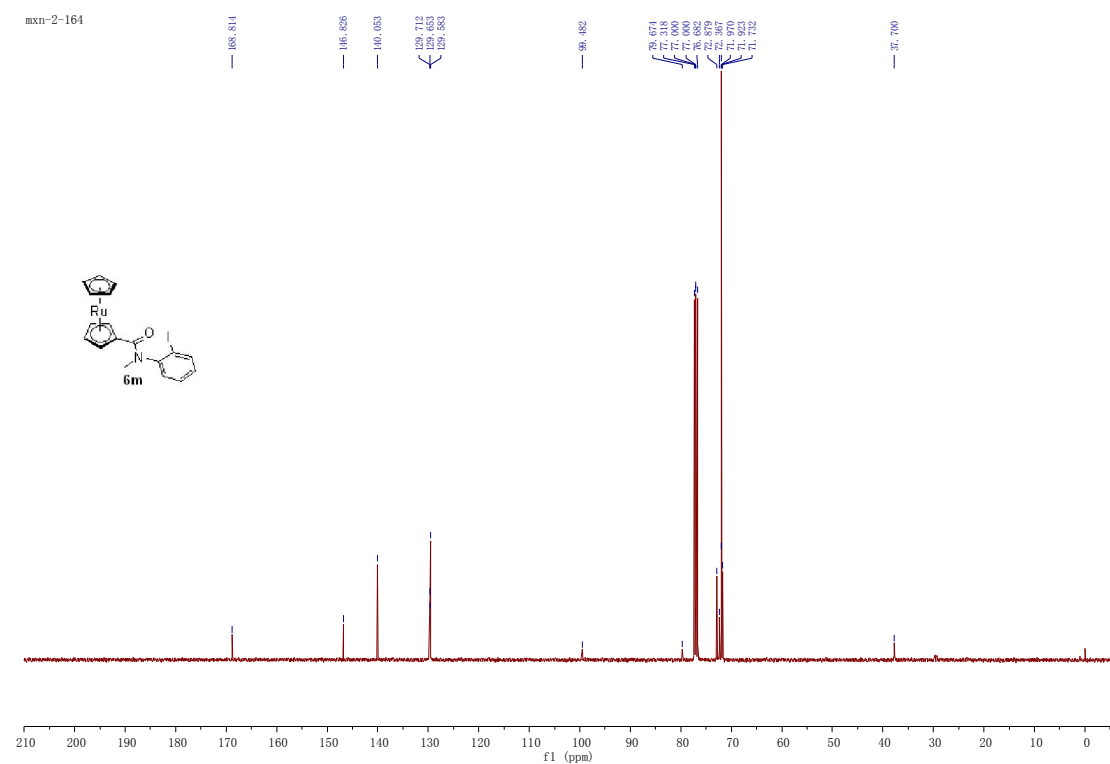
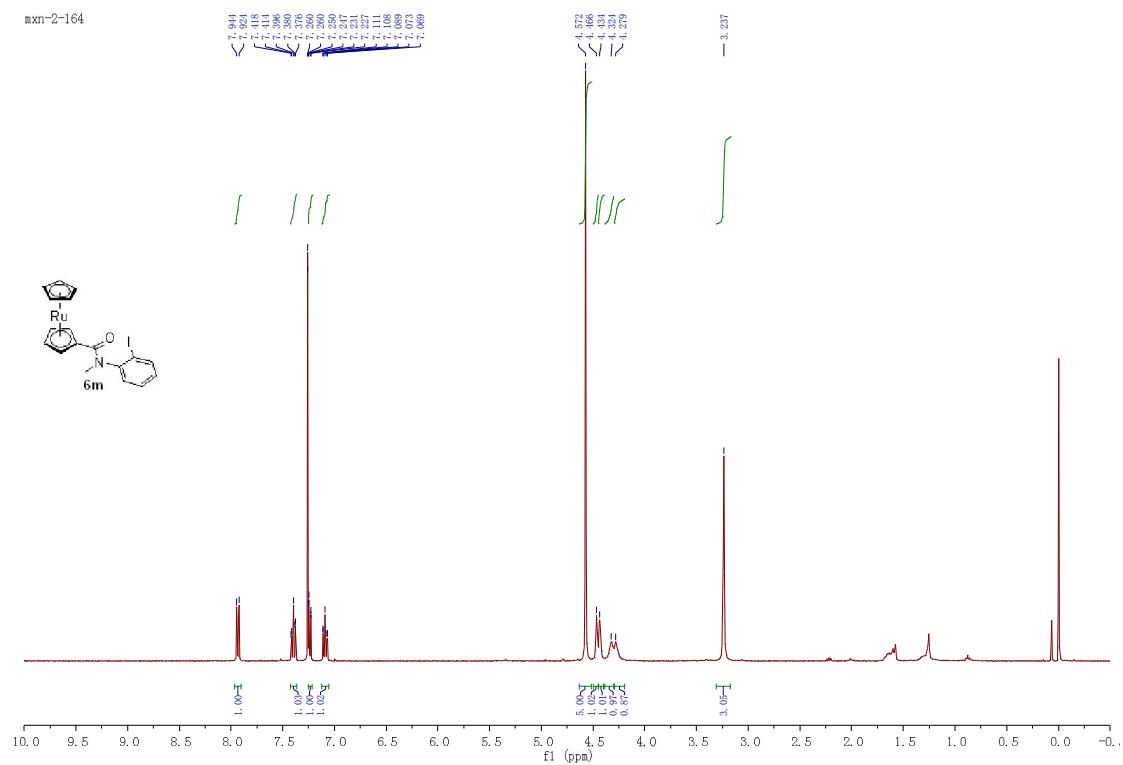


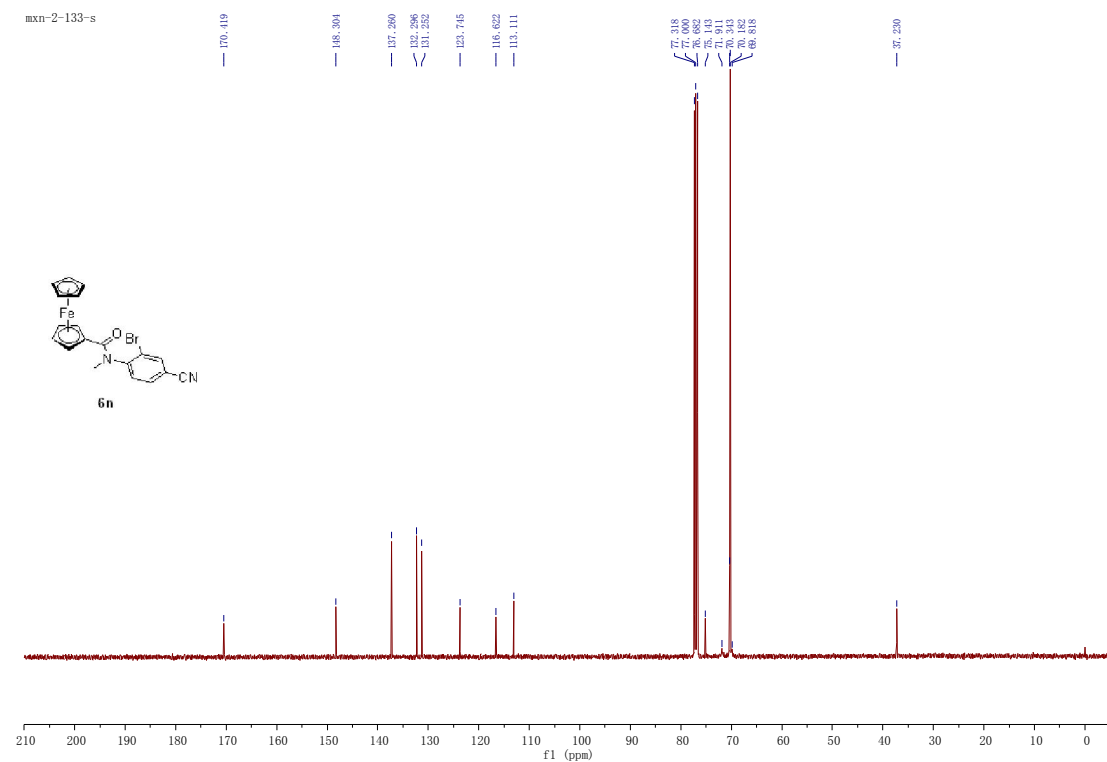
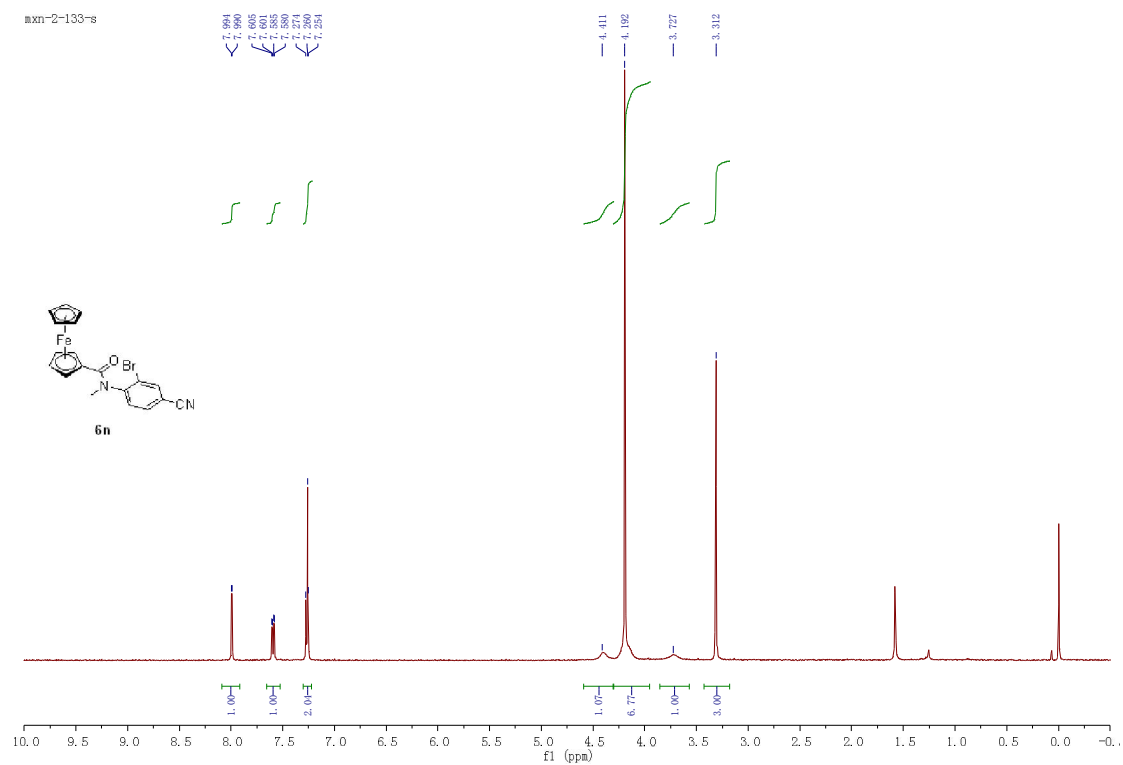
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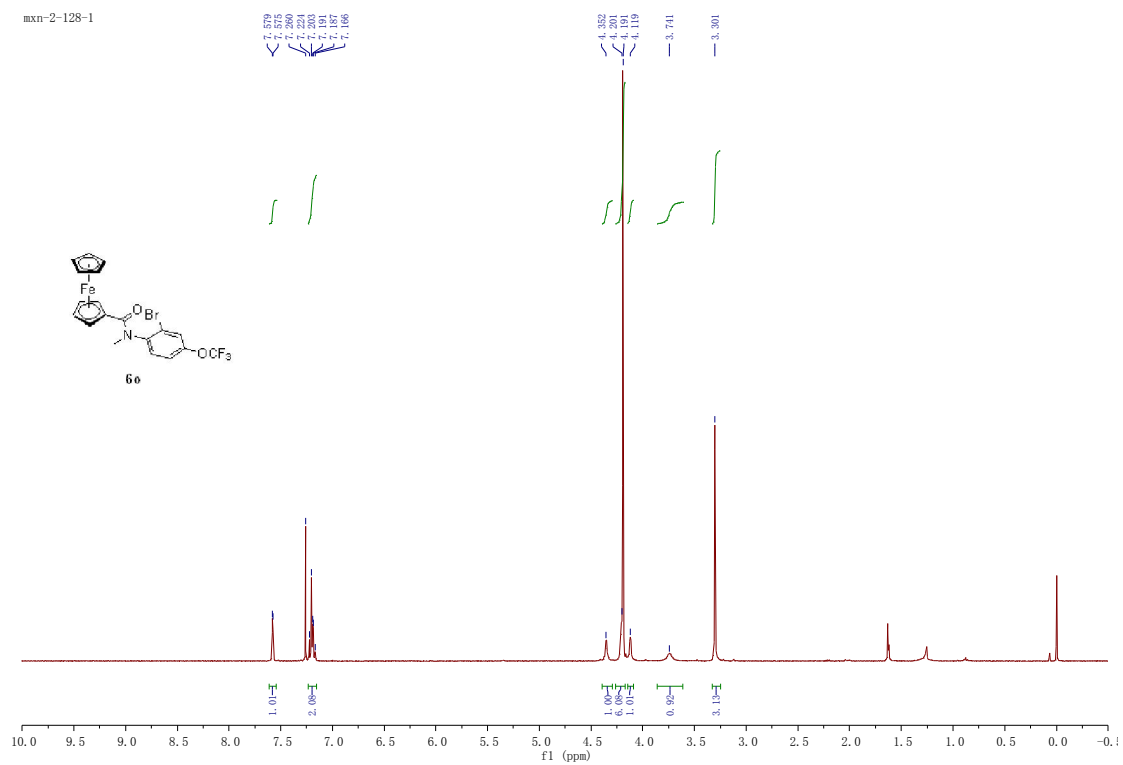
mxn-2-167-2



