

Electronic Supplementary Information (ESI)

Chemoselective synthesis of aryl(pyridinyl)methanol derivatives through Ni-NIXANTPHOS catalyzed α -arylation and tandem arylation/rearrangement of pyridylmethyl ethers

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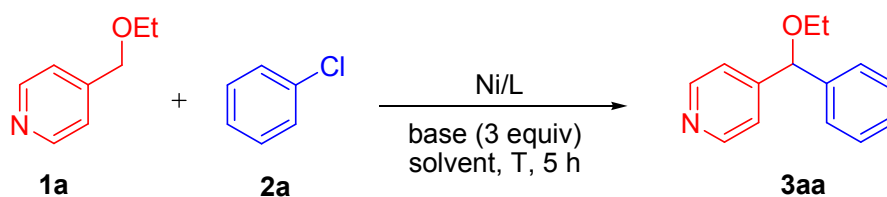
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1. General Information

All air- and moisture-sensitive solutions and chemicals were handled under nitrogen or in a nitrogen filled glovebox and solutions were transferred via syringe. Anhydrous CPME (cyclopentyl methyl ether), DME (Dimethoxyethane), toluene, and tetrahydrofuran (THF) were purchased from Sigma-Aldrich and used without further purification. Unless otherwise stated, all reagents were commercially available and used as received without further purification. Chemicals were obtained from Sigma-Aldrich, Acros, TCI and Alfa-Aesar. TLC was performed with Merck TLC Silica gel60 F₂₅₄ plates with detection under UV light at 254 nm. Silica gel (200-300 mesh, Qingdao) was used for flash chromatography. Proton nuclear magnetic resonance (¹H-NMR) spectra were recorded on Bruker DRX 400 spectrometer at 400 MHz. Carbon-13 nuclear magnetic resonance (¹³C-NMR) was recorded on Bruker DRX 400 spectrometer at 100 MHz. Chemical shifts are reported as δ values in parts per million (ppm) relative to tetramethylsilane (TMS) for all recorded NMR spectra. The infrared (IR) spectra were measured on a Nicolet iS10 FTIR spectrometer with 4 cm⁻¹ resolution and 32 scans between wavenumber of 4000 cm⁻¹ and 400 cm⁻¹. High Resolution Mass spectra were taken on AB QSTAR Pulsar mass spectrometer. Melting points were obtained on a XT-4 melting-point apparatus and were uncorrected.

2. Ni-catalyzed direct arylation of pyridylmethyl ethers: *Lab scale reaction optimization of ligands, Ni sources, solvents, temperature and ratio of Ni/ligand*

Table S1. Optimization of arylation of 4-pyridylmethyl ethyl ether **1a** with chlorobenzene **2a**



	Ni source	L	Ni/L (mol %)	Base	Solvent	Temp. (°C)	Assay Yield (%)
1		L1					70
2		L2					14
3		L3					0
4		L4					0
5		L5					0
6	Ni(COD) ₂	L6	10/15	NaN(SiMe ₃) ₂	DME	65	0
7		L7					0
8		L8					0
9		L9					0
10		L10					0
11		L11					0
13	Ni(acac) ₂						20
14	NiCl ₂ ·gly	L1	10/15	NaN(SiMe ₃) ₂	DME	65	19
15	Ni(OAc) ₂						41
16	NiBr ₂						23
17				LiN(SiMe ₃) ₂			76
18				KN(SiMe ₃) ₂		65	7
19	Ni(COD) ₂	L1	10/15	LiOtBu	DME		7
20				NaOtBu			0
21				KOtBu			0
22					CPME		37
23	Ni(COD) ₂	L1	10/15	LiN(SiMe ₃) ₂	Toluene	65	75
24					THF		8
25	Ni(COD) ₂	L1	10/15	LiN(SiMe ₃) ₂	DME	40	73
26						25	81
27			5/7.5				71
28	Ni(COD) ₂	L1	2.5/3.75	LiN(SiMe ₃) ₂	DME	25	60
29			10/10				76

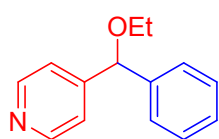
^a Reactions conducted on a 0.1 mmol scale using 1 equiv of **1a**, and 1.5 equiv of **2a**. ^b Assay yield determined by ¹HNMR spectroscopy of the crude reaction mixture.

Ligands: L1 = NIXANTPHOS, L2 = XantPhos, L3 = DavePhos, L4 = DPE-Phos, L5 = Pxy₃, L6 = DCE-Phos, L7 = SPhos, L8 = XPhos, L9 = RuPhos, L10 = DPPE, L11 = P(*o*-tol)₃

3. General Procedure and characterization of Ni-catalyzed arylation of pyridylmethyl ethers (Table 2 and 3)

General Procedure A for the Ni-catalyzed chemoselective C(sp³)-H arylation of pyridylmethyl ethers:

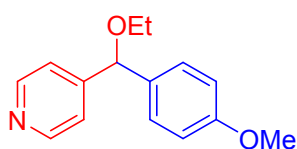
An oven-dried 8 mL reaction vial equipped with a stir bar was charged with pyridylmethyl ether **1a** (0.2 mmol, 1.0 equiv), chlorobenzene **2a** (0.3 mmol, 1.5 equiv) in a glove box under a nitrogen atmosphere at room temperature. A solution (from a stock solution) of Ni(COD)₂ (5.5 mg, 0.02 mmol, 10 mol %) and NIXANTPHOS (16.6 mg, 0.03 mmol, 15 mol %) in 1 mL of dry DME was taken up by syringe and added to the reaction vial. A solution of LiN(SiMe₃)₂ (100.4 mg, 0.6 mmol, 3.0 equiv) in 1 mL of dry DME was added by syringe to the reaction mixture. Total volume of the reaction is 2 mL. The vial was capped, removed from the glove box, and stirred for 5 h in total at 25 °C. The reaction mixture was quenched with 3 drops of H₂O, diluted with 2 mL of ethyl acetate, and filtered over a pad of MgSO₄ and silica. The pad was rinsed with an additional 6 mL of ethyl acetate (3 x 2 mL) and the solution was concentrated in vacuo. The crude material was loaded onto a silica gel column and purified by flash chromatography.



4-(Ethoxy(phenyl)methyl)pyridine (3aa): The reaction was performed following the

General Procedure A with 4-pyridylmethyl ethyl ether **1a** (27.4 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μ L, 0.3 mmol; **2a'**, X = Br, 31.5 μ L, 0.3 mmol). The

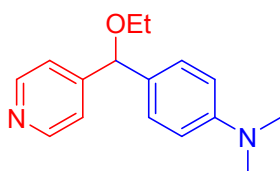
crude product was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:5) to give the product **3aa** (X = Cl, 33.7 mg, 79% yield; X = Br, 36.3 mg, 85% yield) as a yellow oil. R_f = 0.47 (EtOAc : hexanes = 1:1); ¹H NMR (400 MHz, Chloroform-*d*) δ 8.52–8.42 (m, 2H), 7.30–7.25 (m, 4H), 7.24–7.17 (m, 3H), 5.24 (s, 1H), 3.50–3.40 (m, 2H), 1.21 (t, J = 7.0 Hz, 3H) ppm; ¹³C{¹H} NMR (100 MHz, Chloroform-*d*) δ 151.5, 149.9, 141.1, 128.8, 128.2, 127.2, 121.7, 82.4, 64.9, 15.4 ppm; IR (thin film): 3061, 2974, 2870, 1668, 1597, 1561, 1494, 1412, 1308, 1279, 1261, 1188, 1101, 1027, 792, 700 cm⁻¹; HRMS calc'd for C₁₄H₁₆NO⁺ 214.1226, found 214.1226 [M+H]⁺.



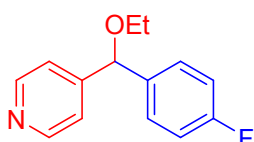
4-(Ethoxy(4-methoxyphenyl)methyl)pyridine (3ab): The reaction was

performed following the General Procedure A with 4-pyridylmethyl ethyl ether **1a** (27.4 mg, 0.20 mmol) and aryl halides **2b** or **2b'** (**2b**, X = Cl, 36.5 μ L, 0.3 mmol;

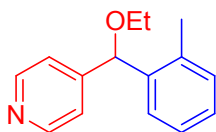
2b', X = Br, 37.5 μ L, 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:5) to give the product **3ab** (X = Cl, 39.4 mg, 81% yield; X = Br, 43.8 mg, 90% yield) as a yellow solid. The ¹H and ¹³C{¹H} NMR data for this compound match the literature data.¹



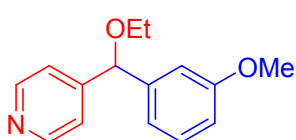
4-(Ethoxy(pyridin-4-yl)methyl)-N,N-dimethylaniline (3ac): The reaction was performed following the General Procedure A with 4-pyridylmethyl ethyl ether **1a** (27.4 mg, 0.20 mmol) and aryl halides **2c'** (X = Br, 60.0 mg, 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:5) to give the product **3ac** (X = Br, 49.2 mg, 96% yield) as a yellow oil. The ^1H and $^{13}\text{C}\{^1\text{H}\}$ NMR data for this compound match the literature data.¹



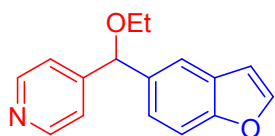
4-(Ethoxy(4-fluorophenyl)methyl)pyridine (3ad): The reaction was performed following the General Procedure A with 4-pyridylmethyl ethyl ether **1a** (27.4 mg, 0.20 mmol) and aryl halides **2d** or **2d'** (**2d**, X = Cl, 32.0 μL , 0.3 mmol; **2d'**, X = Br, 33.0 μL , 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:5) to give the product **3ad** (X = Cl, 36.1 mg, 78% yield; X = Br, 29.1 mg, 63% yield) as a yellow oil. The ^1H and $^{13}\text{C}\{^1\text{H}\}$ NMR data for this compound match the literature data¹. ^{19}F NMR (376 MHz, Chloroform-*d*) δ -114 ppm.