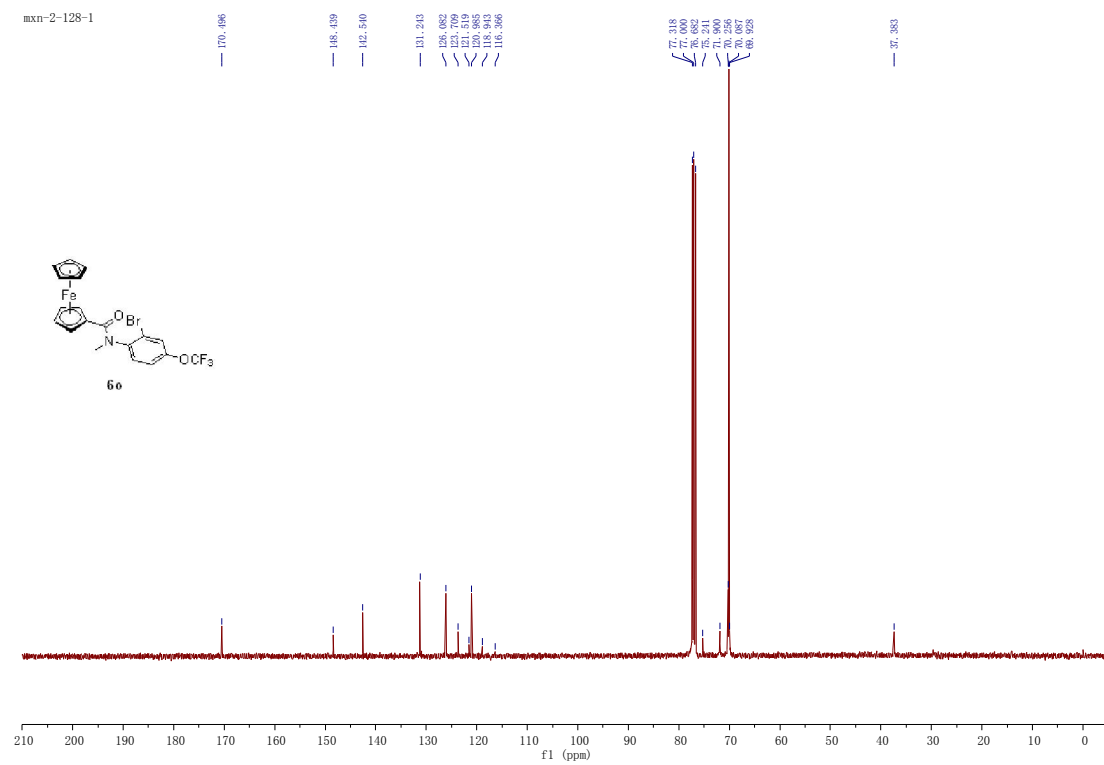


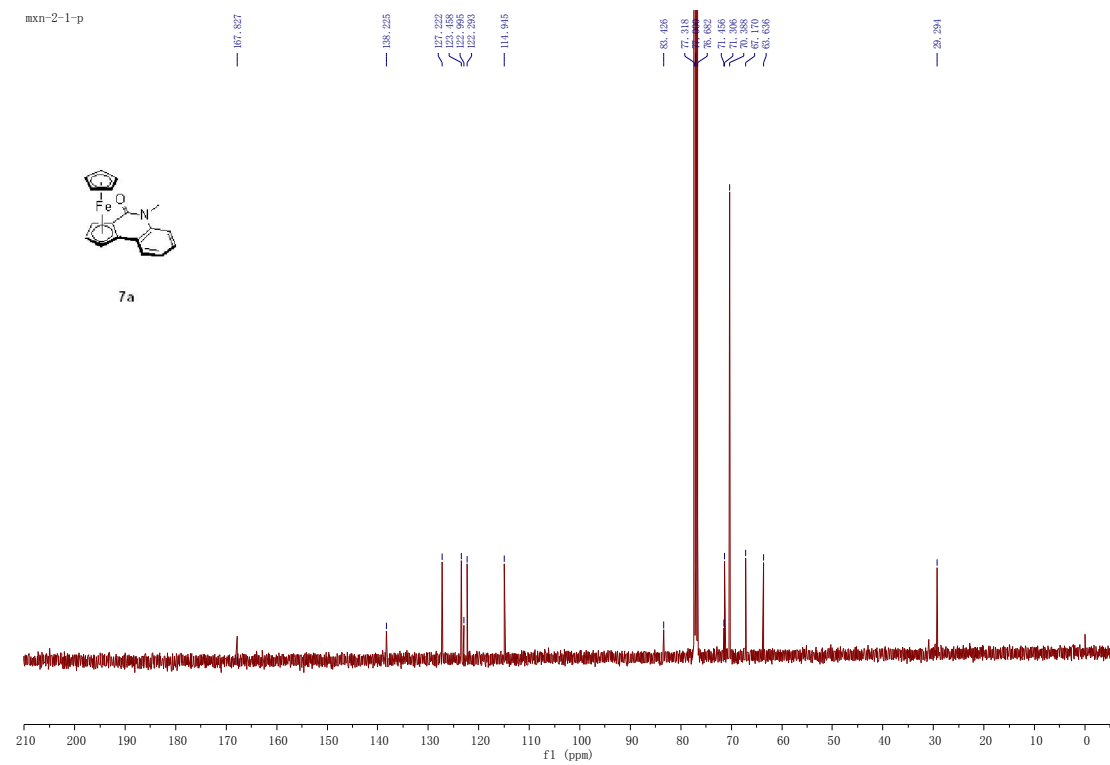
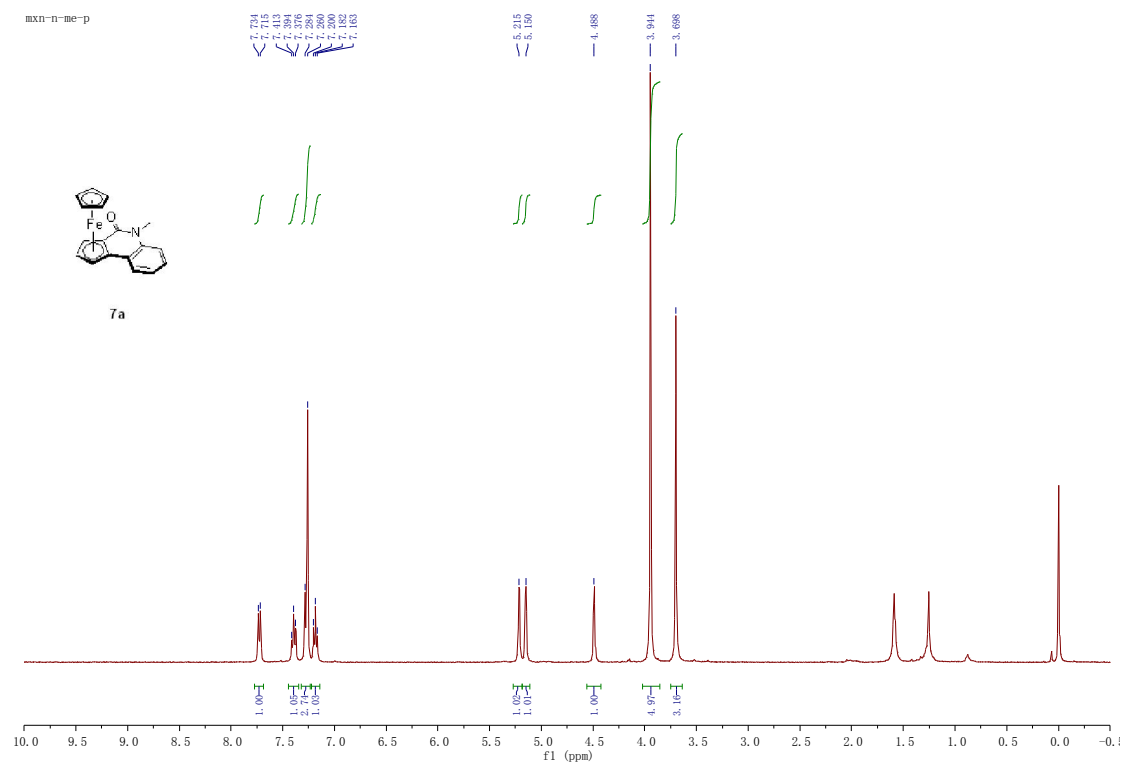


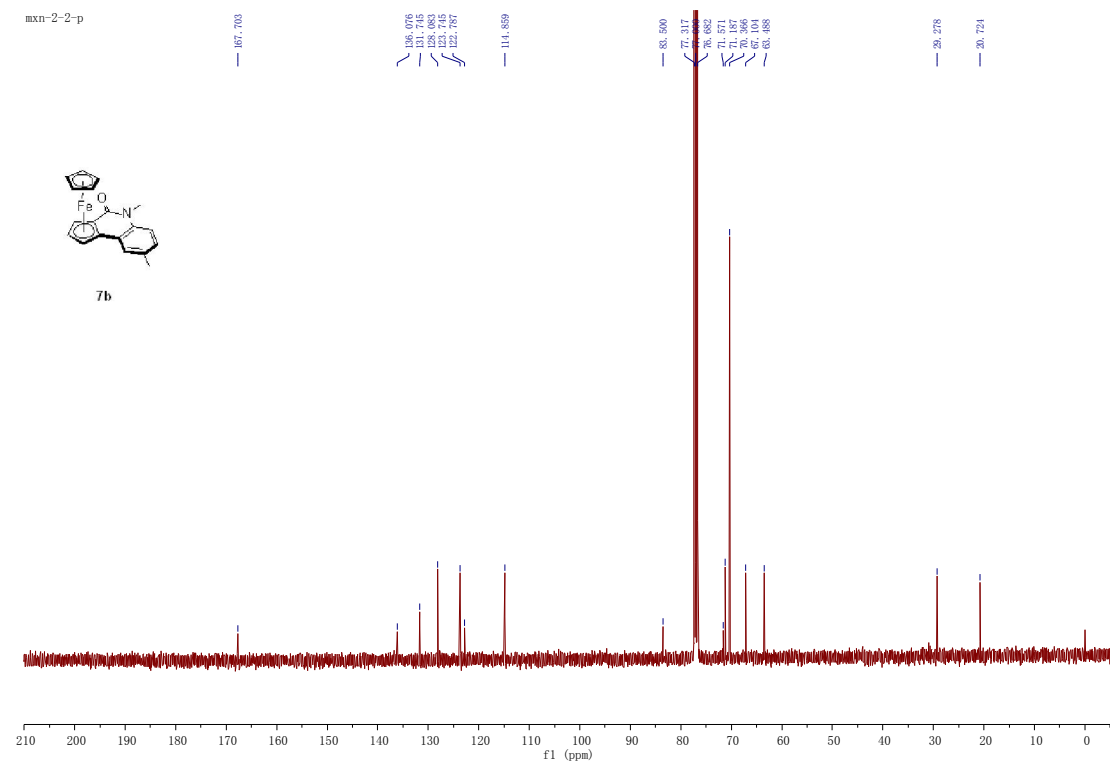
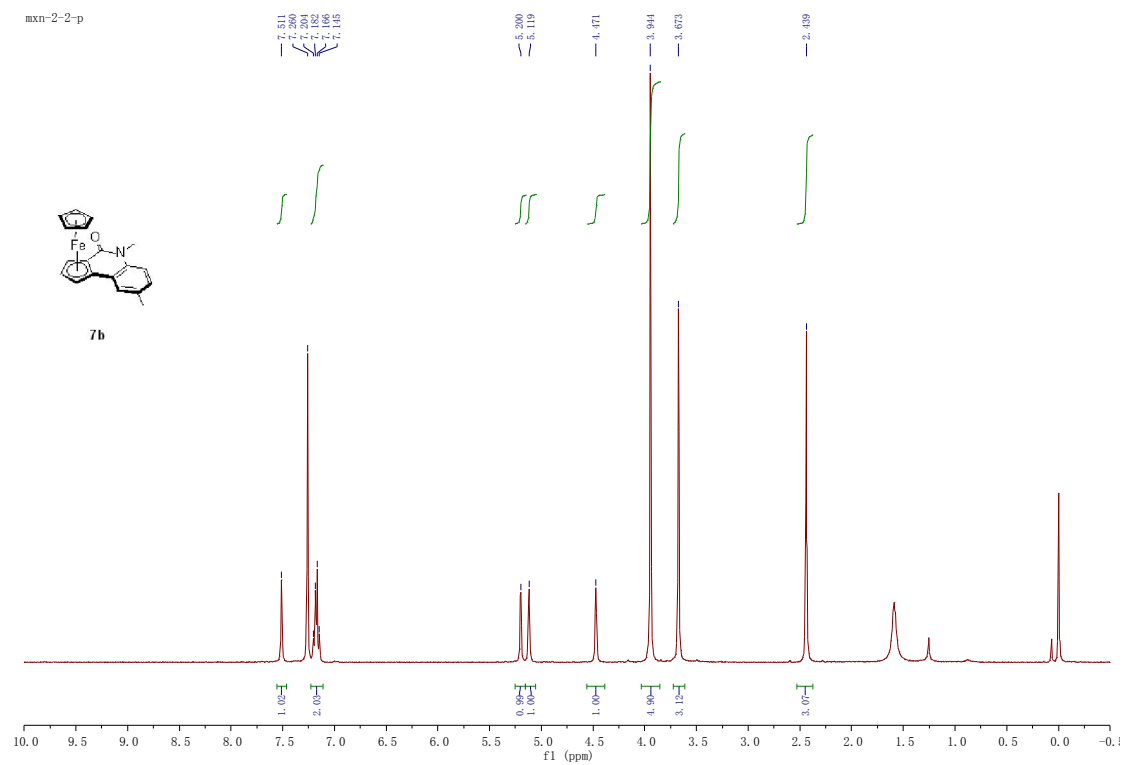
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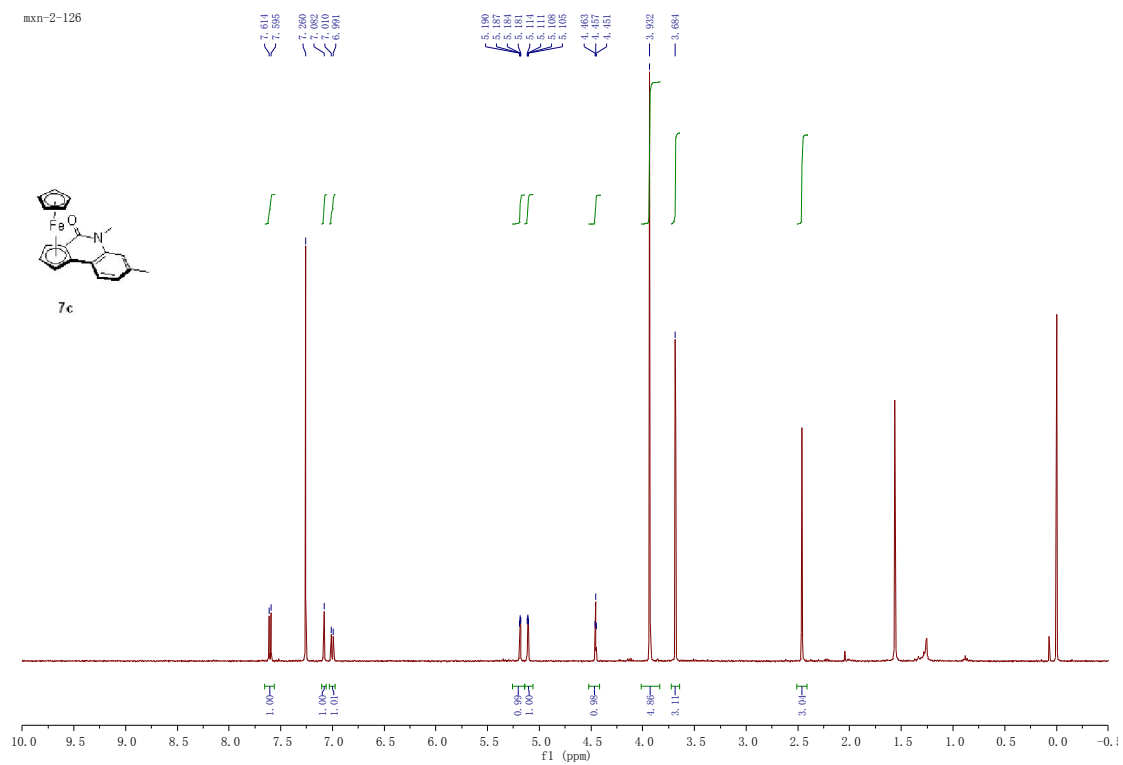
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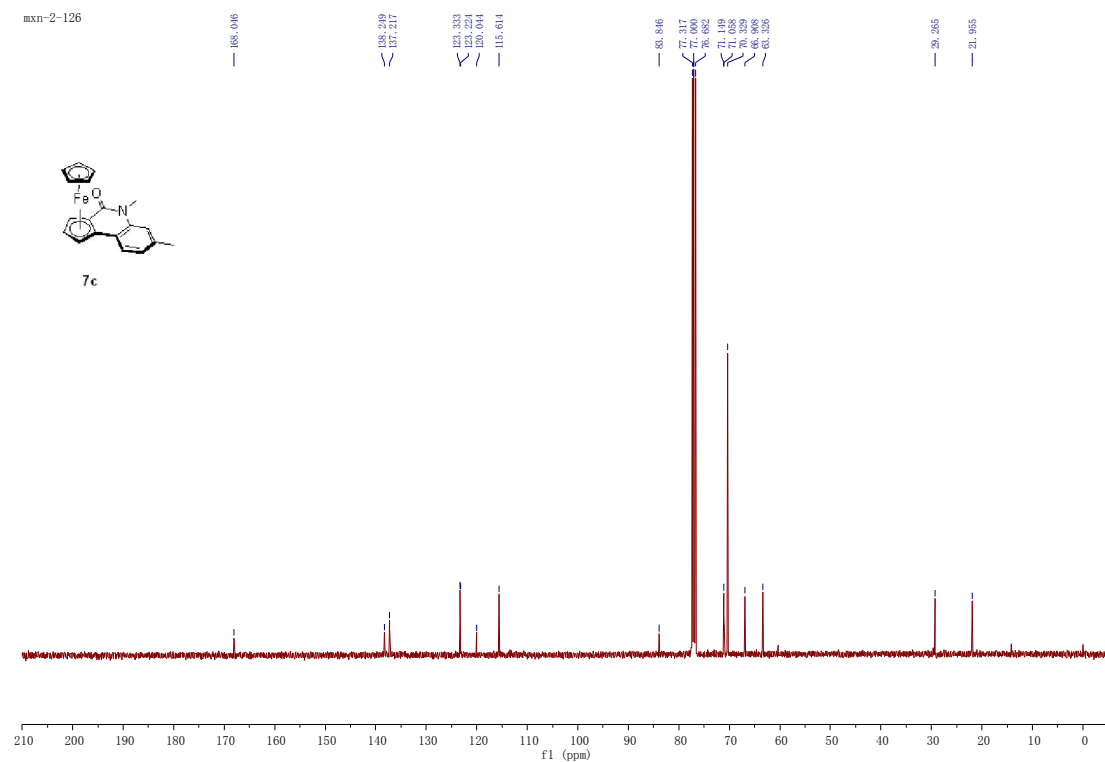