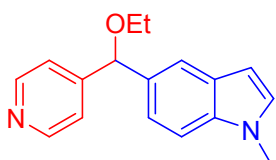
4-(Ethoxy(o-tolyl)methyl)pyridine (3ae): The reaction was performed following the General Procedure A with 4-pyridylmethyl ethyl ether **1a** (27.4 mg, 0.20 mmol) and aryl halides **2e** or **2e'** (**2e**, X = Cl, 35.0 μL , 0.3 mmol; **2e'**, X = Br, 36.0 μL , 0.3 mmol). The crude material was purified by Flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:5) to give the product **3ae** (X = Cl, 32.7 mg, 72% yield; X = Br, 43.2 mg, 95% yield) as a yellow oil. The ^1H and $^{13}\text{C}\{^1\text{H}\}$ NMR data for this compound match the literature data.¹



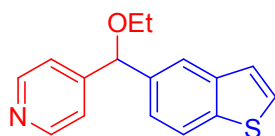
4-(Ethoxy(3-methoxyphenyl)methyl)pyridine (3af): The reaction was performed following the General Procedure A with 4-pyridylmethyl ethyl ether **1a** (27.4 mg, 0.20 mmol) and aryl halides **2f** (X = Br, 37.5 μL , 0.30 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:5) to give the product **3af** (X = Br, 34.6 mg, 71% yield) as a yellow oil. R_f = 0.43 (EtOAc : hexanes = 1:1); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.46 (d, J = 5.2 Hz, 2H), 7.24–7.15 (m, 3H), 6.88–6.78 (m, 2H), 6.75 (dd, J = 8.4, 1.2 Hz, 1H), 5.21 (s, 1H), 3.71 (s, 3H), 3.53–3.39 (m, 2H), 1.20 (t, J = 7.0 Hz, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 160.0, 151.4, 149.9, 142.7, 129.8, 121.7, 119.6, 113.3, 112.9, 82.2, 64.9, 55.4, 15.4 ppm; IR (thin film): 3028, 2974, 2929, 1597, 1488, 1454, 1412, 1314, 1260, 1150, 1103, 1082, 1049, 877, 786, 767, 700 cm^{-1} ; HRMS calc'd for $\text{C}_{15}\text{H}_{18}\text{NO}_2^+$ 244.1332, found 244.1333 $[\text{M}+\text{H}]^+$.



4-(Benzofuran-5-yl(ethoxy)methyl)pyridine (3ag): The reaction was performed following the General Procedure A with 4-pyridylmethyl ethyl ether **1a** (27.4 mg, 0.20 mmol) and aryl halides **2g'** (X = Br, 37.5 μ L, 0.30 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:6) to give the product **3ag** (X = Br, 36.0 mg, 71% yield) as a yellow oil. R_f = 0.52 (EtOAc : hexanes = 1:1); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.50–8.42 (m, 2H), 7.54 (d, J = 5.2 Hz, 1H), 7.49 (d, J = 4.4 Hz, 1H), 7.39 (d, J = 9.2 Hz, 1H), 7.28–7.20 (m, 2H), 7.19–7.15 (m, 1H), 6.66 (dd, J = 5.2, 1.2 Hz, 1H), 5.34 (s, 1H), 3.51–3.40 (m, 2H), 1.20 (t, J = 7.0 Hz, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 154.7, 151.9, 149.9, 145.8, 135.8, 127.7, 123.7, 121.7, 120.1, 111.7, 106.8, 82.4, 64.8, 15.4 ppm; IR (thin film): 3028, 2975, 2928, 2871, 1599, 1468, 1444, 1411, 1331, 1264, 1192, 1098, 1032, 886, 810, 770, 742 cm^{-1} ; HRMS calc'd for $\text{C}_{16}\text{H}_{16}\text{NO}_2^+$ 254.1176, found 254.1176 $[\text{M}+\text{H}]^+$.

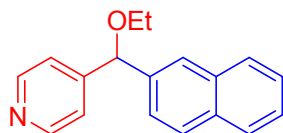


5-(Ethoxy(pyridin-4-yl)methyl)-1-methyl-1H-indole (3ah): The reaction was performed following the General Procedure A with 4-pyridylmethyl ethyl ether **1a** (27.4 mg, 0.20 mmol) and aryl halides **2h'** (X = Br, 63.0 mg, 0.30 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:5) to give the product **3ah** (X = Br, 44.8 mg, 84% yield) as a yellow oil. R_f = 0.47 (EtOAc : hexanes = 1:1); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.57–8.44 (m, 2H), 7.66–7.56 (m, 1H), 7.37–7.27 (m, 3H), 7.15 (dd, J = 8.0, 5.2 Hz, 1H), 7.06 (d, J = 1.6 Hz, 1H), 6.53–6.42 (m, 1H), 5.43 (s, 1H), 3.77 (s, 3H), 3.61–3.46 (m, 2H), 1.28 (t, J = 7.0 Hz, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 152.6, 149.8, 136.5, 132.0, 129.6, 128.4, 121.7, 121.0, 120.2, 109.7, 101.2, 82.9, 64.5, 33.0, 15.4 ppm; IR (thin film): 3025, 2972, 2926, 1578, 1513, 1493, 1447, 1411, 1344, 1307, 1245, 1099, 1023, 993, 892, 799, 723 cm^{-1} ; HRMS calc'd for $\text{C}_{17}\text{H}_{19}\text{N}_2\text{O}^+$ 267.1492, found 267.1492 $[\text{M}+\text{H}]^+$.

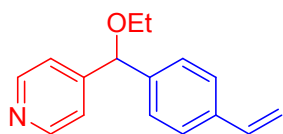


4-(Benzo[b]thiophen-5-yl(ethoxy)methyl)pyridine (3ai): The reaction was performed following the General Procedure A with 4-pyridylmethyl ethyl ether **1a** (27.4 mg, 0.20 mmol) and aryl halides **2i** or **2i'** (**2i**, X = Cl, 50.6 mg, 0.3 mmol; **2i'**, X = Br, 63.9 mg, 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:6) to give the product **3ai** (X = Cl, 39.9 mg, 74% yield; X = Br, 40.4 mg, 75% yield) as a yellow oil. R_f = 0.50 (EtOAc : hexanes = 1:1); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.53–8.38 (m, 2H), 7.75 (d, J = 8.0 Hz, 1H), 7.71 (d, J = 4.0 Hz, 1H), 7.37 (d, J = 8.4 Hz, 1H), 7.26–7.16 (m, 4H), 5.35 (s, 1H), 3.51–3.41 (m, 2H), 1.20 (t, J = 7.0 Hz, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 151.7, 149.9, 139.8, 139.5,

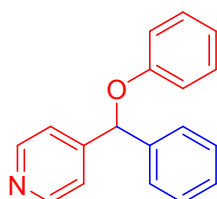
137.3, 127.3, 123.9, 123.5, 122.9, 122.4, 121.7, 82.4, 64.9, 15.4 ppm; IR (thin film): 3028, 2974, 2927, 2870, 1597, 1561, 1410, 1325, 1307, 1257, 1192, 1099, 1026, 993, 805, 777, 704 cm^{-1} ; HRMS calc'd for $\text{C}_{16}\text{H}_{16}\text{NOS}^+$ 270.0947, found 270.0947 $[\text{M}+\text{H}]^+$.



4-(Ethoxy(naphthalen-2-yl)methyl)pyridine (3aj): The reaction was performed following the General Procedure A with 4-pyridylmethyl ethyl ether **1a** (27.4 mg, 0.20 mmol) and aryl halides **2j** or **2j'** (**2j**, X = Cl, 48.8 mg, 0.3 mmol; **2j'**, X = Br, 62.1 mg, 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:6) to give the product **3aj** (X = Cl, 43.7 mg, 83% yield; X = Br, 32.7 mg, 62% yield) as a yellow oil. R_f = 0.52 (EtOAc : hexanes = 1:1); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.54 (d, J = 4.4 Hz, 2H), 7.87–7.78 (m, 4H), 7.49 (td, J = 8.4, 1.6 Hz, 2H), 7.39 (dd, J = 8.0, 1.6 Hz, 1H), 7.37–7.30 (m, 2H), 5.48 (s, 1H), 3.60–3.53 (m, 2H), 1.30 (t, J = 7.0 Hz, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 151.4, 150.0, 138.4, 133.3, 128.9, 128.1, 127.9, 126.5, 126.4, 124.9, 121.8, 82.5, 64.9, 15.4 ppm; IR (thin film): 3056, 2974, 2926, 2854, 1722, 1682, 1597, 1561, 1411, 1370, 1280, 1166, 1099, 992, 851, 747 cm^{-1} ; HRMS calc'd for $\text{C}_{18}\text{H}_{18}\text{NO}^+$ 264.1383, found 264.1382 $[\text{M}+\text{H}]^+$.

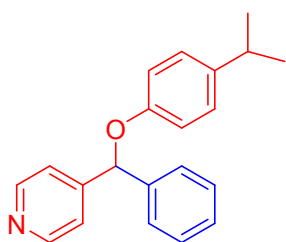


4-(Ethoxy(4-vinylphenyl)methyl)pyridine (3ak): The reaction was performed following the General Procedure A with 4-pyridylmethyl ethyl ether **1a** (27.4 mg, 0.20 mmol) and aryl halides **2k** or **2k'** (**2k**, X = Cl, 36.0 μL , 0.3 mmol; **2k'**, X = Br, 39.0 μL , 0.3 mmol). The crude material was purified by Flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:5) to give the product (X = Cl, 36.4 mg, 76% yield; X = Br, 39.3 mg, 82% yield) as a yellow oil. R_f = 0.48 (EtOAc : hexanes = 1:1); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.52–8.42 (m, 2H), 7.35–7.27 (m, 2H), 7.23–7.18 (m, 4H), 6.62 (dd, J = 20.4, 8.8 Hz, 1H), 5.67 (dd, J = 20.8, 1.2 Hz, 1H), 5.23 (s, 1H), 5.17 (dd, J = 12.4, 1.8 Hz, 1H), 3.48–3.41 (m, 2H), 1.20 (t, J = 7.0 Hz, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 150.0, 140.6, 136.4, 127.5, 126.6, 121.7, 114.4, 82.1, 64.9, 15.4 ppm; IR (thin film): 3025, 2965, 2926, 2854, 1629, 1598, 1509, 1411, 1306, 1261, 1098, 1017, 992, 909, 800, 764 cm^{-1} ; HRMS calc'd for $\text{C}_{16}\text{H}_{18}\text{NO}^+$ 240.1383, found 240.1380 $[\text{M}+\text{H}]^+$.