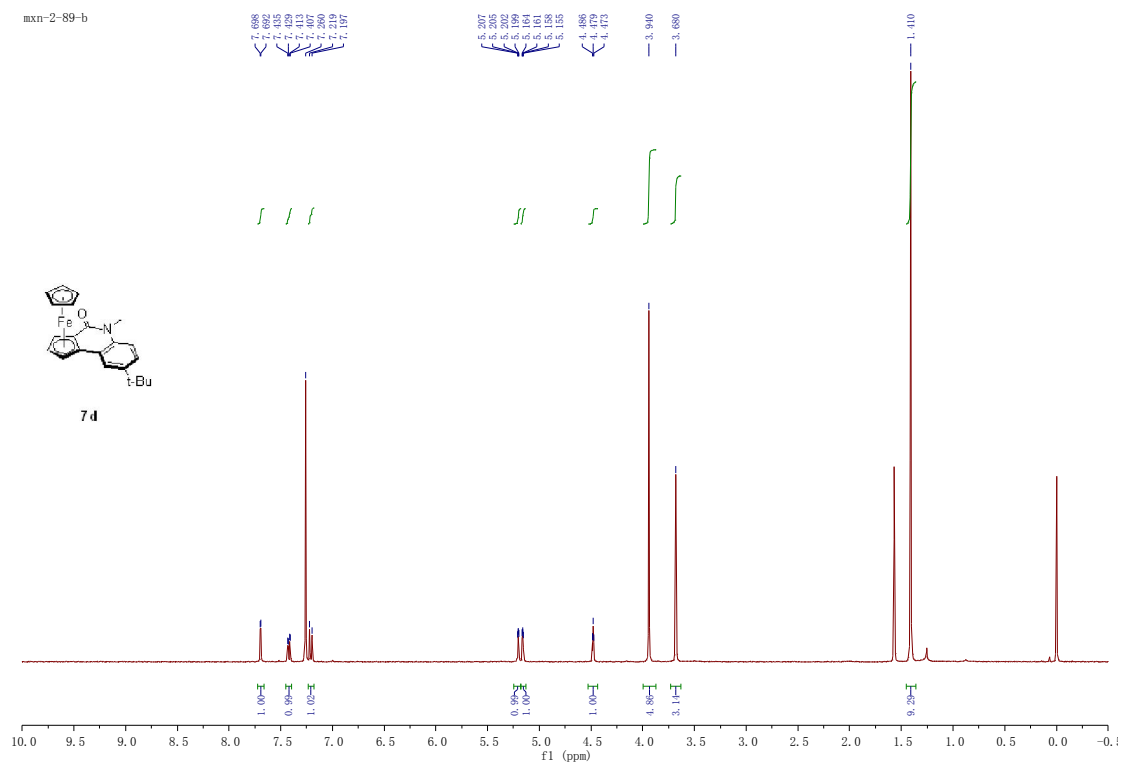
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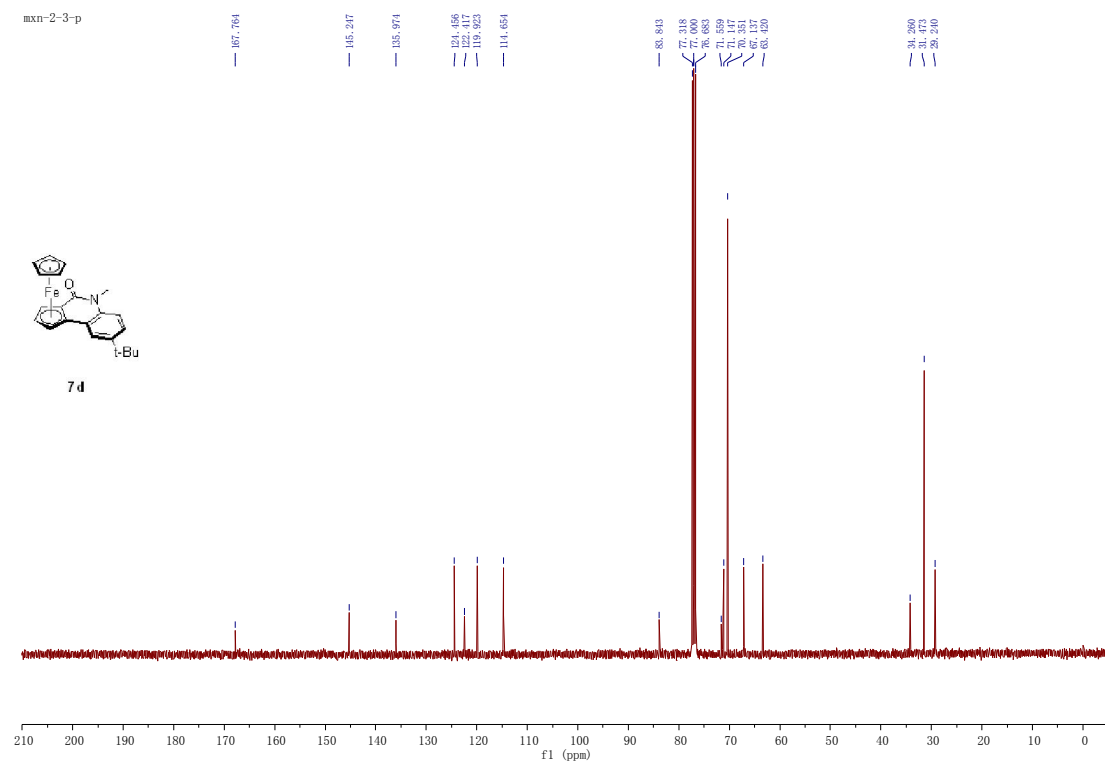
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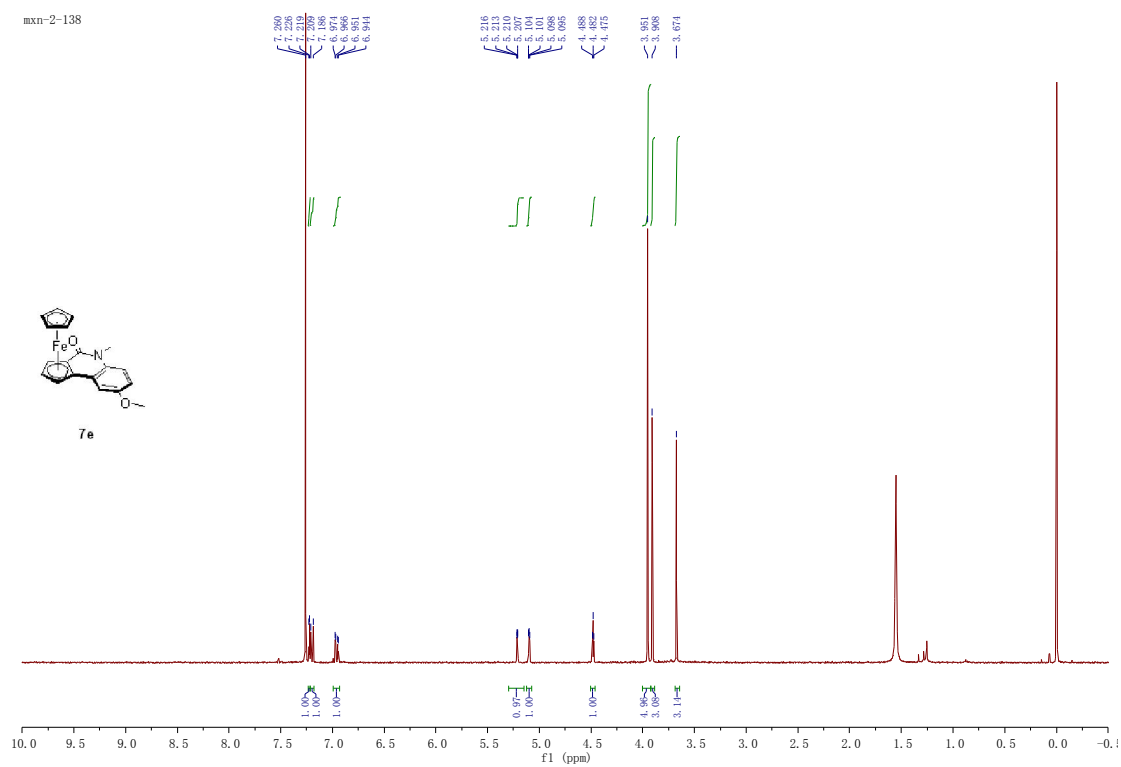
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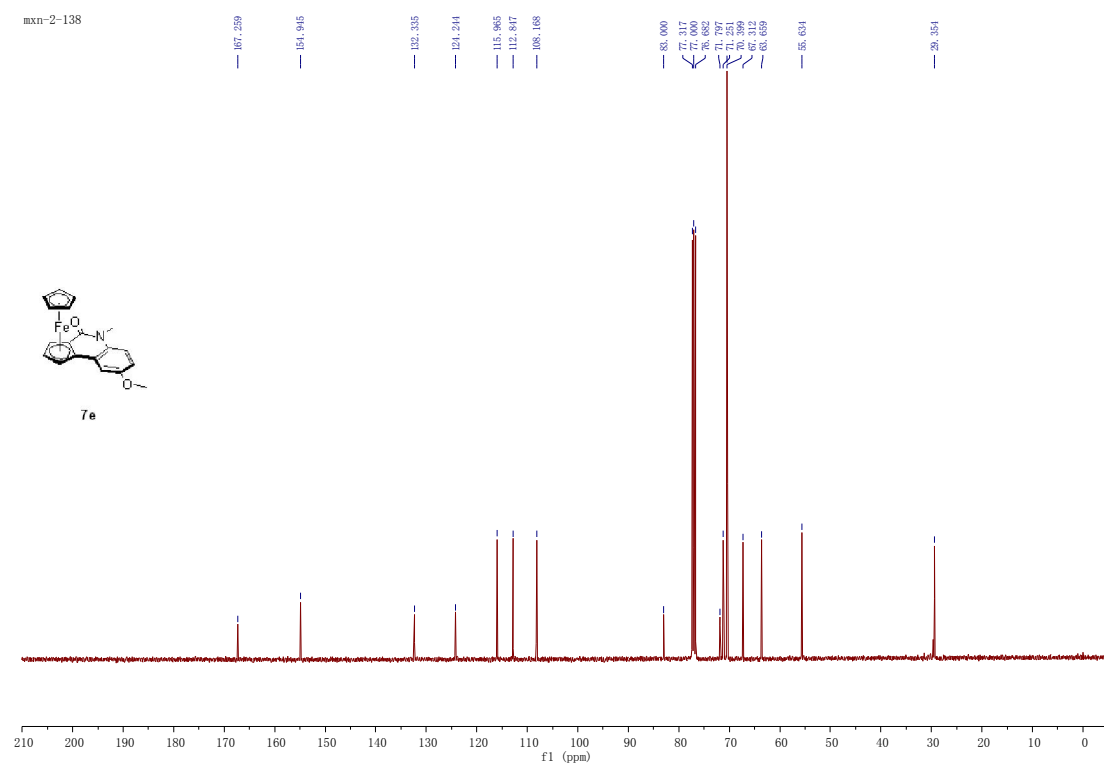
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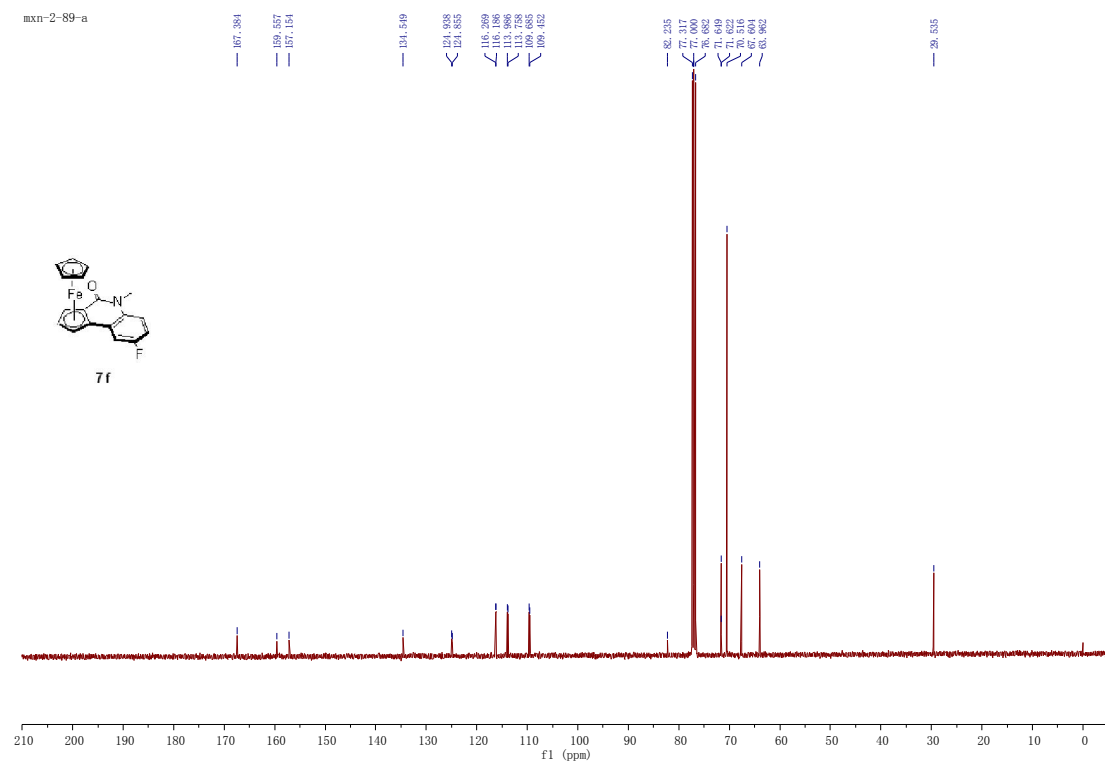
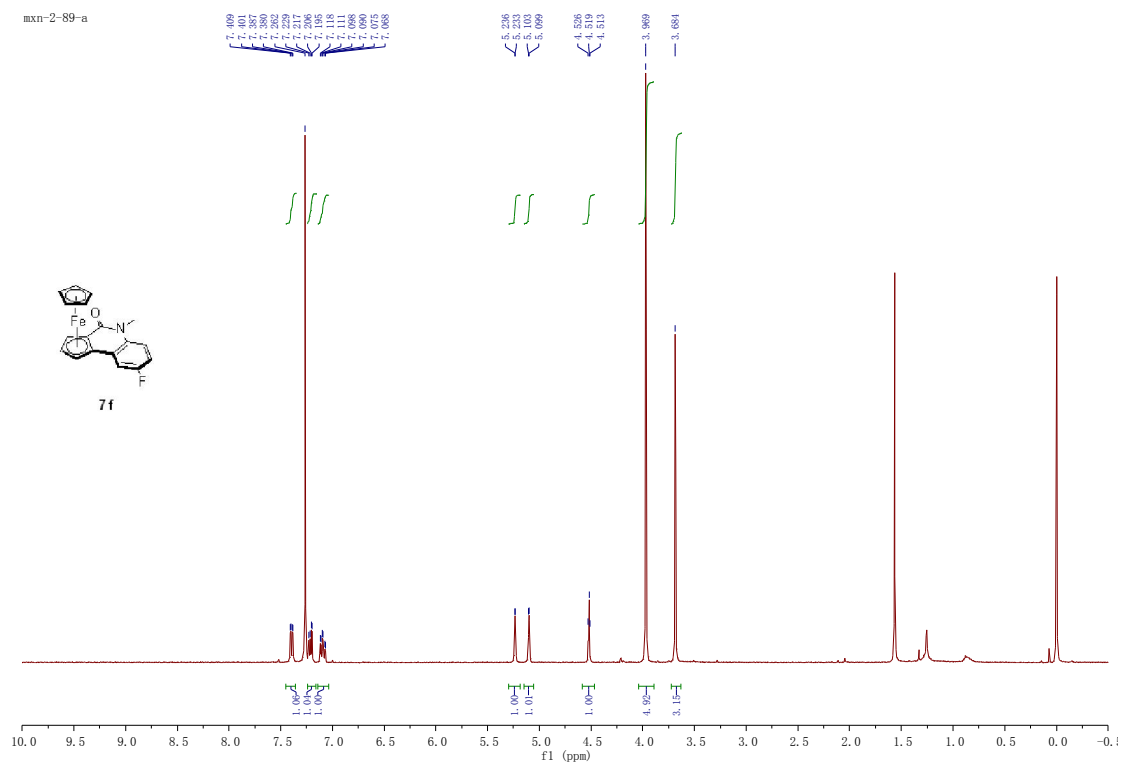