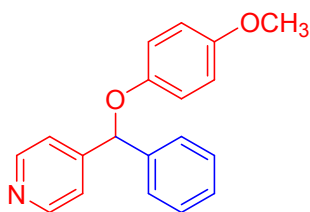


4-(Phenoxy(phenyl)methyl)pyridine (3ba): The reaction was performed following the General Procedure A with **1b** (37.1 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:5) to give the product **3ba** (X = Cl, 45.0 mg, 86% yield; X = Br, 43.4 mg, 83% yield) as a yellow oil. R_f = 0.37 (EtOAc :

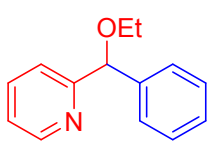
hexanes = 1:1); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.58 (d, J = 8.4 Hz, 2H), 7.50–7.28 (m, 7H), 7.26–7.20 (m, 2H), 7.00–6.85 (m, 3H), 6.16 (s, 1H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 157.5, 150.6, 149.7, 139.7, 129.5, 128.9, 128.5, 127.0, 121.6, 116.1, 105.0, 80.4 ppm; IR (thin film): 3030, 2958, 2920, 2851, 1597, 1493, 1454, 1412, 1232, 1167, 1092, 1092, 1038, 820, 792, 753 cm^{-1} ; HRMS calc'd for $\text{C}_{18}\text{H}_{16}\text{NO}^+$ 262.1226, found 262..1225 $[\text{M}+\text{H}]^+$.



4-((4-Isopropylphenoxy)(phenyl)methyl)pyridine (3ca): The reaction was performed following the General Procedure A with **1c** (45.5 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:6) to give the product **3ca** (X = Cl, 39.4 mg, 65% yield; X = Br, 44.3 mg, 73% yield) as a yellow oil. R_f = 0.52 (EtOAc : hexanes = 1:1); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.57–8.37 (m, 2H), 7.35–7.16 (m, 7H), 7.03–6.92 (m, 2H), 6.84–6.70 (m, 2H), 6.02 (s, 1H), 2.78–2.67 (m, 2H), 1.09 (d, J = 6.8 Hz, 6H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 155.8, 150.5, 150.1, 142.0, 140.1, 128.9, 128.4, 127.4, 127.1, 121.6, 115.9, 80.6, 33.3, 24.2 ppm; IR (thin film): 2958, 2919, 2850, 1632, 1578, 1540, 1509, 1466, 1261, 1231, 1166, 1094, 800, 739 cm^{-1} ; HRMS calc'd for $\text{C}_{21}\text{H}_{22}\text{NO}^+$ 304.1696, found 304.1699 $[\text{M}+\text{H}]^+$.



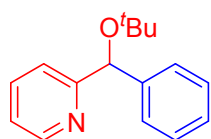
4-((4-Methoxyphenoxy)(phenyl)methyl)pyridine (3da): The reaction was performed following the General Procedure A with **1d** (43.1 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:5) to give the product **3da** (X = Cl, 41.4 mg, 71% yield; X = Br, 52.4 mg, 90% yield) as a yellow oil. R_f = 0.40 (EtOAc : hexanes = 1:1); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.64–8.49 (m, 2H), 7.42–7.27 (m, 7H), 6.90–6.82 (m, 2H), 6.80–6.73 (m, 2H), 6.05 (s, 1H), 3.73 (s, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 154.5, 151.8, 150.5, 150.1, 140.1, 129.0, 128.5, 127.2, 121.6, 117.5, 114.7, 81.6, 55.7 ppm; IR (thin film): 2920, 2851, 1596, 1506, 1454, 1412, 1261, 1224, 1181, 1106, 1038, 825, 797, 761, 738 cm^{-1} ; HRMS calc'd for $\text{C}_{19}\text{H}_{18}\text{NO}_2^+$ 292.1332, found 292.1334 $[\text{M}+\text{H}]^+$.



2-(Ethoxy(phenyl)methyl)pyridine (3ea): The reaction was performed following the General Procedure A with 4-pyridylmethyl ethyl ether **1e** (27.4 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The

crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:10) to give

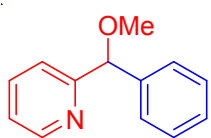
the product **3ea** (X = Cl, 30.7 mg, 72% yield; X = Br, 40.1 mg, 94% yield) as a yellow oil. The ^1H and $^{13}\text{C}\{^1\text{H}\}$ NMR data for this compound match the literature data.¹



2-(*tert*-Butoxy(phenyl)methyl)pyridine (3fa): The reaction was performed following the

General Procedure A with **1f** (33.1 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material was purified by

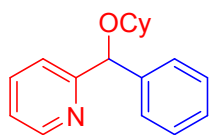
flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:10) to give the product **3fa** (X = Cl, 45.4 mg, 94% yield; X = Br, 41.0 mg, 85% yield) as a yellow oil. R_f = 0.38 (EtOAc : hexanes = 1:5); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.41 (ddd, J = 5.2, 1.6, 1.2 Hz, 1H), 7.63–7.47 (m, 2H), 7.42–7.32 (m, 2H), 7.23–7.18 (m, 2H), 7.14–7.08 (m, 1H), 7.03 (ddd, J = 7.2, 4.8, 1.6 Hz, 1H), 5.65 (s, 1H), 1.15 (s, 9H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 164.8, 148.6, 144.0, 136.7, 128.3, 127.0, 126.7, 122.0, 121.3, 77.3, 75.4, 28.8 ppm; IR (thin film): 2974, 2924, 2851, 1589, 1571, 1470, 1434, 1391, 1367, 1261, 1192, 1097, 1081, 1065, 1024, 905, 800, 744 cm^{-1} ; HRMS calc'd for $\text{C}_{16}\text{H}_{20}\text{NO}^+$ 242.1539, found 242.1539 $[\text{M}+\text{H}]^+$.



2-(Methoxy(phenyl)methyl)pyridine (3ga): The reaction was performed following the

General Procedure A with **1g** (24.6 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material was purified by

flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:8) to give the product **3ga** (X = Cl, 25.1 mg, 63% yield; X = Br, 27.1 mg, 68% yield) as a yellow oil. R_f = 0.20 (EtOAc : hexanes = 1:5); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.47 (ddd, J = 4.8, 1.6, 1.2 Hz, 1H), 7.61 (td, J = 8.4, 2.4 Hz, 1H), 7.42 (dt, J = 8.0, 1.6 Hz, 1H), 7.38–7.34 (m, 2H), 7.29–7.23 (m, 2H), 7.20–7.15 (m, 1H), 7.08 (ddd, J = 8.8, 5.2, 1.6 Hz, 1H), 5.31 (s, 1H), 3.36 (s, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 161.6, 149.2, 140.9, 137.0, 128.6, 127.9, 127.1, 122.5, 120.7, 86.6, 57.3 ppm; IR (thin film): 2959, 2919, 2850, 1588, 1572, 1541, 1470, 1453, 1436, 1261, 1194, 1166, 1102, 977, 800, 745 cm^{-1} ; HRMS calc'd for $\text{C}_{13}\text{H}_{14}\text{NO}^+$ 200.1070, found 200.1070 $[\text{M}+\text{H}]^+$.

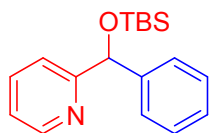


2-((Cyclohexyloxy)(phenyl)methyl)pyridine (3ha): The reaction was performed

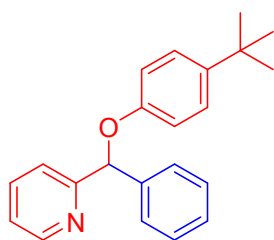
following the General Procedure A with **1h** (38.3 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material

was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:10) to give the product **3ha** (X = Cl, 52.4 mg, 98% yield; X = Br, 52.4 mg, 98% yield) as a colorless oil. R_f = 0.52 (EtOAc : hexanes = 1:5); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.38 (d, J = 4.8 Hz, 1H), 7.53–7.46 (m, 2H), 7.34 (d, J = 8.0 Hz, 2H), 7.18 (t, J = 8.8 Hz, 2H), 7.09 (t, J = 9.2 Hz, 1H), 6.97 (t, J = 6.0 Hz, 1H), 5.58 (s, 1H), 3.30 (tt, J = 9.2, 5.2 Hz, 1H), 1.87–1.79 (m, 2H), 1.65–1.59 (m, 2H), 1.41–1.26 (m, 3H), 1.14–1.05 (m, 2H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz,

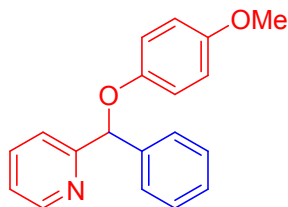
Chloroform-*d*) δ 162.91, 148.68, 142.13, 136.67, 128.26, 127.30, 126.87, 122.10, 120.75, 81.40, 75.51, 32.40, 32.38, 25.80, 24.02 ppm; IR (thin film): 3061, 3029, 2931, 2856, 1588, 1570, 1469, 1450, 1433, 1356, 1259, 1189, 1132, 1099, 1070, 1028, 958, 744, 699, 637 cm^{-1} ; HRMS calc'd for $\text{C}_{18}\text{H}_{21}\text{NONa}^+$ 290.1515, found 290.1504 $[\text{M}+\text{Na}]^+$.



2-(((*tert*-Butyldimethylsilyloxy)(phenyl)methyl)pyridine (3ia): The reaction was performed following the General Procedure A with **1i** (44.7 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:10) to give the product **3ia** (X = Cl, 53.3 mg, 89% yield; X = Br, 57.5 mg, 96% yield) as a colorless oil. The ^1H and $^{13}\text{C}\{^1\text{H}\}$ NMR data for this compound match the literature data.³

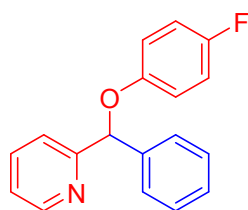


2-((4-(*tert*-Butyl)phenoxy)(phenyl)methyl)pyridine (3ja): The reaction was performed following the General Procedure A with **1j** (48.3 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:10) to give the product **3ja** (X = Cl, 58.4 mg, 92% yield; X = Br, 59.7 mg, 94% yield) as a white solid. m.p. = 88 – 89 $^{\circ}\text{C}$. R_f = 0.38 (EtOAc : hexanes = 1:5); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.58–8.42 (m, 1H), 7.58 (td, J = 8.4, 2.4 Hz, 1H), 7.51 (dt, J = 8.8, 1.2 Hz, 1H), 7.48–7.42 (m, 2H), 7.29–7.23 (m, 2H), 7.20–7.14 (m, 4H), 7.08 (ddd, J = 9.2, 4.8, 1.6 Hz, 1H), 6.86–6.79 (m, 2H), 6.23 (s, 1H), 1.18 (s, 9H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 161.2, 155.7, 149.3, 143.9, 140.6, 137.2, 128.7, 128.0, 126.9, 126.3, 122.7, 120.9, 115.4, 82.7, 34.2, 31.6 ppm; IR (thin film): 2962, 2921, 2852, 1607, 1589, 1511, 1470, 1434, 1364, 1236, 1183, 1095, 1027, 827, 804, 744 cm^{-1} ; HRMS calc'd for $\text{C}_{22}\text{H}_{24}\text{NO}^+$ 318.1852, found 318.1852 $[\text{M}+\text{H}]^+$.



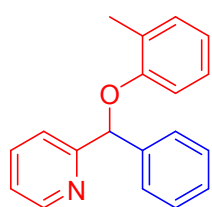
2-((4-Methoxyphenoxy)(phenyl)methyl)pyridine (3ka): The reaction was performed following the General Procedure A with **1k** (43.1 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:8) to give the product **3ka** (X = Cl, 40.8 mg, 70% yield; X = Br, 49.5 mg, 85% yield) as a yellow oil. R_f = 0.20 (EtOAc : hexanes = 1:5); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.49 (ddd, J = 5.2, 2.0, 1.2 Hz, 1H), 7.58 (td, J = 7.2, 1.6 Hz, 1H), 7.50 (dt, J = 8.0, 1.2 Hz, 1H), 7.47–7.40 (m, 2H), 7.32–7.22 (m, 2H), 7.20–7.16 (m, 1H), 7.08 (ddd, J = 7.2, 4.8, 1.2 Hz, 1H), 6.87–6.78 (m, 2H), 6.73–6.62 (m, 2H), 6.18 (s,

1H), 3.64 (s, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 161.1, 154.2, 151.9, 149.3, 140.5, 137.2, 128.7, 128.0, 126.9, 122.7, 120.9, 117.1, 114.7, 83.4, 55.7 ppm; IR (thin film): 2957, 2919, 2850, 1588, 1541, 1506, 1467, 1434, 1224, 1166, 1096, 1040, 823, 800, 742 cm^{-1} ; HRMS calc'd for $\text{C}_{19}\text{H}_{18}\text{NO}_2^+$ 292.1332, found 292.1331 $[\text{M}+\text{H}]^+$.



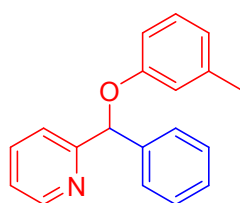
2-((4-Fluorophenoxy)(phenyl)methyl)pyridine (3la): The reaction was performed following the General Procedure A with **1l** (40.6 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material was purified by Flash chromatography on silica gel (eluted with EtOAc :

hexanes = 1:10) to give the product **3la** (X = Cl, 50.8 mg, 91% yield; X = Br, 49.7 mg, 89% yield) as a yellow oil. R_f = 0.30 (EtOAc : hexanes = 1:5); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.59–8.48 (m, 1H), 7.64 (td, J = 7.2, 1.6 Hz, 1H), 7.55–7.47 (m, 3H), 7.32 (t, J = 7.2 Hz, 2H), 7.23 (d, J = 6.0 Hz, 2H), 7.15 (ddd, J = 7.2, 4.8, 1.2 Hz, 1H), 6.88 (d, J = 7.2 Hz, 3H), 6.24 (s, 1H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (101 MHz, Chloroform-*d*) δ 160.7, 157.6 (d, $^1J_{\text{C-F}}$ = 238.0 Hz), 153.92 (d, $^4J_{\text{C-F}}$ = 2.0 Hz), 149.4, 140.1, 137.2, 128.8, 128.1 (d, $^3J_{\text{C-F}}$ = 8.0 Hz), 126.9, 122.8, 120.9, 117.2, 117.1, 116.1 (d, $^2J_{\text{C-F}}$ = 23.0 Hz), 83.4 ppm; IR (thin film): 2959, 2920, 2851, 1588, 1572, 1542, 1504, 1467, 1435, 1202, 1166, 1096, 826, 797, 781, 742 cm^{-1} ; HRMS calc'd for $\text{C}_{18}\text{H}_{15}\text{FNO}^+$ 280.1132, found 280.1137 $[\text{M}+\text{H}]^+$. ^{19}F NMR (376 MHz, Chloroform-*d*) δ -123 ppm.



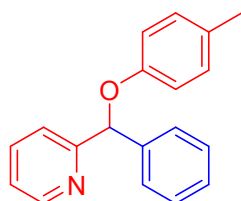
2-(Phenyl(o-tolyl)oxy)methylpyridine (3ma): The reaction was performed following the General Procedure A with **1m** (39.9 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material was purified

by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:10) to give the product **3ma** (X = Cl, 44.6 mg, 81% yield; X = Br, 45.2 mg, 82% yield) as a yellow oil. R_f = 0.37 (EtOAc : hexanes = 1:5); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.48 (ddd, J = 5.2, 2.4, 1.6 Hz, 1H), 7.57 (td, J = 7.2, 1.6 Hz, 1H), 7.53–7.42 (m, 3H), 7.30–7.21 (m, 2H), 7.20–7.14 (m, 1H), 7.07 (ddd, J = 7.2, 4.8, 1.2 Hz, 2H), 6.92 (ddd, J = 8.8, 7.2, 1.6 Hz, 1H), 6.80–6.62 (m, 2H), 6.29 (s, 1H), 2.33 (s, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 161.3, 155.8, 149.1, 140.7, 137.3, 130.9, 128.7, 127.9, 127.3, 126.8, 126.6, 122.7, 120.9, 120.6, 113.1, 82.2, 16.9 ppm; IR (thin film): 3063, 2960, 2919, 2851, 1588, 1541, 1493, 1465, 1434, 1235, 1185, 1165, 1120, 1095, 1049, 1027, 883, 800, 745 cm^{-1} ; HRMS calc'd for $\text{C}_{19}\text{H}_{18}\text{NO}^+$ 276.1383, found 276.1382 $[\text{M}+\text{H}]^+$.



2-(Phenyl(*m*-tolyl)oxy)methylpyridine (3na): The reaction was performed following the General Procedure A with **1n** (39.9 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material was

purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:10) to give the product **3na** (X = Cl, 50.1 mg, 91% yield; X = Br, 45.2 mg, 82% yield) as a yellow oil. R_f = 0.37 (EtOAc : hexanes = 1:5); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.47 (ddd, J = 4.8, 1.6, 1.2 Hz, 1H), 7.55 (td, J = 7.2, 1.6 Hz, 1H), 7.50–7.40 (m, 3H), 7.29–7.21 (m, 2H), 7.20–7.13 (m, 1H), 7.05 (ddd, J = 7.2, 4.8, 1.2 Hz, 1H), 7.00 (t, J = 8.0 Hz, 1H), 6.75 (t, J = 2.0 Hz, 1H), 6.71–6.60 (m, 2H), 6.26 (s, 1H), 2.18 (s, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 161.0, 157.8, 149.2, 140.4, 139.6, 137.2, 129.2, 128.7, 128.0, 126.9, 122.7, 122.1, 120.9, 117.0, 112.8, 82.4, 21.6 ppm; IR (thin film): 2957, 2919, 2850, 1579, 1540, 1489, 1467, 1434, 1289, 1257, 1166, 1095, 1047, 770, 742 cm^{-1} ; HRMS calc'd for $\text{C}_{19}\text{H}_{18}\text{NO}^+$ 276.1383, found 276.1382 $[\text{M}+\text{H}]^+$.

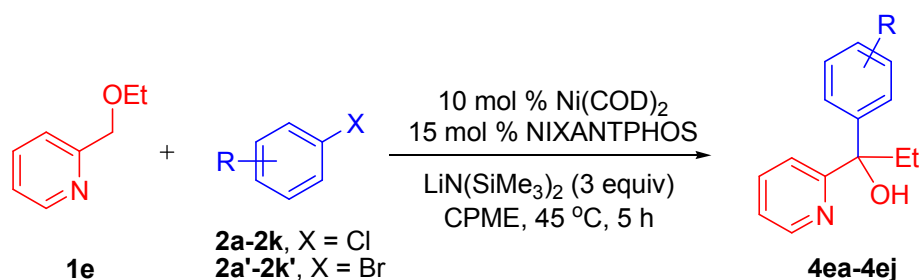


2-(Phenyl(*p*-tolylloxy)methyl)pyridine (3oa**):** The reaction was performed following the General Procedure A with **1o** (39.9 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μL , 0.3 mmol; **2a'**, X = Br, 31.5 μL , 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:10) to

give the product **3oa** (X = Cl, 45.7 mg, 83% yield; X = Br, 47.9 mg, 87% yield) as a yellow oil. R_f = 0.33 (EtOAc : hexanes = 1:5); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.49 (ddd, J = 4.8, 1.6, 1.2 Hz, 1H), 7.57 (td, J = 7.2, 1.6 Hz, 1H), 7.49 (dt, J = 8.0, 1.2 Hz, 1H), 7.47–7.42 (m, 2H), 7.29–7.23 (m, 2H), 7.20–7.16 (m, 1H), 7.07 (ddd, J = 7.2, 4.8, 1.2 Hz, 1H), 6.97–6.91 (m, 2H), 6.82–6.76 (m, 2H), 6.23 (s, 1H), 2.16 (s, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 161.1, 155.7, 149.3, 140.5, 137.2, 130.5, 130.0, 128.7, 128.0, 126.9, 122.7, 120.9, 115.9, 82.7, 20.6 ppm; IR (thin film): 3030, 2960, 2920, 2851, 1612, 1586, 1509, 1470, 1452, 1434, 1228, 1174, 1095, 1027, 922, 807, 742 cm^{-1} ; HRMS calc'd for $\text{C}_{19}\text{H}_{18}\text{NO}^+$ 276.1383, found 276.1383 $[\text{M}+\text{H}]^+$.