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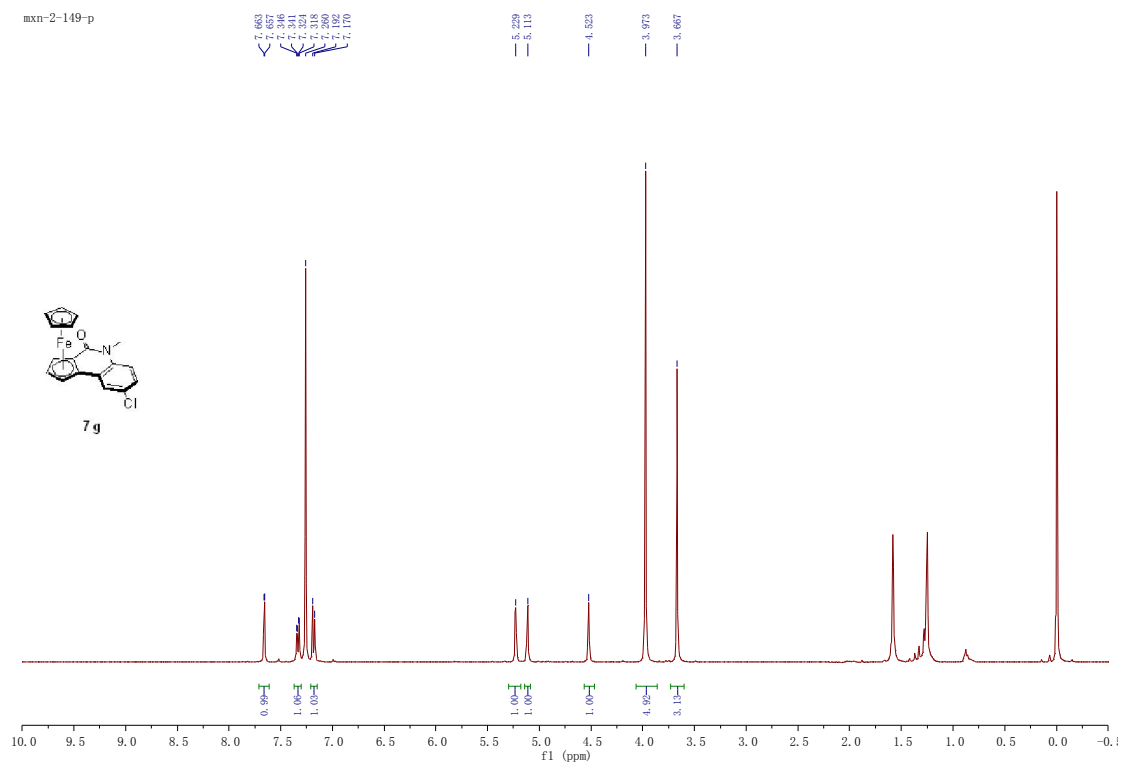


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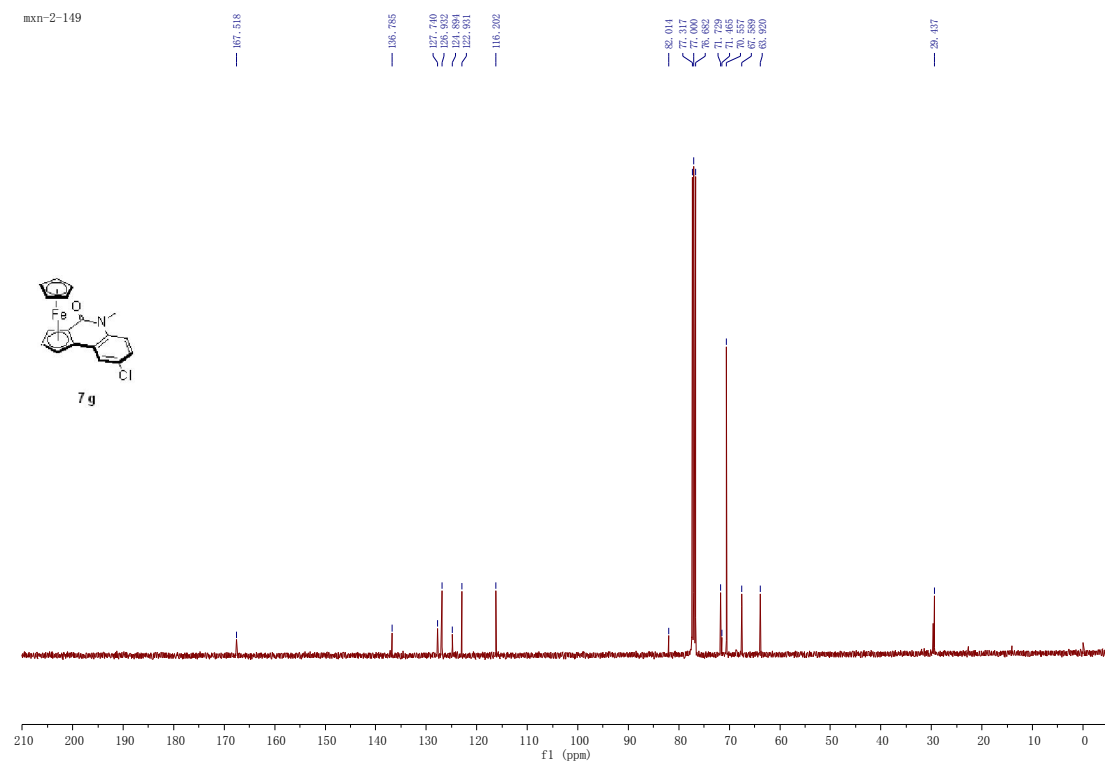


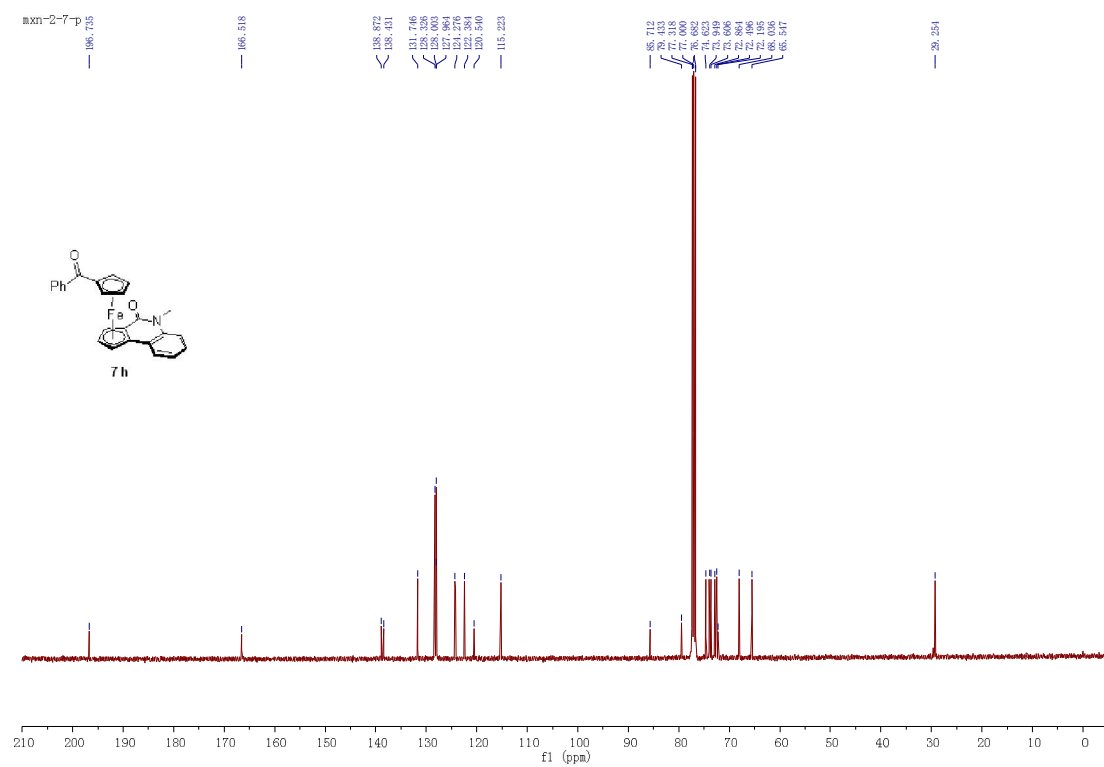
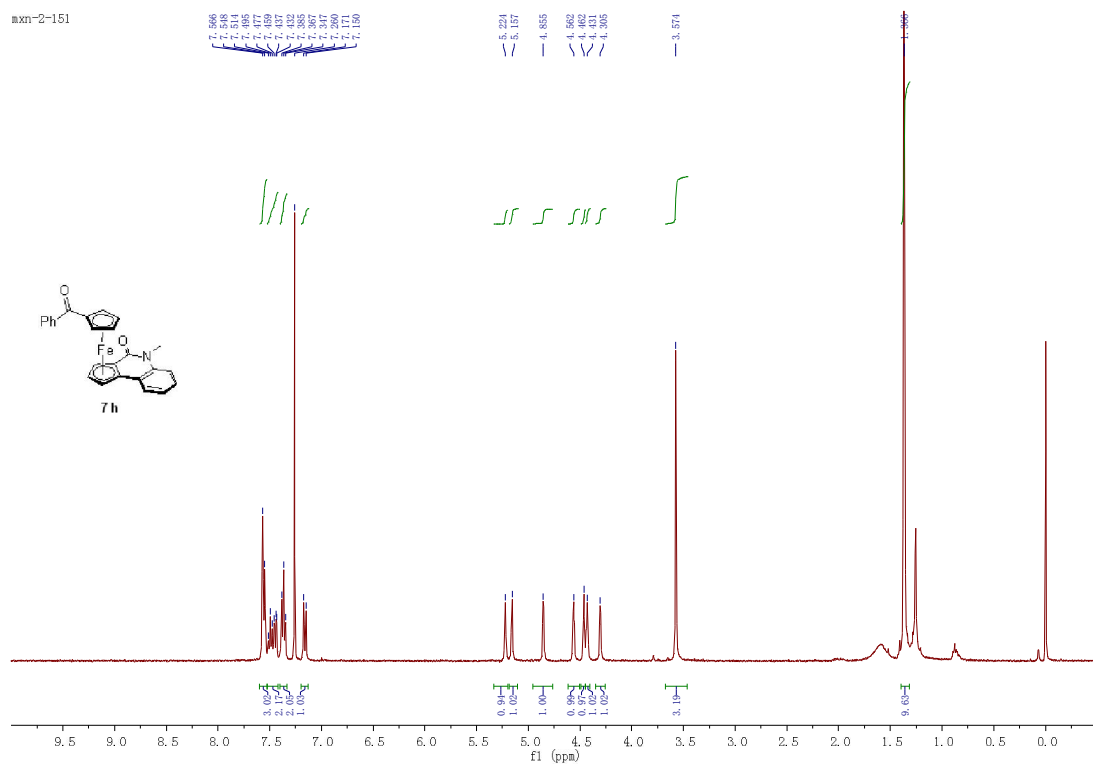


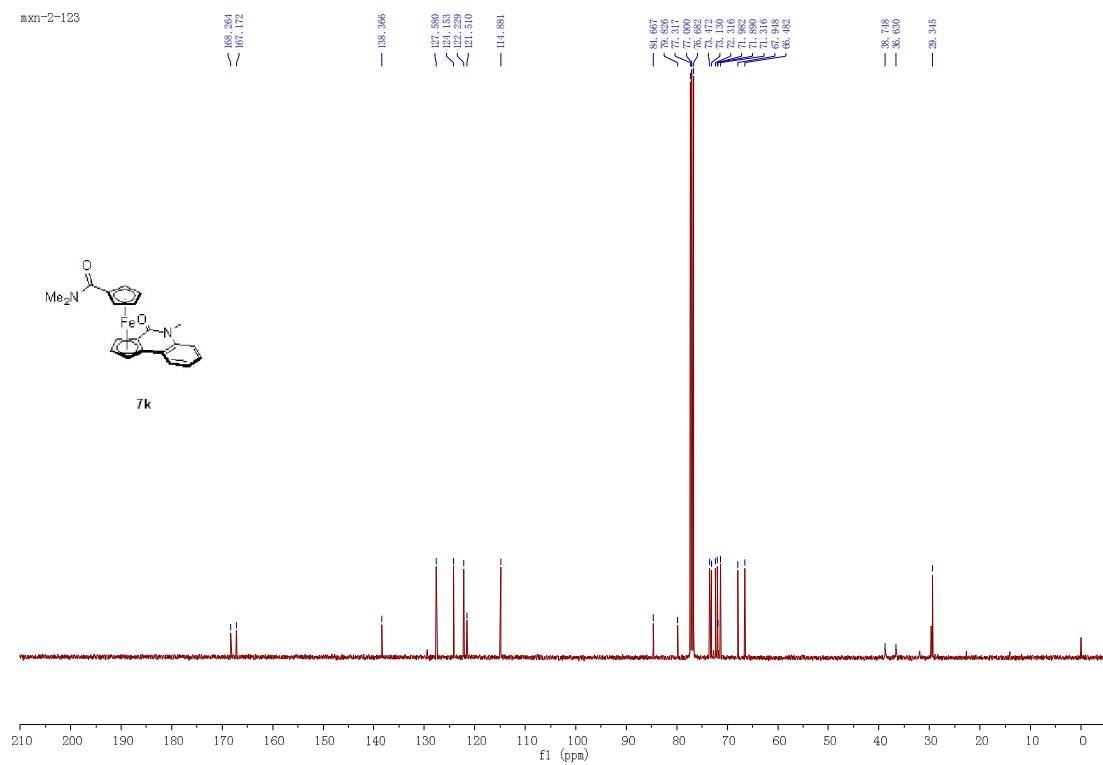
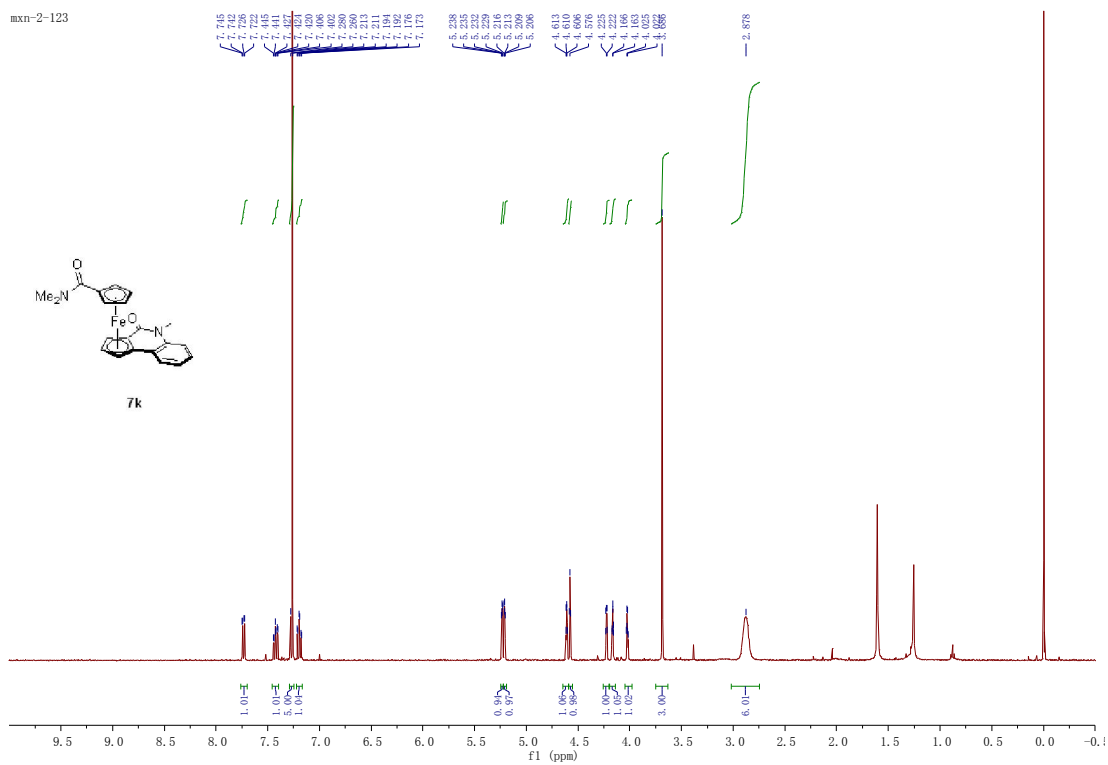
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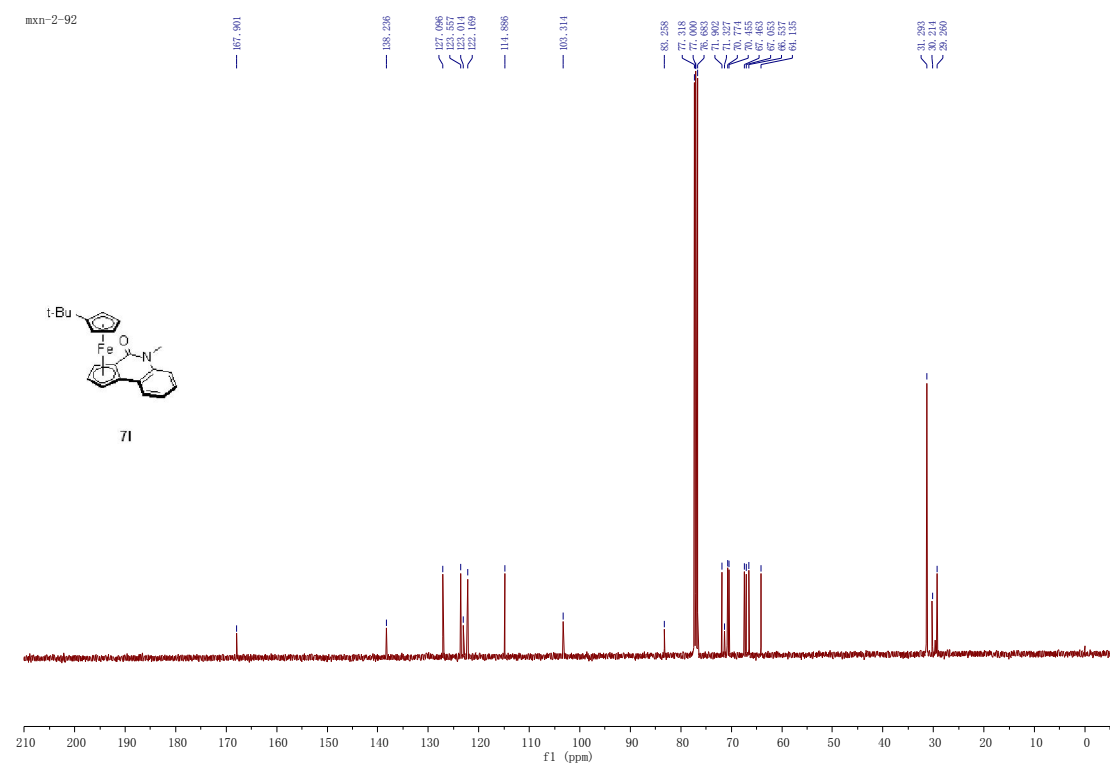
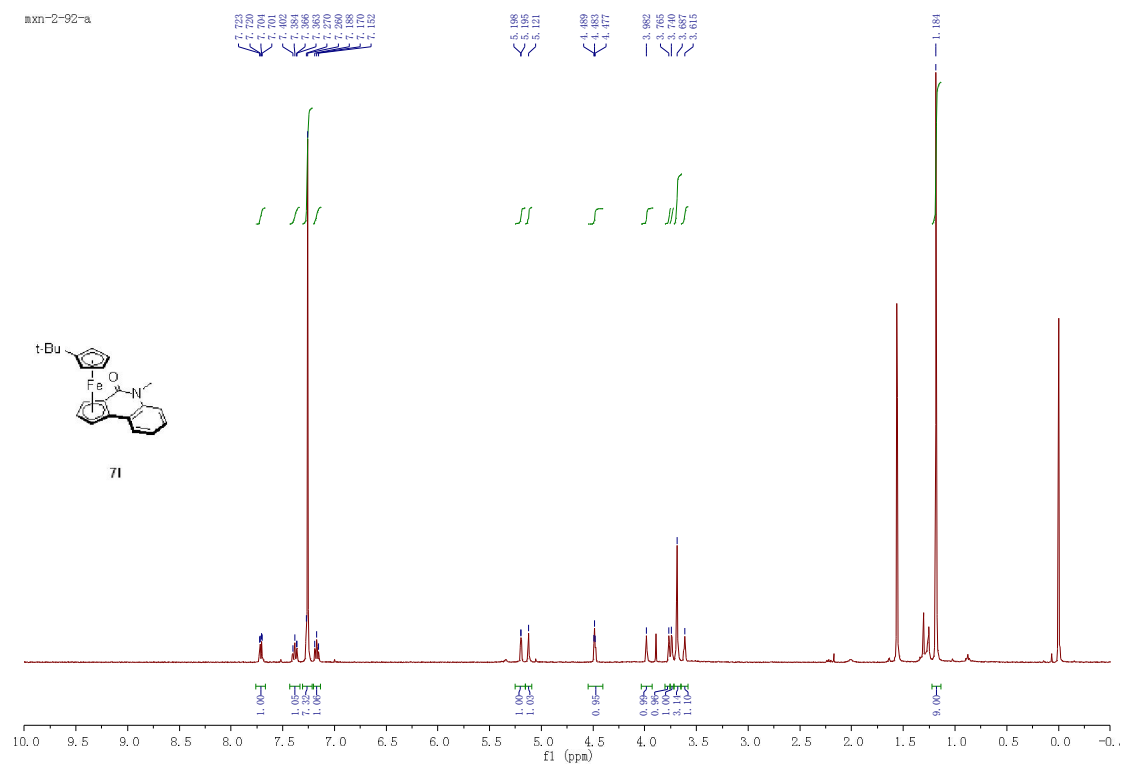