4. Tandem arylation/[1,2]-Wittig rearrangement of 2-pyridylmethyl ethyl ether: *Lab scale reaction optimization of Ni source, solvents and temperature*

Table S2. Optimization of tandem arylation/[1,2]-Wittig rearrangement of 2-pyridylmethyl ethyl ether **1e** with aryl halides.

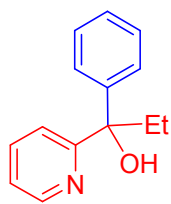


	Ni source	Base	Solvent	Temp. (°C)	Assay Yield (%)
1	Ni(COD) ₂				56
2	Ni(acac) ₂				4
3	NiCl ₂ •gly	NaN(SiMe ₃) ₂	CPME	65	3
4	Ni(OAc) ₂				10
5	NiBr ₂				2
6		LiN(SiMe ₃) ₂			71
7		KN(SiMe ₃) ₂			23
8	Ni(COD) ₂	LiO ^t Bu	CPME	65	0
9		NaO ^t Bu			0
10		KO ^t Bu			0
11			DME		21
12	Ni(COD) ₂	LiN(SiMe ₃) ₂	Toluene	65	60
13			THF		65
14	Ni(COD) ₂	LiN(SiMe ₃) ₂	CPME	45	72
15				25	68

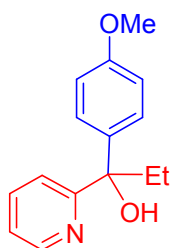
^a Reactions conducted on a 0.1 mmol scale using 1 equiv of **1e**, and 1.5 equiv of **2a**. ^b Assay yield determined by ¹HNMR spectroscopy of the crude reaction mixture.

5. Procedure and characterization of Ni-catalyzed chemoselective tandem arylation/[1,2]-Wittig rearrangement reaction of 2-pyridylmethyl ethers (Table 5).

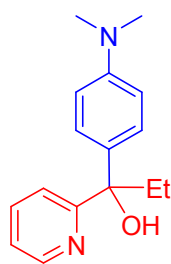
General Procedure B: An oven-dried 8 mL reaction vial equipped with a stir bar was charged with 2-pyridylmethyl ethyl ether **1e** (0.20 mmol, 1.0 equiv) and chlorobenzene **2a** (0.3 mmol, 1.5 equiv) in a glove box under a nitrogen atmosphere at room temperature. A solution (from a stock solution) of Ni(COD)₂ (5.5 mg, 0.02 mmol, 10 mol %) and NIXANTPHOS (16.6 mg, 0.03 mmol, 15 mol %) in 1 mL of dry CPME was taken up by syringe and added to the reaction vial. A solution of LiN(SiMe₃)₂ (100.4 mg, 0.6 mmol, 3.0 equiv) in 1 mL of dry CPME was added by syringe to the reaction mixture. Total volume of the reaction is 2 mL. The vial was capped, removed from the glove box, and stirred for 5 h in total at 45 °C. The reaction mixture was quenched with 3 drops of H₂O, diluted with 2 mL of ethyl acetate, and filtered over a pad of MgSO₄ and silica. The pad was rinsed with an additional 6 mL of ethyl acetate (3 x 2 mL) and the solution was concentrated in vacuo. The crude material was loaded onto a deactivated silica gel column and purified by flash chromatography.



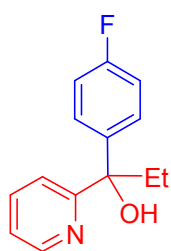
1-Phenyl-1-(pyridin-2-yl)propan-1-ol (4ea): The reaction was performed following the General Procedure B with **1e** (27.4 mg, 0.20 mmol) and aryl halides **2a** or **2a'** (**2a**, X = Cl, 30.5 μ L, 0.3 mmol; **2a'**, X = Br, 31.5 μ L, 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:8) to give the product **4ea** (X = Cl, 29.9 mg, 70% yield; X = Br, 26.9 mg, 63% yield) as a white solid. The ^1H and $^{13}\text{C}\{^1\text{H}\}$ NMR data for this compound match the literature data.²



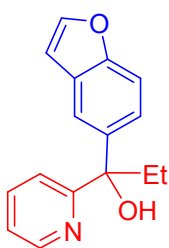
1-(4-Methoxyphenyl)-1-(pyridin-2-yl)propan-1-ol (4eb): The reaction was performed following the General Procedure B with **1e** (27.4 mg, 0.20 mmol) and aryl halides **2b** or **2b'** (**2b**, X = Cl, 36.5 μ L, 0.3 mmol; **2b'**, X = Br, 37.5 μ L, 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:8) to give the product **4eb** (X = Cl, 33.6 mg, 69% yield; X = Br, 27.3 mg, 56% yield) as a yellow solid. The ^1H and $^{13}\text{C}\{^1\text{H}\}$ NMR data for this compound match the literature data.¹



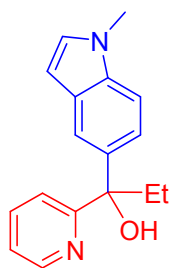
1-(4-(Dimethylamino)phenyl)-1-(pyridin-2-yl)propan-1-ol (4ec): The reaction was performed following the General Procedure B with **1e** (27.4 mg, 0.20 mmol) and aryl halides **2c** (X = Br, 60.0 mg, 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:4) to give the product **4ec** (X = Br, 28.2 mg, 55% yield) as a yellow solid. The ^1H and $^{13}\text{C}\{^1\text{H}\}$ NMR data for this compound match the literature data.¹



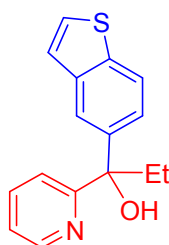
1-(4-Fluorophenyl)-1-(pyridin-2-yl)propan-1-ol (4ed): The reaction was performed following the General Procedure B with **1e** (27.4 mg, 0.20 mmol) and aryl halides **2d** or **2d'** (**2d**, X = Cl, 32.0 μ L, 0.3 mmol; **2d'**, X = Br, 33.0 μ L, 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:8) to give the product **4ed** (X = Cl, 31.0 mg, 67% yield; X = Br, 26.4 mg, 57% yield) as a yellow oil. The ^1H and $^{13}\text{C}\{^1\text{H}\}$ NMR data for this compound match the literature data¹. ^{19}F NMR (376 MHz, Chloroform-*d*) δ -116 ppm.



1-(Benzofuran-5-yl)-1-(pyridin-2-yl)propan-1-ol (4eg): The reaction was performed following the General Procedure B with **1e** (27.4 mg, 0.20 mmol) and aryl halides **2g** (X = Br, 37.5 μ L, 0.30 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:6) to give the product **4eg** (X = Br, 30.4 mg, 60% yield) as a white solid. The ^1H and $^{13}\text{C}\{^1\text{H}\}$ NMR data for this compound match the literature data.¹

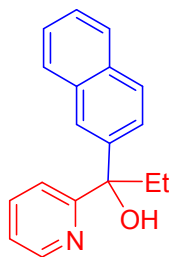


1-(1-Methyl-1H-indol-5-yl)-1-(pyridin-2-yl)propan-1-ol (4eh): The reaction was performed following the General Procedure B with **1e** (27.4 mg, 0.20 mmol) and aryl halides **2h** (X = Br, 63.0 mg, 0.30 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:6) to give the product **4eh** (X = Br, 22.9 mg, 43% yield) as a white solid. The ^1H and $^{13}\text{C}\{^1\text{H}\}$ NMR data for this compound match the literature data.¹



1-(Benzo[b]thiophen-5-yl)-1-(pyridin-2-yl)propan-1-ol (4ei): The reaction was performed following the General Procedure B with **1e** (27.4 mg, 0.20 mmol) and aryl halides **2i** or **2i'** (**2i**, X = Cl, 50.6 mg, 0.3 mmol; **2i'**, X = Br, 63.9 mg, 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:8) to give the product **4ei** (X = Cl, 37.7 mg, 70% yield; X = Br, 36.6 mg, 68% yield) as a yellow oil. R_f = 0.33

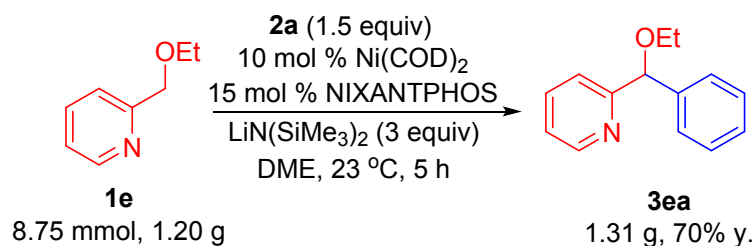
(EtOAc : hexanes = 1:5); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.56–8.48 (m, 1H), 8.04 (d, J = 1.6 Hz, 1H), 7.80 (d, J = 8.4 Hz, 1H), 7.64 (td, J = 7.2, 1.6 Hz, 1H), 7.49 (dd, J = 8.4, 1.6 Hz, 1H), 7.41 (d, J = 5.2 Hz, 1H), 7.36 (d, J = 8.0 Hz, 1H), 7.31 (d, J = 5.2 Hz, 1H), 7.17 (ddd, J = 7.2, 4.8, 1.2 Hz, 1H), 6.02 (s, 1H), 2.49–2.29 (m, 2H), 0.90 (t, J = 7.2 Hz, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 163.8, 147.3, 142.9, 139.8, 138.4, 137.2, 126.7, 124.3, 123.1, 122.3, 122.1, 121.0, 120.7, 77.6, 34.1, 8.2 ppm; IR (thin film): 3072, 2965, 2925, 2876, 1591, 1570, 1467, 1433, 1393, 1261, 1215, 1152, 1089, 1066, 1048, 988, 813, 776, 704 cm^{-1} ; HRMS calc'd for $\text{C}_{16}\text{H}_{16}\text{NOS}^+$ 270.0947, found 270.0949 $[\text{M}+\text{H}]^+$.