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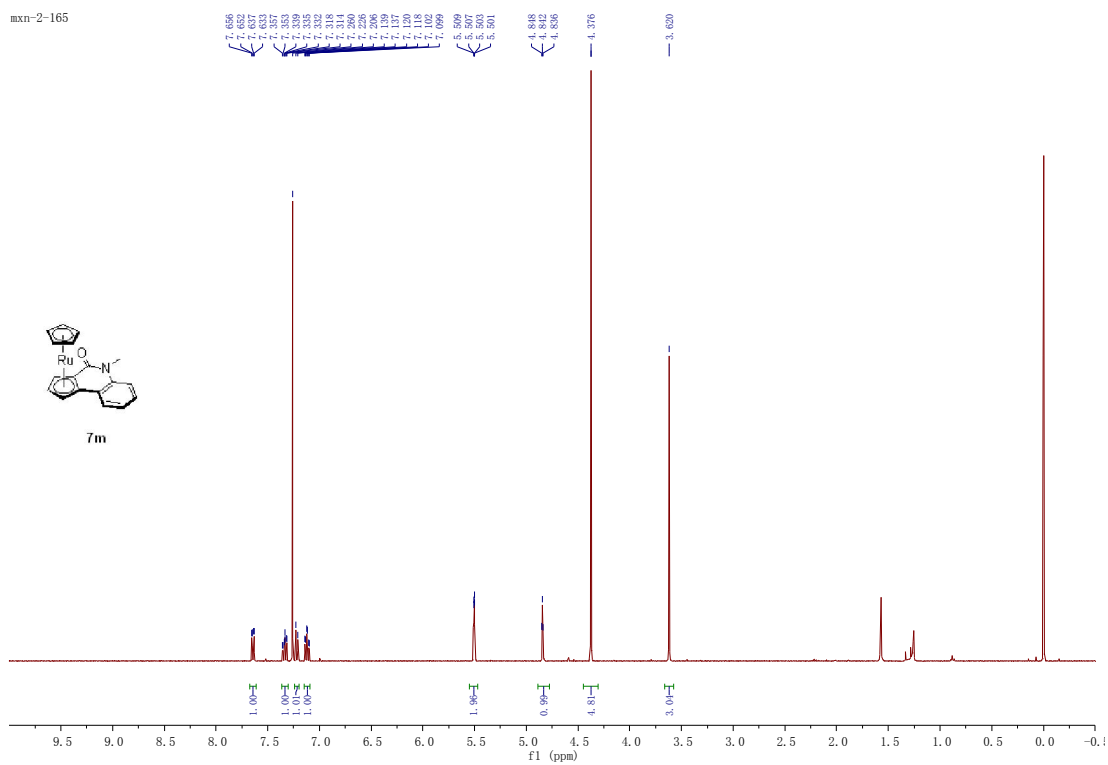




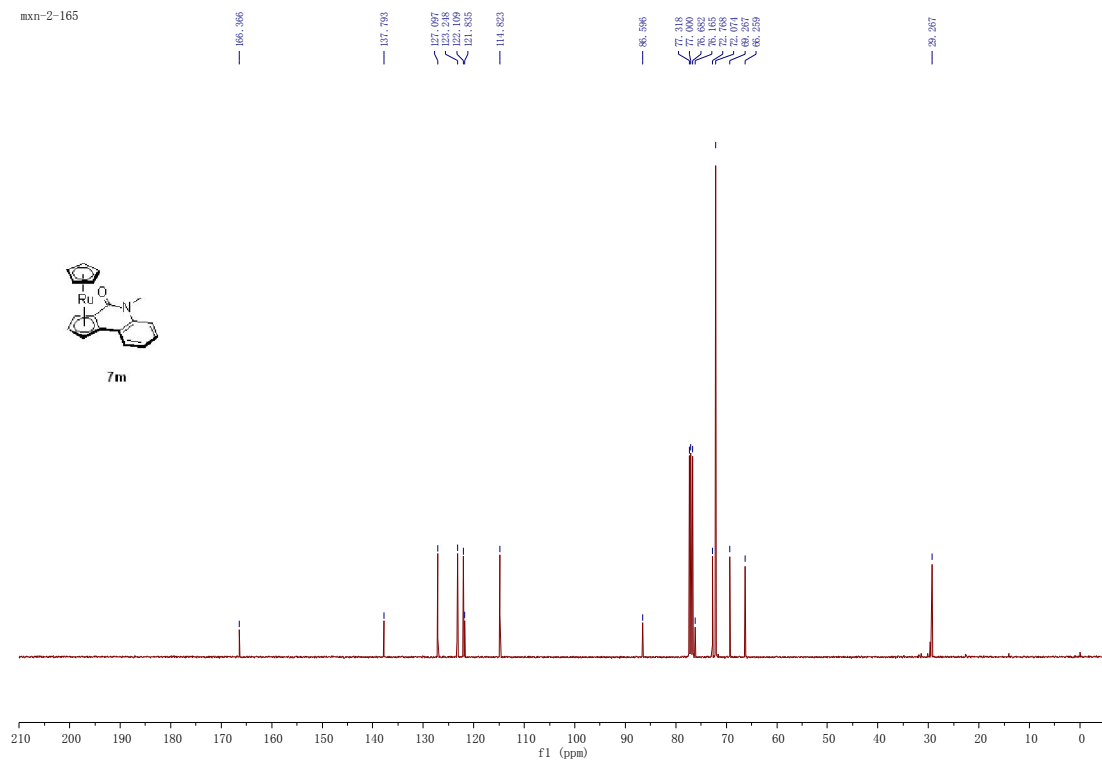




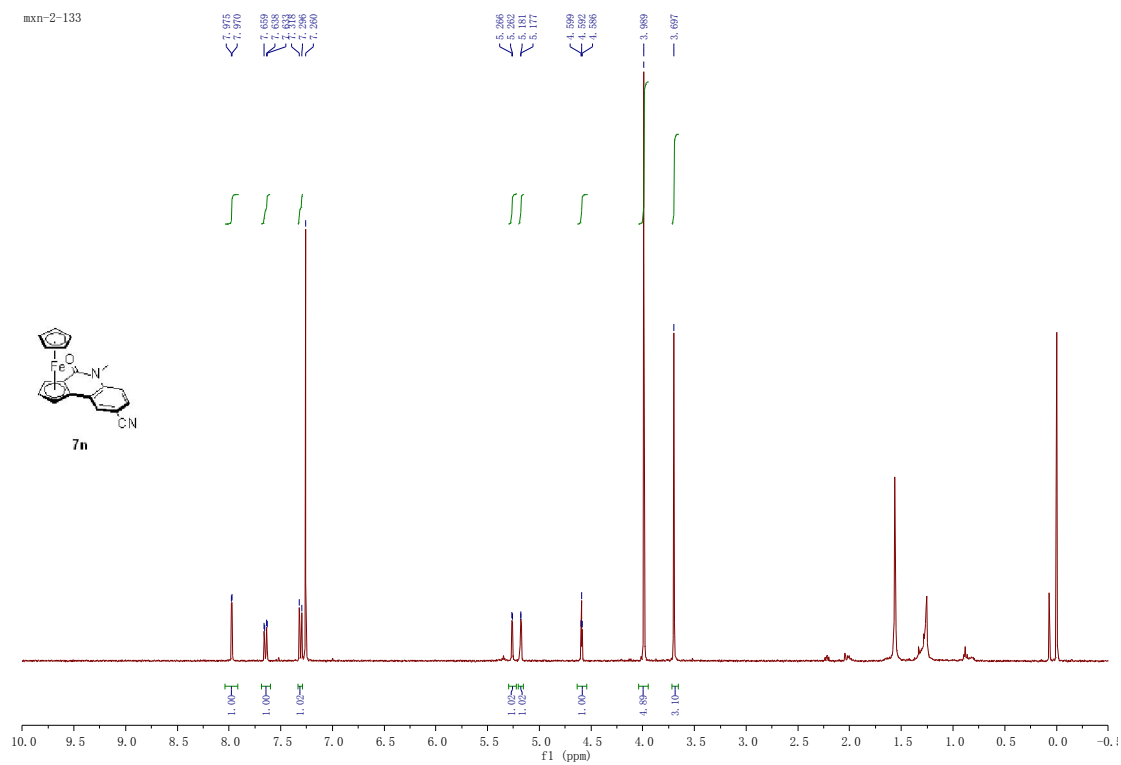
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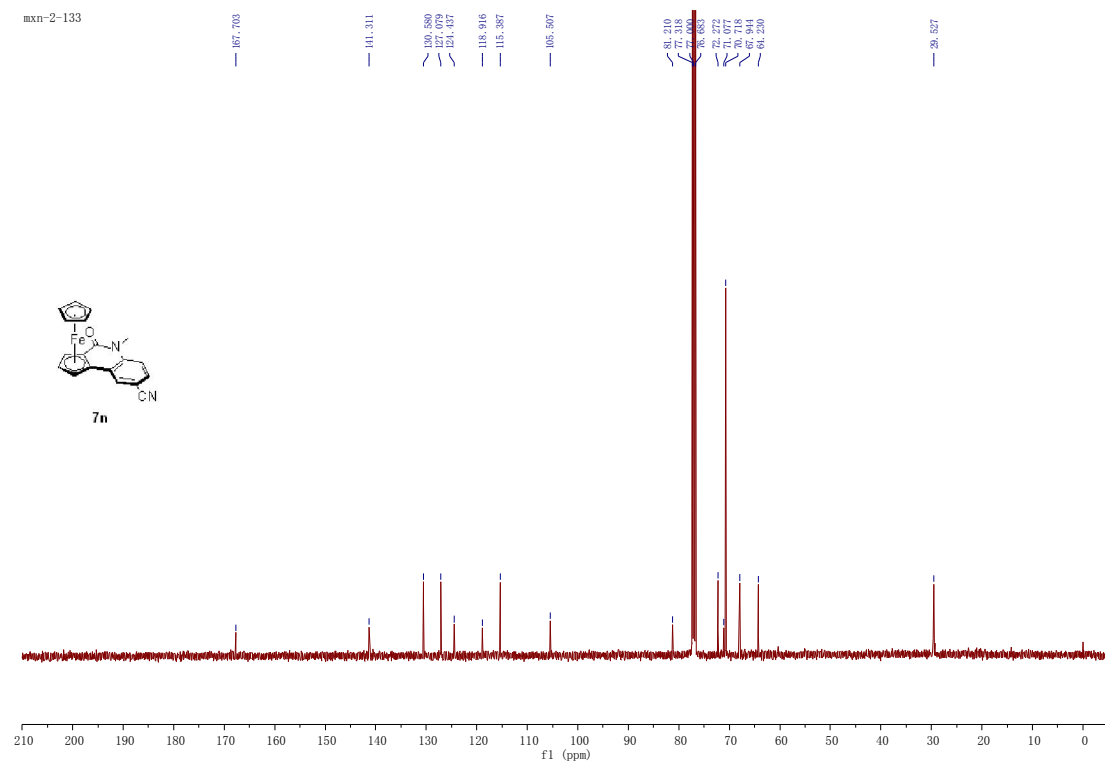
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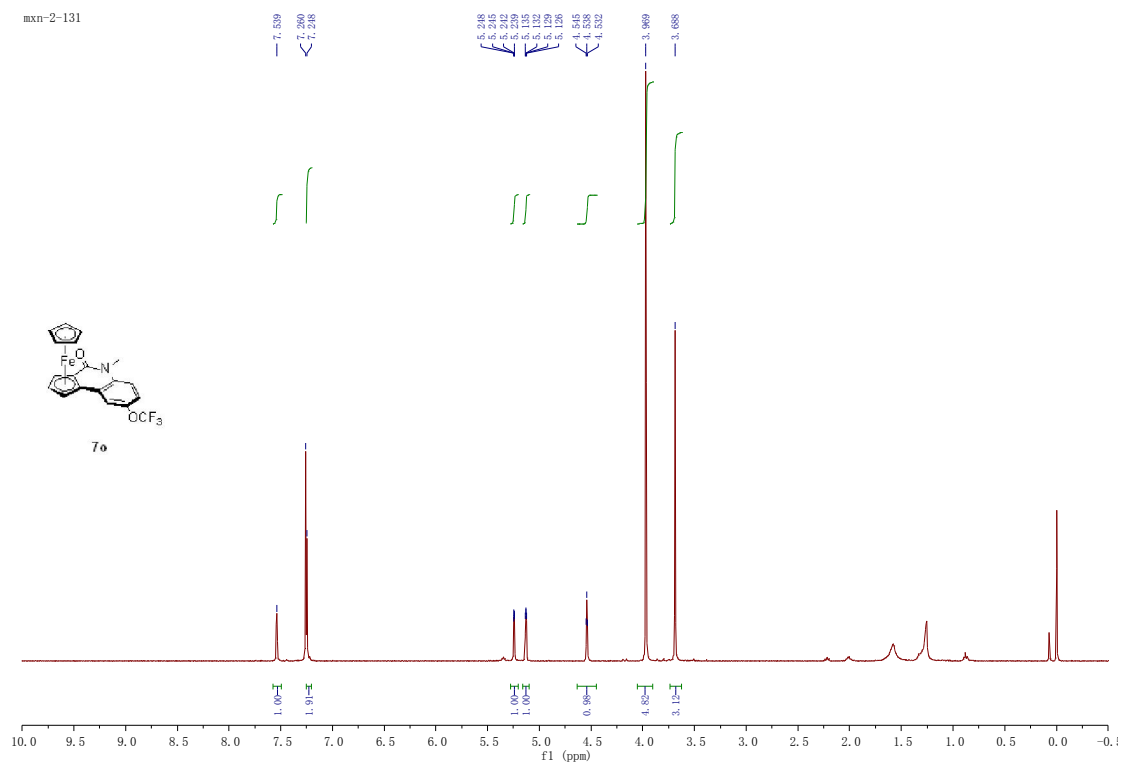
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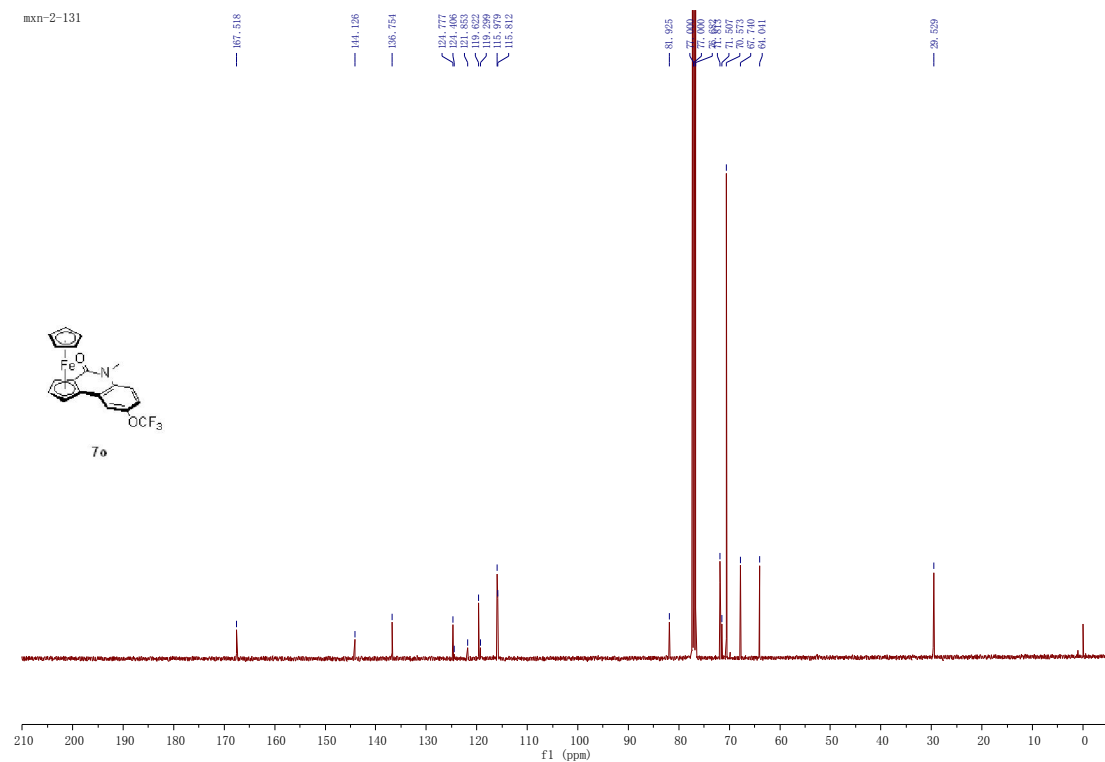
mxn-2-133