1-(Naphthalen-2-yl)-1-(pyridin-2-yl)propan-1-ol (4ej): The reaction was performed following the General Procedure B with **1e** (27.4 mg, 0.20 mmol) and aryl halides **2j** or **2j'** (**2j**, X = Cl, 48.8 mg, 0.3 mmol; **2j'**, X = Br, 62.1 mg, 0.3 mmol). The crude material was purified by flash chromatography on silica gel (eluted with EtOAc : hexanes = 1:8) to give the product **4ej** (X = Cl, 38.5 mg, 73% yield; X = Br, 35.3 mg, 67% yield) as a white oil. R_f = 0.35

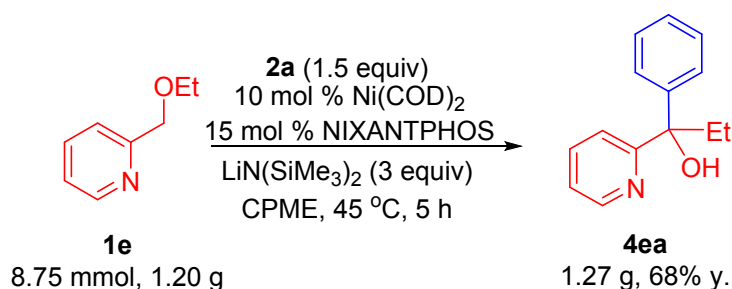
(EtOAc : hexanes = 1:5); ^1H NMR (400 MHz, Chloroform-*d*) δ 8.54 (d, J = 4.8 Hz, 1H), 8.05 (s, 1H), 7.88–7.81 (m, 1H), 7.77 (dd, J = 8.4, 4.8 Hz, 2H), 7.65 (t, J = 7.2 Hz, 1H), 7.58 (dd, J = 8.8, 1.6 Hz, 1H), 7.45 (ddd, J = 6.4, 4.0, 1.6 Hz, 2H), 7.37 (d, J = 8.0 Hz, 1H), 7.18 (dd, J = 7.2, 4.8 Hz, 1H), 6.05 (s, 1H), 2.53–2.33 (m, 2H), 0.92 (t, J = 7.2 Hz, 3H) ppm; $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, Chloroform-*d*) δ 163.6, 147.3, 143.8, 137.3, 133.3, 132.5, 128.4, 128.1, 127.6, 126.1, 126.0, 124.9, 124.6, 122.2, 120.9, 77.7, 33.9, 8.2 ppm; IR (thin film): 2959, 2920, 2851, 1661, 1590, 1541, 1466, 1431, 1374, 1166, 1127, 988, 818, 778, 746 cm^{-1} ; HRMS calc'd for $\text{C}_{18}\text{H}_{18}\text{NO}^+$ 264.1383, found 264.1383 $[\text{M}+\text{H}]^+$.

6. Gram scale synthesis of 3ea



An oven-dried 250 mL round bottom flask equipped with a stir bar was charged with 2-pyridylmethyl ethyl ether **1e** (1.20 g, 8.75 mmol, 1.0 equiv), Ni(COD)_2 (240.7 mg, 0.875 mmol, 10 mol %), NIXANTPHOS (724 mg, 1.31 mmol, 15 mol %) and chlorobenzene **2a** (1.33 mL, 13.12 mmol, 1.5 equiv) in a glove box under a nitrogen atmosphere at room temperature. 44 mL of dry DME was taken up by syringe and added to the reaction vial under nitrogen. Then, a solution of $\text{LiN(SiMe}_3)_2$ (4.39 g, 26.24 mmol, 3.0 equiv) in 44 mL of dry DME was added to the reaction mixture. The flask was capped with a rubber septum, removed from the glove box, and stirred for 5 h in total at 25 $^\circ\text{C}$, opened to air, and quenched with 10 mL of H_2O . The layers were separated and the aqueous layer was extracted with DCM (3X5 mL). The combined organic layers were concentrated *in vacuo*. The crude material was loaded onto a silica gel column and purified by flash chromatography using 10:1 hexanes/ethyl acetate as eluent to afford desired product **3ea** (1.31 g, 70% yield) as a yellow oil. The ^1H and $^{13}\text{C}\{^1\text{H}\}$ NMR data for this compound match the literature data.¹

7. Gram scale synthesis of 4ea



An oven-dried 250 mL round bottom flask equipped with a stir bar was charged with 2-pyridylmethyl ethyl ether **1e** (1.20 g, 8.75 mmol, 1.0 equiv), Ni(COD)_2 (240.7 mg, 0.875 mmol, 10 mol %), NIXANTPHOS (724 mg, 1.31 mmol, 15 mol %) and chlorobenzene **2a** (1.33 mL, 13.12 mmol, 1.5 equiv) in a glove box under a nitrogen atmosphere at room temperature. 44 mL of dry CPME was taken up by syringe and added to the reaction vial under nitrogen. Then, a solution of $\text{LiN(SiMe}_3)_2$ (4.39 g, 26.24 mmol, 3.0 equiv) in 44 mL of dry CPME was added to the reaction mixture. The flask was capped with a rubber septum, removed from the glove box, and

stirred for 5 h in total at 45 °C, opened to air, and quenched with 10 mL of H₂O. The layers were separated and the aqueous layer was extracted with DCM (3X5 mL). The combined organic layers were concentrated in *vacuo*. The crude material was loaded onto a deactivated silica gel column and purified by flash chromatography using 8 :1 hexanes/ethyl acetate as eluent to afford desired product **4ea** (1.27 g, 48% yield) as a white solid. The ¹H and ¹³C{¹H} NMR data for this compound match the literature data.²

8. X-ray crystal structures of compound 4ea

CCDC 1569085 contains the supplementary crystallographic data for compound **4ea**. The data can be obtained free of charge from The Cambridge Crystallographic Data Center via www.ccdc.cam.ac.uk/data_request/cif.

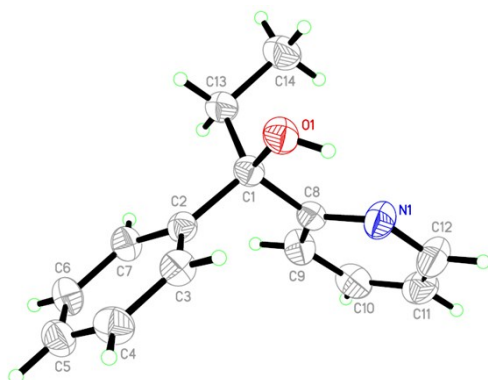


Figure 1 X-ray crystal structures of compound **4ea**

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C2 0.0473(10) 0.0416(10) 0.0487(10) -0.0032(8) 0.0082(8) -0.0013(8)
C3 0.0597(11) 0.0539(12) 0.0569(11) 0.0061(9) 0.0121(9) 0.0028(9)
C4 0.0738(14) 0.0608(13) 0.0680(12) 0.0077(10) 0.0281(11) -0.0052(11)
C5 0.0545(12) 0.0672(13) 0.0808(14) -0.0076(11) 0.0235(10) -0.0094(10)
C6 0.0466(11) 0.0653(13) 0.0768(13) 0.0064(11) 0.0093(9) 0.0037(9)
C7 0.0498(10) 0.0565(12) 0.0594(11) 0.0089(9) 0.0100(8) -0.0004(9)
C8 0.0454(10) 0.0502(11) 0.0498(10) 0.0015(8) 0.0152(8) 0.0005(8)
C9 0.0586(11) 0.0482(12) 0.0828(14) 0.0035(11) 0.0111(10) 0.0012(9)
C10 0.0737(14) 0.0495(13) 0.1045(17) -0.0064(12) 0.0330(13) -0.0055(11)
C11 0.0828(16) 0.0743(17) 0.0771(15) -0.0228(13) 0.0230(12) -0.0216(13)
C12 0.0835(15) 0.0845(18) 0.0698(14) -0.0120(13) -0.0083(12) -0.0047(13)
C13 0.0583(11) 0.0694(14) 0.0562(11) -0.0023(10) 0.0164(9) 0.0027(9)
C14 0.0732(14) 0.1080(19) 0.0765(14) 0.0002(13) 0.0324(11) 0.0041(13)
```

_geom_special_details

```
;
All esds (except the esd in the dihedral angle between two l.s. planes)
are estimated using the full covariance matrix. The cell esds are taken
into account individually in the estimation of esds in distances, angles
and torsion angles; correlations between esds in cell parameters are only
used when they are defined by crystal symmetry. An approximate (isotropic)
treatment of cell esds is used for estimating esds involving l.s. planes.
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;

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O1 H1 0.8200 . ?
N1 C8 1.331(2) . ?
N1 C12 1.337(3) . ?
C1 C2 1.533(2) . ?
C1 C8 1.537(2) . ?
C1 C13 1.538(2) . ?
C2 C3 1.387(2) . ?
C2 C7 1.391(2) . ?
C3 C4 1.391(3) . ?
C3 H3 0.9300 . ?
C4 C5 1.369(3) . ?
C4 H4 0.9300 . ?
C5 C6 1.372(3) . ?
C5 H5 0.9300 . ?
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C6 C7 1.383(2) . ?
 C6 H6 0.9300 . ?
 C7 H7 0.9300 . ?
 C8 C9 1.380(2) . ?
 C9 C10 1.375(3) . ?
 C9 H9 0.9300 . ?
 C10 C11 1.364(3) . ?
 C10 H10 0.9300 . ?
 C11 C12 1.363(3) . ?
 C11 H11 0.9300 . ?
 C12 H12 0.9300 . ?
 C13 C14 1.527(3) . ?
 C13 H13A 0.9700 . ?
 C13 H13B 0.9700 . ?
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 C14 H14B 0.9600 . ?
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 C8 N1 C12 117.55(17) . . ?
 O1 C1 C2 108.72(13) . . ?
 O1 C1 C8 109.02(13) . . ?
 C2 C1 C8 108.78(13) . . ?
 O1 C1 C13 107.86(14) . . ?
 C2 C1 C13 111.94(14) . . ?
 C8 C1 C13 110.46(14) . . ?
 C3 C2 C7 117.40(16) . . ?
 C3 C2 C1 120.07(15) . . ?
 C7 C2 C1 122.51(15) . . ?
 C2 C3 C4 120.94(17) . . ?
 C2 C3 H3 119.5 . . ?
 C4 C3 H3 119.5 . . ?
 C5 C4 C3 120.39(18) . . ?
 C5 C4 H4 119.8 . . ?
 C3 C4 H4 119.8 . . ?
 C4 C5 C6 119.73(18) . . ?
 C4 C5 H5 120.1 . . ?
 C6 C5 H5 120.1 . . ?
 C5 C6 C7 120.03(18) . . ?
 C5 C6 H6 120.0 . . ?
 C7 C6 H6 120.0 . . ?
 C6 C7 C2 121.52(18) . . ?
 C6 C7 H7 119.2 . . ?
 C2 C7 H7 119.2 . . ?
 N1 C8 C9 121.60(17) . . ?
 N1 C8 C1 115.30(15) . . ?
 C9 C8 C1 123.11(16) . . ?
 C10 C9 C8 119.51(19) . . ?
 C10 C9 H9 120.2 . . ?
 C8 C9 H9 120.2 . . ?
 C11 C10 C9 119.2(2) . . ?
 C11 C10 H10 120.4 . . ?