mxn-2-131



mxn-2-131



D-2000: maxinna Series: 0038

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-01-22

Reported Date and Time: 2014-06-02

17:36

20:38

Processed Date and Time: 2014-06-02

20:38

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0038\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0038

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-l-p-rac

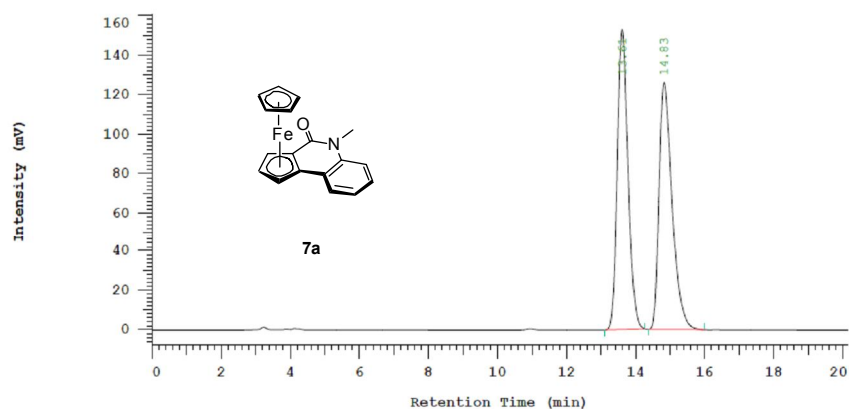
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	13.61	3316666	50.051	BB
2	14.83	3309967	49.949	BB
		6626633	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0039

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-01-22
18:05

Reported Date and Time: 2014-06-02
20:36

Processed Date and Time: 2014-06-02
20:35

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0039\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0039

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-l-p-rac

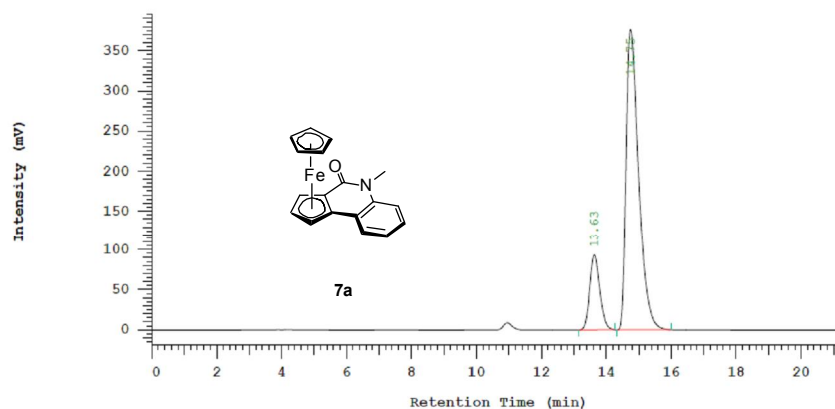
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	13.63	2014185	16.770	BB
2	14.75	9996677	83.230	BB
		12010862	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-04-14

Reported Date and Time: 2014-06-02

22:16

20:43

Processed Date and Time: 2014-06-02

20:43

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0055\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0055

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-2-rac

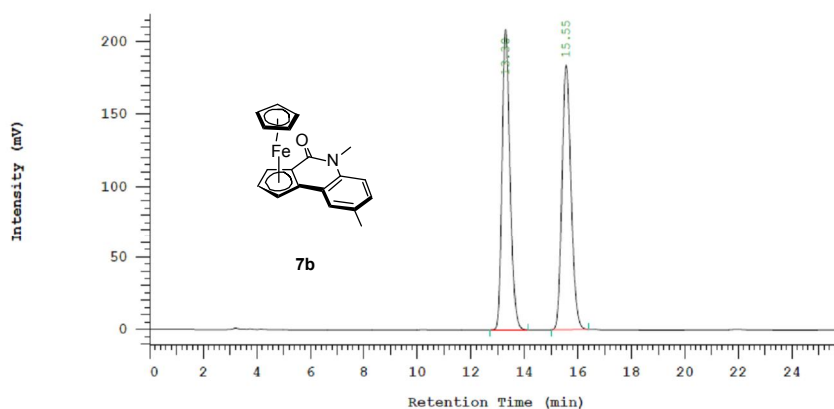
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	13.30	4321282	50.047	BB
2	15.55	4313176	49.953	BB
		8634458	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0056 Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-04-14 22:44 Reported Date and Time: 2014-06-02 20:41

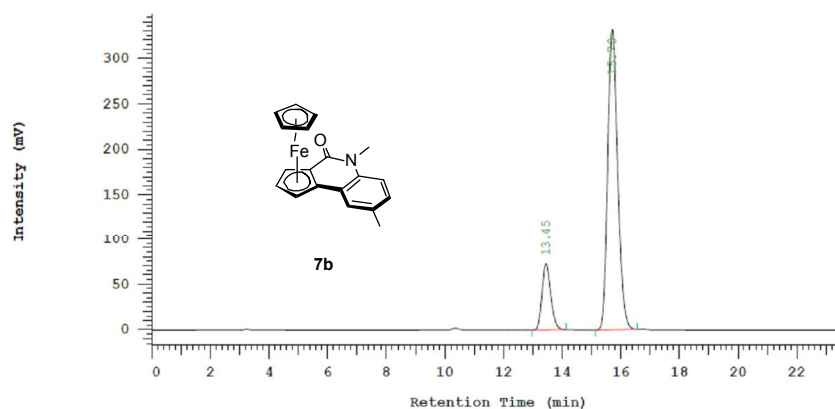
Processed Date and Time: 2014-06-02 20:40

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0056\

Processing Method: isopropanol

System (acquisition): LCE Series: 0056
Application(data): maxinna Vial Number: 1
Sample Name: mxn-2-2-rac Vial Type: UNK
Injection from this vial: 1 of 1 Volume: 10.0 ul
Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	13.45	1485346	16.038	BB
2	15.70	7776032	83.962	BB
		9261378	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-03-26

Reported Date and Time: 2014-06-02

11:04

21:17

Processed Date and Time: 2014-06-02

21:17

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0048\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0048

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-126-rac-1

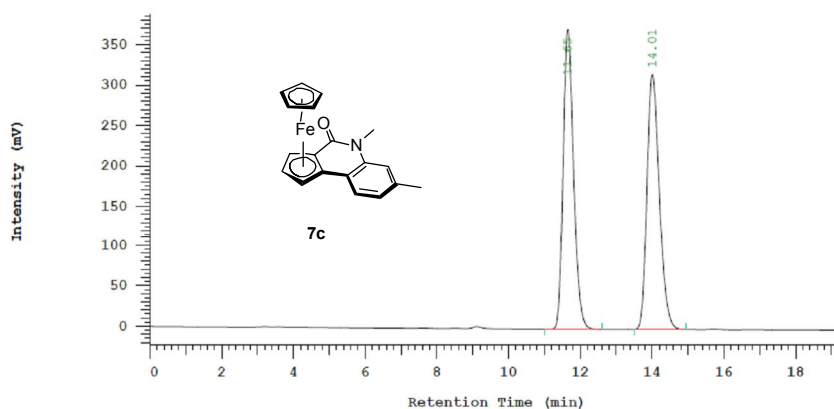
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	11.65	7568147	50.015	BB
2	14.01	7563459	49.985	BB
		15131606	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0049

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-03-26

Reported Date and Time: 2014-06-02

11:27

21:15

Processed Date and Time: 2014-06-02

21:15

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0049\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0049

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-126-rac-1

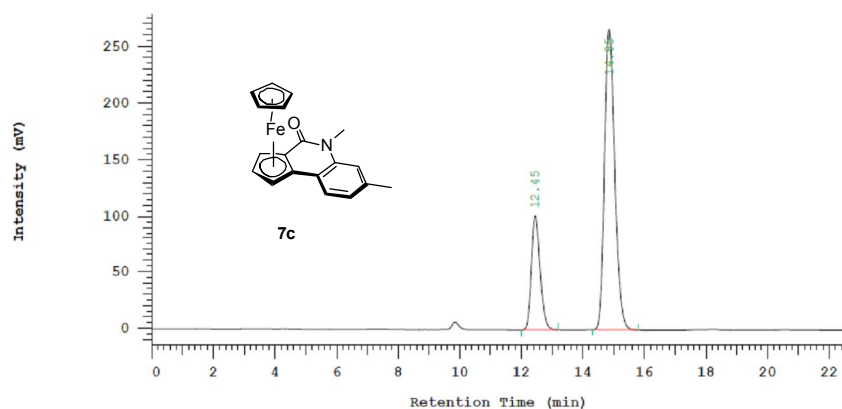
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	12.45	1978784	24.540	BB
2	14.85	6084881	75.460	BB
		8063665	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0036

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2013-12-30

Reported Date and Time: 2014-06-02

21:50

20:45

Processed Date and Time: 2014-06-02

20:45

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0036\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0036

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-89-b-rac

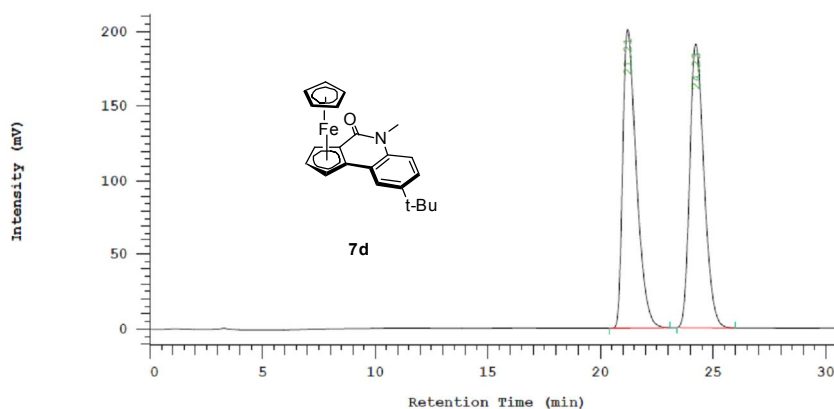
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	21.21	8383408	50.039	BB
2	24.23	8370294	49.961	BB
		16753702	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0037

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2013-12-30

Reported Date and Time: 2014-06-02

22:24

20:48

Processed Date and Time: 2014-06-02

20:47

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0037\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0037

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-89-b-rac

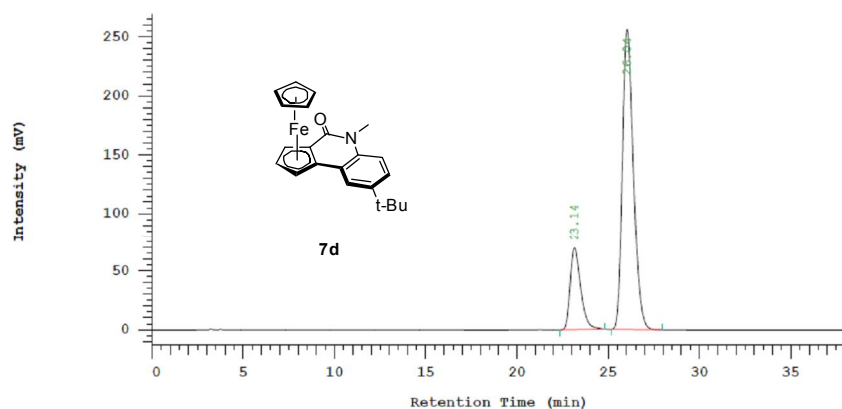
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	23.14	2760809	20.526	BB
2	26.04	10689787	79.474	BB
		13450596	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-04-15 22:53 Reported Date and Time: 2014-06-02 21:28