C9 C10 H10 120.4 . . ?
 C12 C11 C10 118.0(2) . . ?
 C12 C11 H11 121.0 . . ?
 C10 C11 H11 121.0 . . ?
 N1 C12 C11 124.2(2) . . ?
 N1 C12 H12 117.9 . . ?
 C11 C12 H12 117.9 . . ?
 C14 C13 C1 113.91(16) . . ?
 C14 C13 H13A 108.8 . . ?
 C1 C13 H13A 108.8 . . ?
 C14 C13 H13B 108.8 . . ?
 C1 C13 H13B 108.8 . . ?
 H13A C13 H13B 107.7 . . ?
 C13 C14 H14A 109.5 . . ?
 C13 C14 H14B 109.5 . . ?
 H14A C14 H14B 109.5 . . ?
 C13 C14 H14C 109.5 . . ?
 H14A C14 H14C 109.5 . . ?
 H14B C14 H14C 109.5 . . ?

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 _geom_torsion_site_symmetry_3
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 C8 C1 C2 C3 91.76(18) ?
 C13 C1 C2 C3 -145.89(16) ?
 O1 C1 C2 C7 154.81(15) ?
 C8 C1 C2 C7 -86.58(18) ?
 C13 C1 C2 C7 35.8(2) ?
 C7 C2 C3 C4 0.0(3) ?
 C1 C2 C3 C4 -178.40(16) ?
 C2 C3 C4 C5 0.2(3) ?
 C3 C4 C5 C6 -0.1(3) ?
 C4 C5 C6 C7 -0.3(3) ?
 C5 C6 C7 C2 0.6(3) ?
 C3 C2 C7 C6 -0.4(3) ?
 C1 C2 C7 C6 177.97(17) ?
 C12 N1 C8 C9 0.0(3) ?
 C12 N1 C8 C1 -179.68(17) ?
 O1 C1 C8 N1 3.5(2) ?
 C2 C1 C8 N1 -114.92(16) ?
 C13 C1 C8 N1 121.84(16) ?
 O1 C1 C8 C9 -176.15(15) ?
 C2 C1 C8 C9 65.4(2) ?
 C13 C1 C8 C9 -57.8(2) ?
 N1 C8 C9 C10 -0.8(3) ?
 C1 C8 C9 C10 178.86(16) ?
 C8 C9 C10 C11 1.1(3) ?
 C9 C10 C11 C12 -0.6(3) ?
 C8 N1 C12 C11 0.5(3) ?
 C10 C11 C12 N1 -0.2(4) ?
 O1 C1 C13 C14 58.9(2) ?

C2 C1 C13 C14 178.50(16) ?
 C8 C1 C13 C14 -60.1(2) ?

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 _geom_hbond_angle_DHA
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 _refine_diff_density_min -0.134
 _refine_diff_density_rms 0.031

_shelxl_version_number 2014-3

_shelx_res_file

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CELL 0.71073 9.139 8.2606 16.2236 90 103.14 90
ZERR 4 0.0016 0.0015 0.0029 0 0.002 0
LATT 1
SYMM 0.5-X,0.5+Y,0.5-Z
SFAC C H N O
UNIT 56 60 4 4
OMIT 1 0 5
OMIT 4 1 0
OMIT -2 2 7
OMIT 1 4 4
OMIT -1 3 2
OMIT 0 3 2
OMIT 0 0 2
OMIT -1 4 4
OMIT 3 3 1
OMIT -2 0 12
OMIT 0 50
L.S. 20
ACTA
BOND $H
FMAP 2
PLAN 2
HTAB O1 N1
HTAB
CONF
SIZE 0.32 0.26 0.20
TEMP 20
WGHT 0.065400
FVAR 0.71185
O1 4 -0.047769 0.474841 0.150273 11.00000 0.05820 0.04840 =
0.07600 -0.00300 0.00823 0.01161
AFIX 147
H1 2 -0.118580 0.457926 0.110035 11.00000 -1.50000
AFIX 0
N1 3 -0.173102 0.232126 0.062212 11.00000 0.06558 0.06241 =
0.05925 -0.00212 -0.00344 0.00301
C1 1 0.030565 0.327268 0.174968 11.00000 0.04761 0.04341 =
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0.05299 0.00194 0.00943 0.00700
 C2 1 0.188216 0.340460 0.157725 11.00000 0.04730 0.04156 =
 0.04870 -0.00318 0.00818 -0.00131
 C3 1 0.213805 0.442720 0.094644 11.00000 0.05972 0.05389 =
 0.05693 0.00614 0.01208 0.00280
 AFIX 43
 H3 2 0.135309 0.505027 0.063967 11.00000 -1.20000
 AFIX 0
 C4 1 0.355249 0.453391 0.076618 11.00000 0.07377 0.06080 =
 0.06805 0.00771 0.02813 -0.00523
 AFIX 43
 H4 2 0.370512 0.522847 0.034242 11.00000 -1.20000
 AFIX 0
 C5 1 0.472047 0.362267 0.120889 11.00000 0.05452 0.06723 =
 0.08082 -0.00757 0.02349 -0.00943
 AFIX 43
 H5 2 0.566395 0.369646 0.108615 11.00000 -1.20000
 AFIX 0
 C6 1 0.449370 0.259958 0.183468 11.00000 0.04655 0.06532 =
 0.07680 0.00640 0.00925 0.00370
 AFIX 43
 H6 2 0.528329 0.197452 0.213504 11.00000 -1.20000
 AFIX 0
 C7 1 0.309171 0.249781 0.201876 11.00000 0.04982 0.05647 =
 0.05944 0.00890 0.01000 -0.00041
 AFIX 43
 H7 2 0.295400 0.180736 0.244756 11.00000 -1.20000
 AFIX 0
 C8 1 -0.053740 0.188165 0.121576 11.00000 0.04544 0.05025 =
 0.04982 0.00149 0.01516 0.00055
 C9 1 -0.010886 0.028123 0.133725 11.00000 0.05860 0.04825 =
 0.08281 0.00345 0.01106 0.00119
 AFIX 43
 H9 2 0.073583 0.000086 0.175121 11.00000 -1.20000
 AFIX 0
 C10 1 -0.094028 -0.089595 0.084156 11.00000 0.07375 0.04947 =
 0.10447 -0.00642 0.03299 -0.00550
 AFIX 43
 H10 2 -0.067415 -0.198070 0.092284 11.00000 -1.20000
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 C11 1 -0.215965 -0.045219 0.022949 11.00000 0.08277 0.07431 =
 0.07707 -0.02284 0.02298 -0.02163
 AFIX 43
 H11 2 -0.273678 -0.122098 -0.011853 11.00000 -1.20000
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 C12 1 -0.250673 0.115299 0.014339 11.00000 0.08348 0.08451 =
 0.06978 -0.01198 -0.00825 -0.00473
 AFIX 43
 H12 2 -0.333675 0.145476 -0.027543 11.00000 -1.20000
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 C13 1 0.035961 0.299803 0.269396 11.00000 0.05829 0.06939 =
 0.05624 -0.00230 0.01636 0.00266
 AFIX 23
 H13A 2 0.089761 0.199971 0.287235 11.00000 -1.20000
 H13B 2 0.092128 0.387478 0.301740 11.00000 -1.20000
 AFIX 0
 C14 1 -0.118639 0.290224 0.290139 11.00000 0.07317 0.10804 =
 0.07648 0.00018 0.03236 0.00413
 AFIX 137
 H14A 2 -0.173340 0.200320 0.260649 11.00000 -1.50000

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H14B 2 -0.172849 0.388603 0.272809 11.00000 -1.50000
H14C 2 -0.106918 0.275726 0.349997 11.00000 -1.50000
AFIX 0
HKLf 4

REM 1 in P2(1)/n
REM R1 = 0.0441 for 1374 Fo > 4sig(Fo) and 0.0724 for all 2084 data
REM 147 parameters refined using 0 restraints

END

WGHT 0.0654 0.0000

REM Instructions for potential hydrogen bonds
HTAB O1 N1

REM Highest difference peak 0.117, deepest hole -0.134, 1-sigma level 0.031
Q1 1 -0.5060 0.0777 0.0134 11.00000 0.05 0.12
Q2 1 0.0775 0.0187 0.2368 11.00000 0.05 0.11
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_shelx_res_checksum 85945