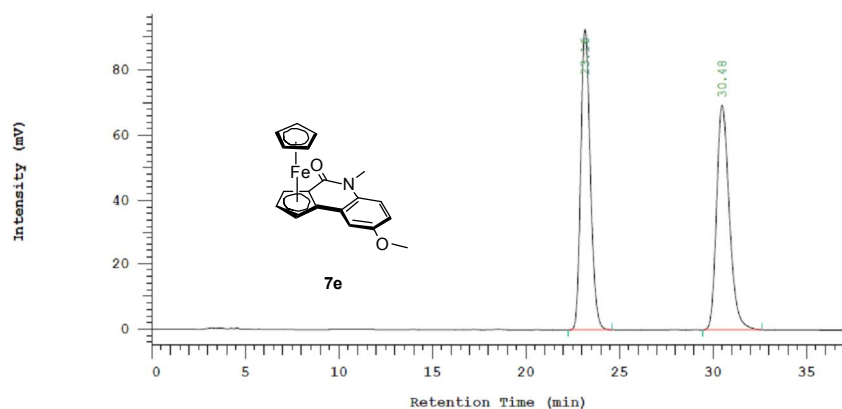
Processed Date and Time: 2014-06-02 21:28

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0065\

Processing Method: isopropanol

System (acquisition): LCE Series: 0065
Application(data): maxinna Vial Number: 1
Sample Name: mxn-2-138 Vial Type: UNK
Injection from this vial: 1 of 1 Volume: 10.0 ul
Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	23.16	3207038	49.772	BB
2	30.48	3236478	50.228	BB
		6443516	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-04-15

Reported Date and Time: 2014-06-02

23:33

21:28

Processed Date and Time: 2014-06-02

21:27

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0066\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0066

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-138

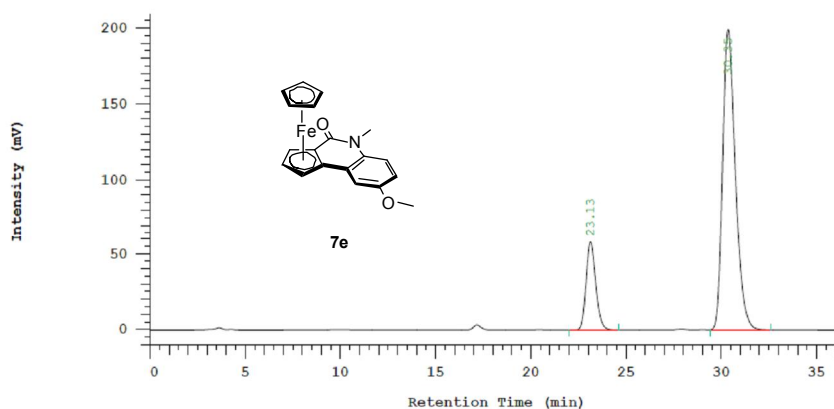
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	23.13	2020216	17.921	BB
2	30.35	9252836	82.079	BB
		11273052	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0032

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2013-12-30

Reported Date and Time: 2014-06-02

17:09

20:54

Processed Date and Time: 2014-06-02

20:53

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0032\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0032

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-89-a-rac

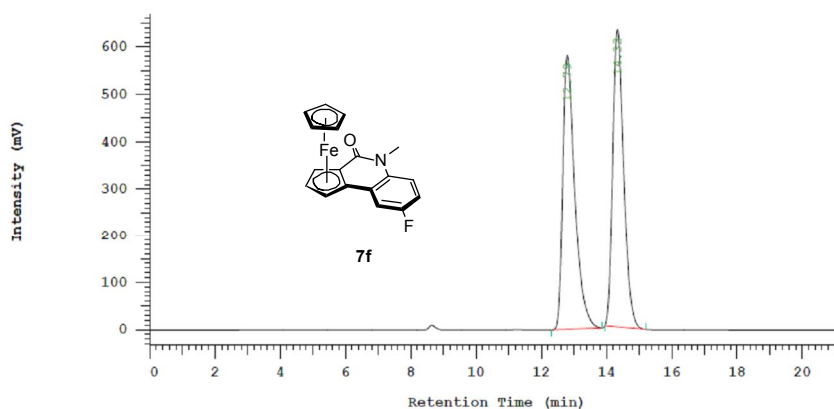
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	12.79	14945224	50.167	BB
2	14.33	14845587	49.833	BB
		29790811	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2013-12-30 17:36 Reported Date and Time: 2014-06-02 20:53

Processed Date and Time: 2014-06-02 20:51

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0033_001\

Processing Method: isopropanol

System (acquisition): LCE Series: 0033_001

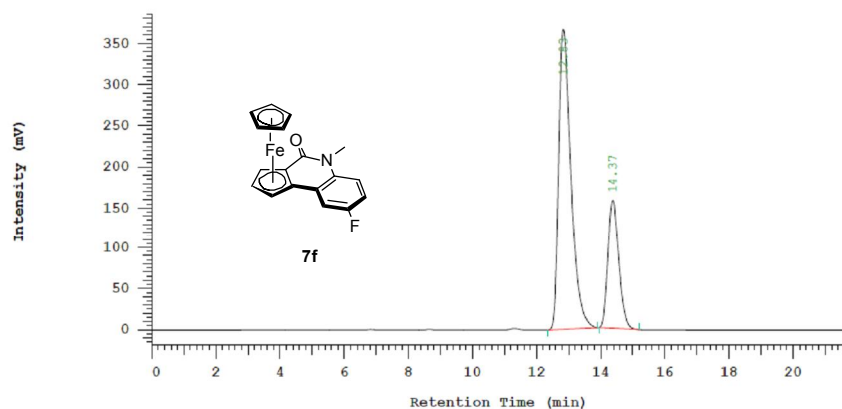
Application(data): maxinna Vial Number: 1

Sample Name: mxn-2-89-a-rac Vial Type: UNK

Injection from this vial: 1 of 1 Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	12.83	9650331	72.328	BB
2	14.37	3692179	27.672	BB
		13342510	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0058

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-04-14

Reported Date and Time: 2014-06-02

23:47

20:58

Processed Date and Time: 2014-06-02

20:57

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0058\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0058

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-5

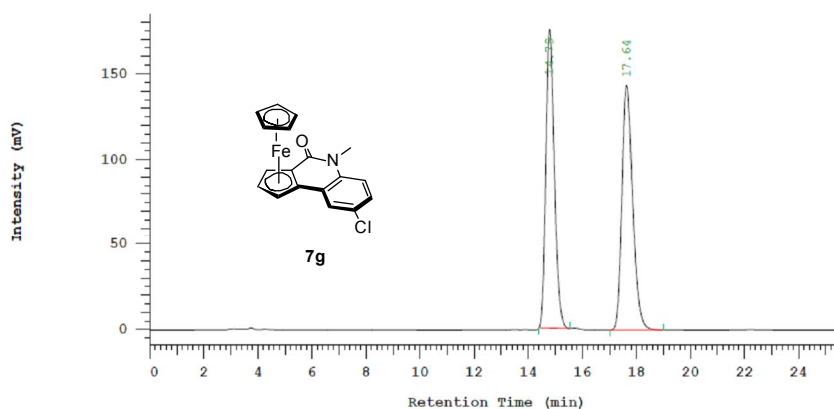
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	14.79	3838877	49.469	BB
2	17.64	3921306	50.531	BB
		7760183	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-04-15 11:35 Reported Date and Time: 2014-06-02 20:56

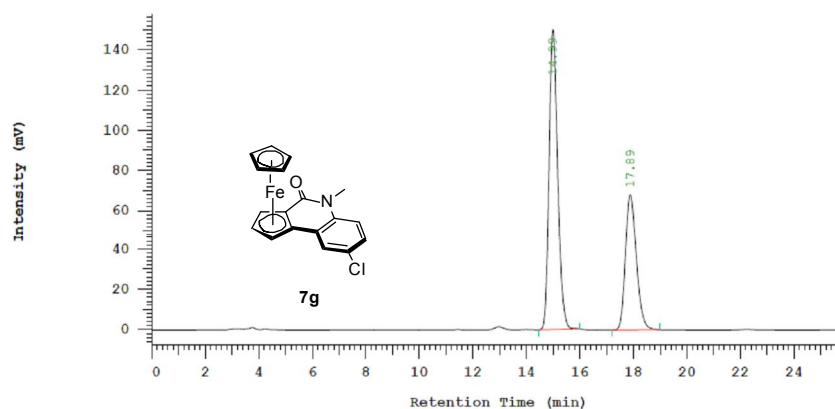
Processed Date and Time: 2014-06-02 20:55

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0061\

Processing Method: isopropanol

System (acquisition): LCE Series: 0061
 Application(data): maxinna Vial Number: 1
 Sample Name: mxn-2-5 Vial Type: UNK
 Injection from this vial: 1 of 1 Volume: 10.0 ul
 Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	14.99	3359905	63.928	BB
2	17.89	1895860	36.072	BB
		5255765	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-04-17 12:18 Reported Date and Time: 2014-06-02 21:07

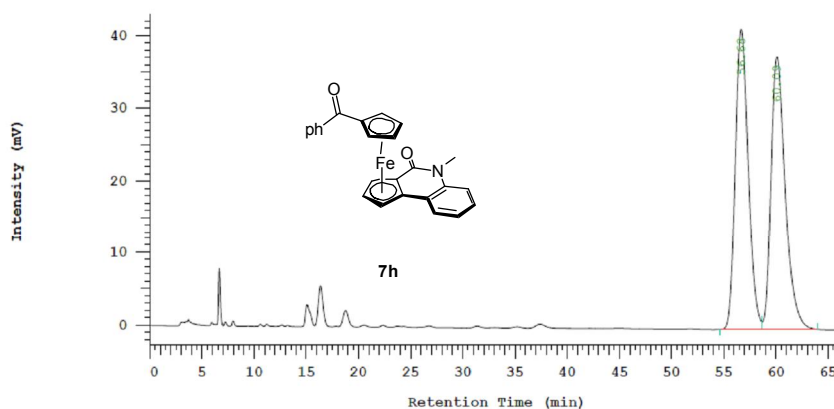
Processed Date and Time: 2014-06-02 21:07

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0071\

Processing Method: isopropanol

System (acquisition): LCE Series: 0071
 Application(data): maxinna Vial Number: 1
 Sample Name: mxn-2-7 Vial Type: UNK
 Injection from this vial: 1 of 1 Volume: 10.0 ul
 Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	56.68	3523887	49.839	BV
2	60.09	3546699	50.161	VB
		7070586	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0072

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-04-17

Reported Date and Time: 2014-06-02

13:38

21:05

Processed Date and Time: 2014-06-02

21:04

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0072\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0072

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-7

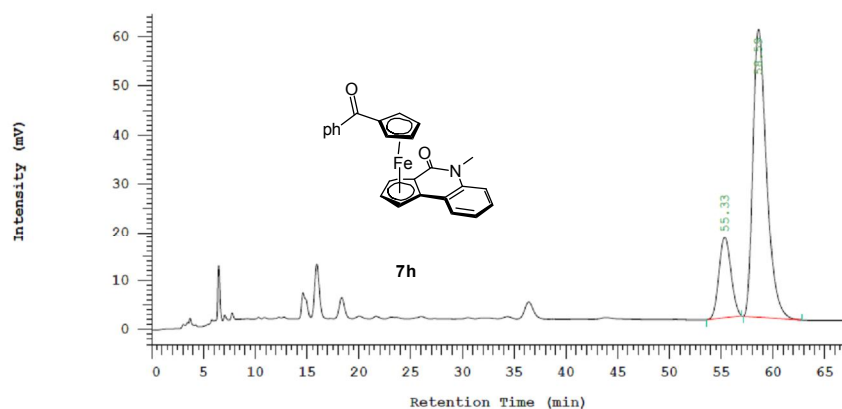
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	55.33	1301946	19.479	BB
2	58.59	5381908	80.521	BB
		6683854	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0115

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-05-06

Reported Date and Time: 2014-06-02

16:46

21:47

Processed Date and Time: 2014-06-02

21:46

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0115\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0115

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-151

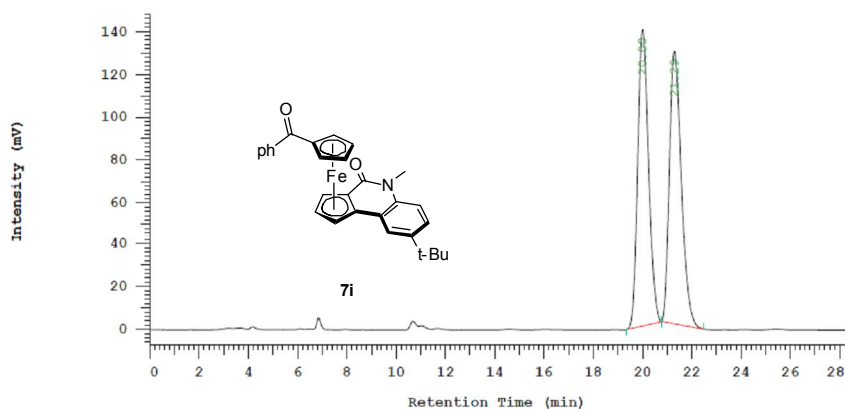
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	20.00	4264319	50.055	BB
2	21.29	4254999	49.945	BB
		8519318	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-05-06 17:15 Reported Date and Time: 2014-06-02 21:45

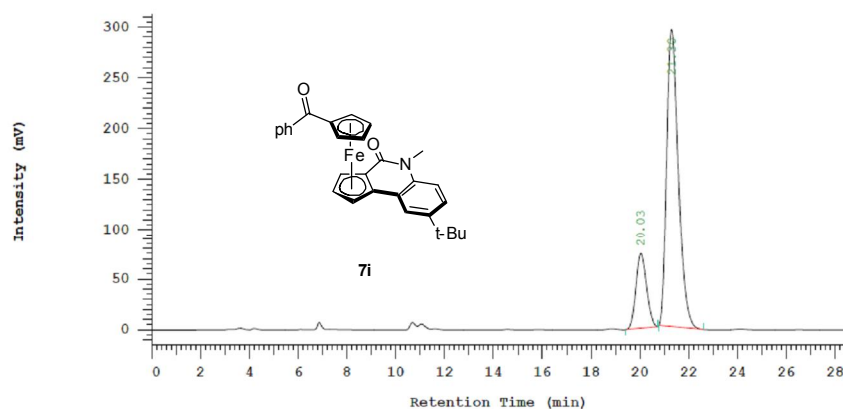
Processed Date and Time: 2014-06-02 21:44

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0116\

Processing Method: isopropanol

System (acquisition): LCE Series: 0116
 Application(data): maxinna Vial Number: 1
 Sample Name: mxn-2-151 Vial Type: UNK
 Injection from this vial: 1 of 1 Volume: 10.0 ul
 Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	20.03	2216919	18.501	BB
2	21.30	9765971	81.499	BB
		11982890	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-04-15

Reported Date and Time: 2014-06-02

16:52

21:10

Processed Date and Time: 2014-06-02

21:10

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0062\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0062

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-108-b

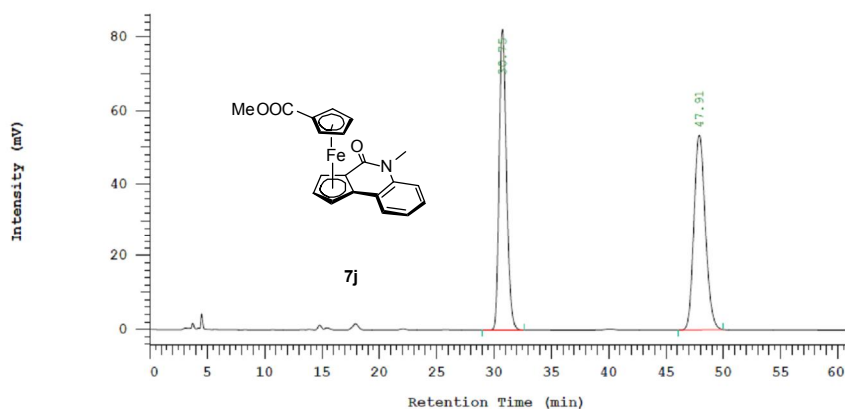
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	30.75	3616373	50.061	BB
2	47.91	3607591	49.939	BB
		7223964	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-04-15 18:48 Reported Date and Time: 2014-06-02 21:09

Processed Date and Time: 2014-06-02
21:09

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0064\

Processing Method: isopropanol

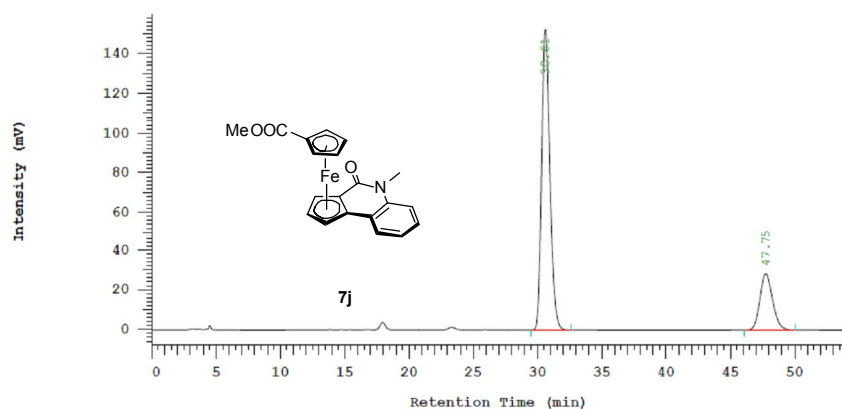
Series: 0064

Vial Number: 1

Vial Type: UNK

Volume: 10.0 ul

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	30.61	6656194	77.689	BB
2	47.75	1911501	22.311	BB
		8567695	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-03-15

Reported Date and Time: 2014-06-02

17:46

21:14

Processed Date and Time: 2014-06-02

21:13

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0046\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0046

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-122-rac-4

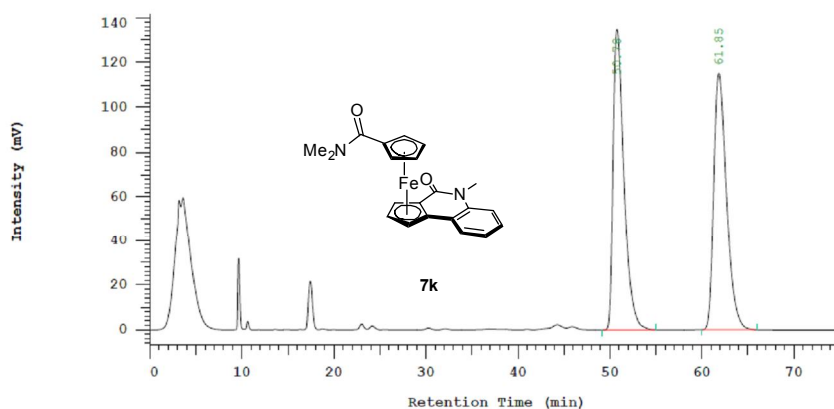
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	50.78	11020169	50.187	BB
2	61.85	10937936	49.813	BB
		21958105	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0047

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-03-15
19:09

Reported Date and Time: 2014-06-02
21:12

Processed Date and Time: 2014-06-02
21:12

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0047\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0047

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-122-rac-4

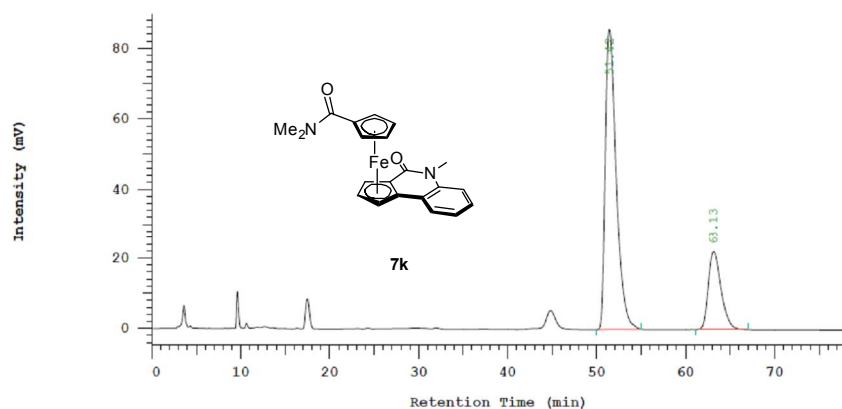
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	51.42	7151881	76.639	BB
2	63.13	2180042	23.361	BB
		9331923	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0040

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-01-22

Reported Date and Time: 2014-06-02

18:37

21:02

Processed Date and Time: 2014-06-02

21:02

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0040\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0040

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-1-p-rac

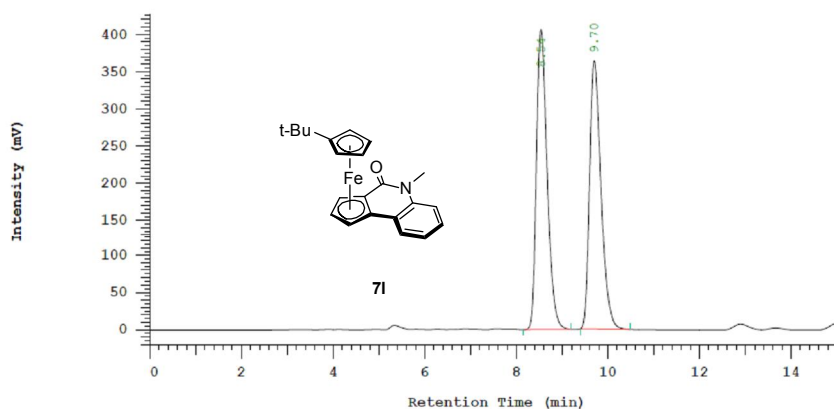
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	8.54	6511258	50.611	BB
2	9.70	6353925	49.389	BB
		12865183	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-01-22 18:58 Reported Date and Time: 2014-06-02 21:01

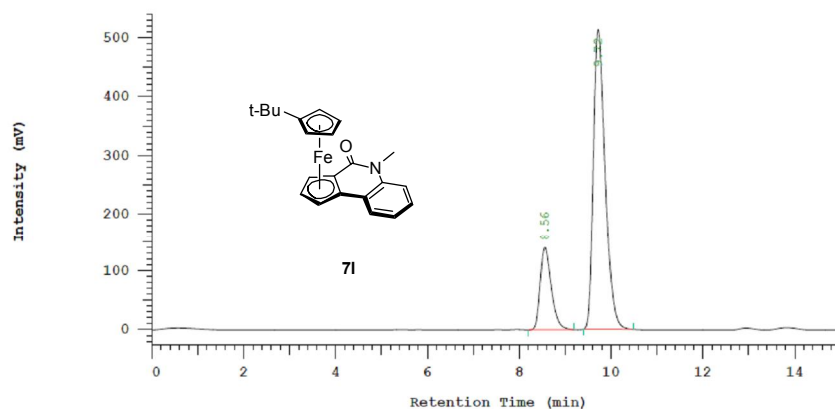
Processed Date and Time: 2014-06-02 21:00

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0041\

Processing Method: isopropanol

System (acquisition): LCE Series: 0041
Application(data): maxinna Vial Number: 1
Sample Name: mxn-2-l-p-rac Vial Type: UNK
Injection from this vial: 1 of 1 Volume: 10.0 ul
Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	8.56	2337420	20.238	BB
2	9.72	9212408	79.762	BB
		11549828	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-05-07 18:19 Reported Date and Time: 2014-06-02 21:54

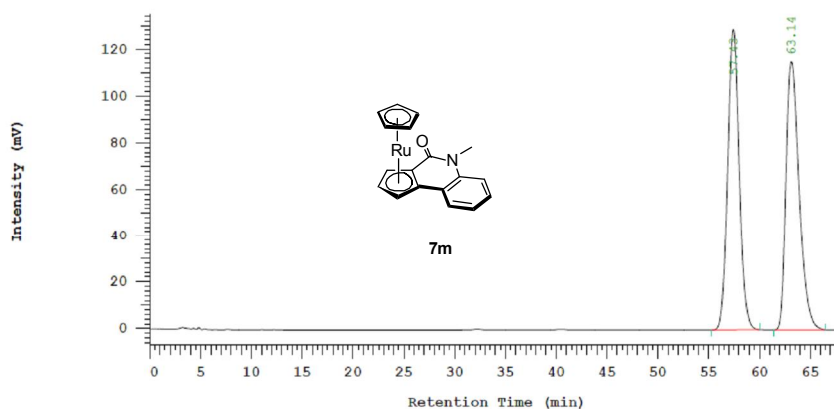
Processed Date and Time: 2014-06-02 21:53

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0121\

Processing Method: isopropanol

System (acquisition): LCE Series: 0121
 Application(data): maxinna Vial Number: 1
 Sample Name: mxn-2-165 Vial Type: UNK
 Injection from this vial: 1 of 1 Volume: 10.0 ul
 Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	57.43	10158084	50.001	BB
2	63.14	10157879	49.999	BB
		20315963	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-05-07 20:08 Reported Date and Time: 2014-06-02 21:53

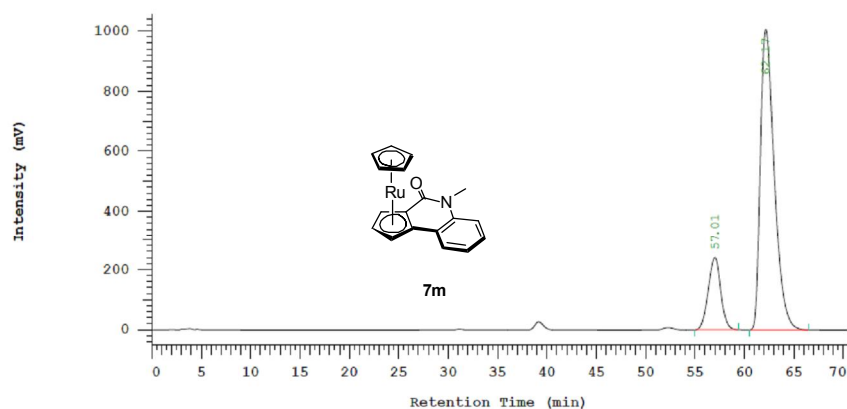
Processed Date and Time: 2014-06-02 21:52

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0122\

Processing Method: isopropanol

System (acquisition): LCE Series: 0122
 Application(data): maxinna Vial Number: 1
 Sample Name: mxn-2-165 Vial Type: UNK
 Injection from this vial: 1 of 1 Volume: 10.0 ul
 Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	57.01	20874481	17.654	BB
2	62.17	97364953	82.346	BB
		1.182E+08	100.000	

Peak rejection level: 0

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-05-21

Reported Date and Time: 2014-06-02

11:04

21:26

Processed Date and Time: 2014-06-02

21:25

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0135\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0135

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-133

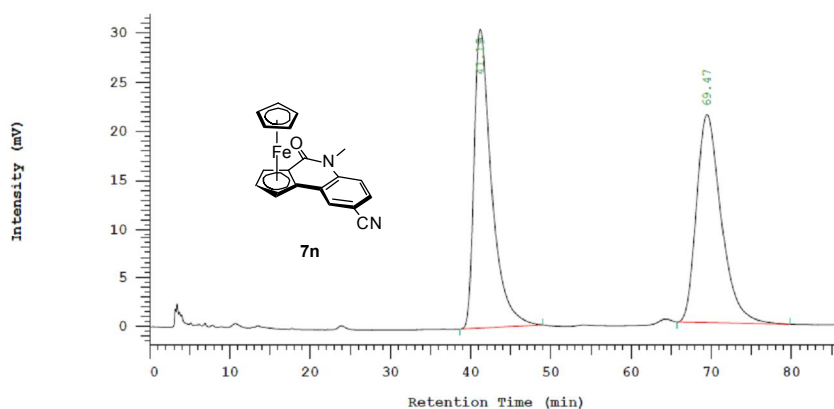
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	41.19	4619472	50.509	BB
2	69.47	4526402	49.491	BB
		9145874	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0136

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-05-21
12:32

Reported Date and Time: 2014-06-02
21:24

Processed Date and Time: 2014-06-02
21:24

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0136\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0136

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-133

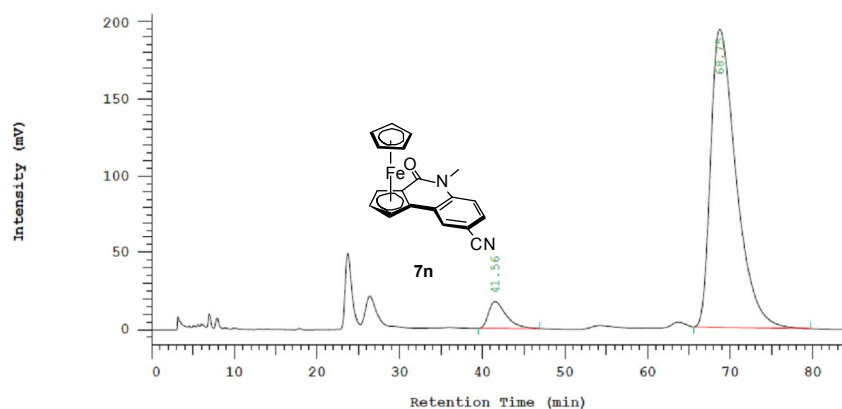
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	41.56	2526523	5.785	BB
2	68.75	41146339	94.215	BB
		43672862	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0050

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-03-26

Reported Date and Time: 2014-06-02

11:55

21:22

Processed Date and Time: 2014-06-02

21:21

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0050\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0050

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-131

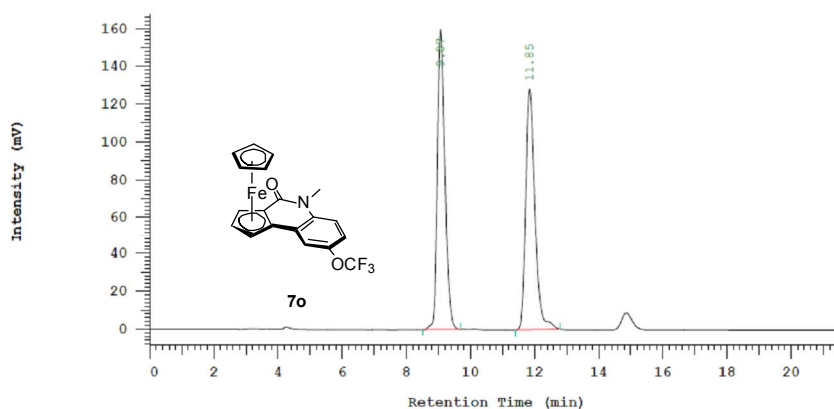
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	9.07	2595461	50.108	BB
2	11.85	2584226	49.892	BB
		5179687	100.000	

Peak rejection level: 0

D-2000: maxinna Series: 0051

Report Name: modified System: LCE

D-2000 Elite HPLC System Manager Report

Analyzed Date and Time: 2014-03-26

Reported Date and Time: 2014-06-02

12:22

21:20

Processed Date and Time: 2014-06-02

21:19

Data Path: C:\WIN32APP\D2000HSM\maxinna\DATA\0051\

Processing Method: isopropanol

System (acquisition): LCE

Series: 0051

Application(data): maxinna

Vial Number: 1

Sample Name: mxn-2-131

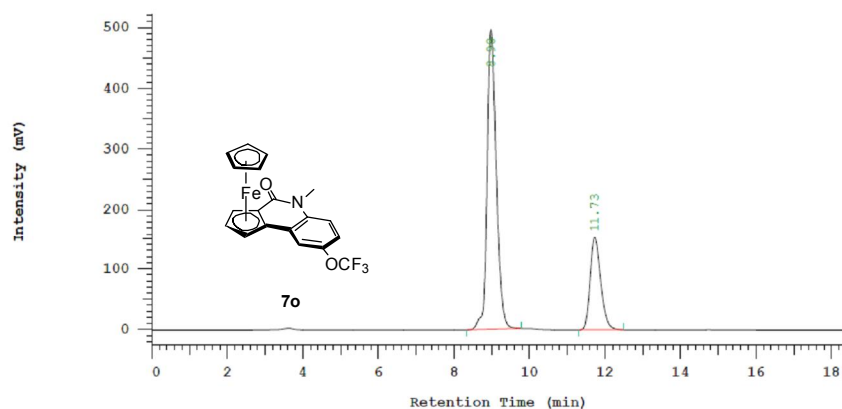
Vial Type: UNK

Injection from this vial: 1 of 1

Volume: 10.0 ul

Sample Description: xleiliu04-113-OJ-12

Chrom Type: HPLC Channel : 1



Processing Method: isopropanol

Column Type: Column

Method Developer:

Method Description:

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA

Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	8.98	8323963	73.771	BB
2	11.73	2959532	26.229	BB
		11283495	100.000	

Peak rejection level: 0