_shelx_hkl_file
;
-1 0 0 -0.67 0.15
-1 0 0 -0.21 0.19
-2 0 0 640.88 14.78
-2 0 0 642.60 15.42
-3 0 0 0.52 0.41
-3 0 0 1.27 0.52
-4 0 0 480.93 12.11
-4 0 0 494.08 12.83
-5 0 0 0.51 0.62
-5 0 0 -0.12 0.58
-6 0 0 35.80 2.58
-6 0 0 38.14 2.37
-7 0 0 0.10 1.08
-7 0 0 -0.55 0.67
-8 0 0 63.41 4.06
-8 0 0 59.96 3.62
-9 0 0 -0.49 1.07
-9 0 0 0.64 1.52
-10 0 0 2.65 1.56
10 0 0 5.15 2.01
-10 0 0 3.56 2.12
11 0 0 0.40 1.60
0 1 0 0.59 0.24
0 1 0 0.36 0.24
1 1 0 120.67 3.36
1 -1 0 128.27 3.11
-1 1 0 129.85 3.41
-2 1 0 794.93 18.70
-2 1 0 792.54 17.96
2 1 0 756.89 18.57
-3 -1 0 1.10 0.55
-3 1 0 1.99 0.64
-3 -1 0 2.57 0.62
3 -1 0 2.14 0.35
-3 1 0 2.06 0.64
-4 -1 0 13.28 1.26

```

-4	1	0	13.46	1.23
4	-1	0	15.18	0.77
-4	-1	0	14.51	1.38
-4	1	0	13.59	1.35
-5	-1	0	16.10	1.56
5	-1	0	15.93	0.92
-5	1	0	14.87	1.59
-5	1	0	14.14	1.42
-5	-1	0	15.50	1.62
-6	-1	0	0.58	0.73
-6	1	0	0.28	0.64
-6	1	0	0.37	0.65
-6	-1	0	-0.33	0.89
-7	-1	0	34.29	2.82
-7	1	0	31.78	2.42
-7	-1	0	31.21	2.47
7	1	0	30.82	2.77
8	1	0	9.57	1.97
-8	1	0	7.44	1.81
-8	-1	0	10.48	1.84
-8	1	0	7.00	1.57
9	1	0	7.33	1.99
-9	1	0	5.39	1.99
-9	1	0	4.62	1.53
-10	1	0	3.96	2.22
10	1	0	3.39	1.87
-10	1	0	6.28	1.91
-11	1	0	0.47	2.14
11	1	0	1.11	1.85
0	2	0	348.07	8.45
0	2	0	337.98	8.53
-1	2	0	540.08	12.89
1	2	0	541.38	13.07
-1	2	0	544.55	13.12
1	2	0	535.56	12.81
2	2	0	35.70	1.63
-2	2	0	30.00	1.72
-2	2	0	29.77	1.63
-3	2	0	40.33	2.03
-3	2	0	42.83	2.08
3	-2	0	37.29	1.32
3	2	0	38.10	1.97
-4	2	0	26.51	1.87
4	2	0	26.84	1.80
4	-2	0	25.68	0.98
-4	2	0	24.58	1.78
5	2	0	114.64	4.58
-5	2	0	118.27	4.72
-5	2	0	121.64	4.39
-6	2	0	30.61	2.51
-6	2	0	31.27	2.28
6	2	0	29.56	2.48
-7	2	0	120.77	4.89
-7	2	0	112.50	5.34
7	2	0	120.86	5.37
-8	2	0	46.77	3.50
8	2	0	47.29	3.55
-8	2	0	46.25	3.07
-9	2	0	17.26	2.26
9	2	0	12.02	2.39

-9	2	0	17.67	2.69
-10	2	0	-0.11	1.41
10	2	0	1.63	1.70
-10	2	0	-0.93	1.59
-11	2	0	3.90	2.55
11	2	0	2.65	2.13
0	3	0	0.77	0.58
0	3	0	0.16	0.56
-1	3	0	16.96	1.27
-1	3	0	20.48	1.55
1	3	0	13.86	1.22
1	3	0	17.23	1.35
-2	3	0	616.75	15.13
2	3	0	613.26	15.26
2	3	0	619.24	15.08
-2	3	0	607.09	15.34
3	3	0	116.50	4.14
-3	3	0	113.47	4.20
-3	3	0	122.42	4.29
-4	3	0	184.22	6.22
-4	3	0	185.96	5.94
4	3	0	195.38	6.18
-5	3	0	53.77	3.06
5	3	0	56.63	3.12
-5	3	0	58.76	3.04
-6	3	0	28.20	2.54
-6	3	0	20.05	2.00
6	3	0	22.51	2.21
-7	3	0	28.05	2.82
-7	3	0	28.29	2.46
7	3	0	29.90	2.78
-8	3	0	33.08	2.78
8	3	0	29.11	2.96
-8	3	0	34.98	3.27
9	3	0	17.27	2.66
-9	3	0	12.93	2.20
-9	3	0	9.49	2.28
10	3	0	4.91	2.11
-10	3	0	1.93	1.77
11	3	0	2.43	2.40
0	4	0	498.77	12.75
0	4	0	501.38	12.85
-1	4	0	74.24	3.07
1	4	0	78.39	3.25
1	4	0	74.67	3.07
-1	4	0	78.30	3.33
-2	4	0	43.35	2.33
2	4	0	43.30	2.41
-2	4	0	45.52	2.65
2	4	0	41.51	2.30
-3	4	0	211.26	6.73
3	4	0	213.47	6.83
-3	4	0	212.87	6.89
3	4	0	207.42	6.61
-4	4	0	1.43	0.89
-4	4	0	1.91	1.21
4	4	0	1.05	0.98
-5	4	0	155.21	5.77
-5	4	0	157.95	6.09
5	4	0	163.94	6.02

6	4	0	22.60	2.38
-6	4	0	22.69	2.28
-6	4	0	25.53	2.54
-7	4	0	87.07	4.81
-7	4	0	95.93	4.52
7	4	0	92.22	4.82
-8	4	0	9.01	2.14
8	4	0	9.79	2.16
-8	4	0	9.59	1.92
9	4	0	-0.54	1.47
-9	4	0	-1.42	1.24
-10	4	0	2.30	1.91
10	4	0	-2.12	1.78
0	5	0	0.66	0.86
0	5	0	0.50	0.82
-1	5	0	86.01	3.89
1	5	0	85.49	3.58
-1	5	0	80.19	3.47
1	5	0	79.11	3.71
2	5	0	244.88	7.75
-2	5	0	246.52	7.55
2	5	0	236.01	7.48
-2	5	0	248.31	7.92
3	5	0	22.40	2.18
3	5	0	19.59	2.04
-3	5	0	23.41	2.02
-3	5	0	18.27	2.21
-4	5	0	-0.98	0.77
-4	5	0	2.98	1.45
4	5	0	0.91	1.08
4	5	0	-0.03	0.95
5	5	0	4.07	1.50
-5	5	0	4.24	1.40
-5	5	0	5.42	1.62
-6	5	0	10.31	1.96
-6	5	0	9.64	1.80
6	5	0	10.38	2.03
-7	5	0	4.22	1.68
-7	5	0	-0.56	1.05
7	5	0	0.53	1.33
-8	5	0	29.35	3.25
8	5	0	23.48	3.00
9	5	0	-0.07	1.92
-9	5	0	-1.25	1.38
-10	5	0	3.02	2.34
10	5	0	-0.58	2.37
0	6	0	3.46	1.17
0	6	0	4.94	1.46
1	6	0	0.48	1.09
1	6	0	0.81	1.08
-1	6	0	0.54	1.22
-1	6	0	-0.82	0.88
2	6	0	140.52	5.76
-2	6	0	141.25	5.93
2	6	0	138.32	5.32
-2	6	0	135.38	5.30
3	6	0	3.83	1.51
3	6	0	3.86	1.43
-3	6	0	3.16	1.24
-3	6	0	3.07	1.84

-4	6	0	47.22	3.22
-4	6	0	48.16	3.82
4	6	0	51.98	3.58
4	6	0	49.67	3.34
5	6	0	4.21	1.65
-5	6	0	0.98	1.19
5	6	0	3.96	1.64
-5	6	0	2.58	1.71
-6	6	0	4.46	1.67
6	6	0	4.13	1.77
-6	6	0	5.09	1.83
6	6	0	4.67	1.91
-7	6	0	10.04	2.27
7	6	0	10.13	2.52
-8	6	0	1.36	1.71
-9	6	0	-0.63	1.66
0	7	0	0.71	1.12
0	7	0	-1.65	1.33
1	7	0	8.81	1.93
-1	7	0	10.89	1.74
-1	7	0	7.55	2.10
-2	7	0	1.03	1.17
2	7	0	3.48	1.50
-2	7	0	-0.13	1.58
3	7	0	17.61	2.55
-3	7	0	21.83	2.35
-3	7	0	20.61	3.04
-4	7	0	55.02	3.62
4	7	0	51.74	3.92
-4	7	0	46.41	4.09
-5	7	0	4.66	2.10
5	7	0	1.73	1.67
-5	7	0	0.14	1.35
-6	7	0	18.73	2.71
6	7	0	18.71	3.06
-7	7	0	3.33	1.84
7	7	0	3.37	2.24
-8	7	0	1.19	1.93
8	7	0	-0.04	1.79
-9	7	0	4.13	2.67
0	8	0	5.53	1.95
1	8	0	0.75	1.42
-1	8	0	0.02	1.80
2	8	0	-1.08	1.48
-2	8	0	-0.56	1.85
3	8	0	-0.23	1.60
-3	8	0	3.35	2.41
4	8	0	-0.53	1.47
-4	8	0	0.39	2.16
5	8	0	7.81	2.51
6	8	0	2.95	2.11
7	8	0	1.09	2.05
0	9	0	0.81	1.90
-1	9	0	1.98	2.23
1	9	0	-0.76	1.71
-2	9	0	13.06	3.25
2	9	0	13.44	2.91
-3	9	0	0.44	2.49
3	9	0	1.91	2.21
4	9	0	3.62	2.36

5	9	0	5.95	2.65
0	10	0	6.09	2.72
-1	10	0	1.98	2.86
1	10	0	3.71	2.73
-2	10	0	8.90	3.28
2	10	0	12.75	3.31
3	10	0	2.07	2.40
4	10	0	1.85	2.71
0	11	0	-2.20	2.86
-11	0	1	17.31	3.32
11	0	-1	18.88	3.26
-10	0	1	-0.27	1.13
-10	0	1	1.28	1.97
-9	0	1	1.68	1.52
-9	0	1	-0.29	1.07
-8	0	1	0.47	1.02
-8	0	1	0.59	1.29
-7	0	1	7.03	1.39
-7	0	1	7.53	1.64
-6	0	1	0.54	0.67
-6	0	1	0.28	0.75
-5	0	1	277.33	7.78
-5	0	1	281.73	8.43
-4	0	1	1.32	0.64
-4	0	1	0.01	0.41
-3	0	1	431.95	10.55
-3	0	1	441.68	11.18
-2	0	1	0.19	0.37
-2	0	1	-0.29	0.46
-1	0	1	2662.02	61.37
-1	0	1	2672.70	59.72
-1	0	-1	202.01	5.11
-2	0	-1	0.50	0.47
-2	0	-1	-0.48	0.31
-3	0	-1	109.35	3.53
-4	0	-1	-0.25	0.38
-4	0	-1	-0.40	0.48
-5	0	-1	61.58	3.07
-5	0	-1	54.42	2.70
-6	0	-1	-0.14	0.82
-6	0	-1	0.05	0.69
-7	0	-1	289.07	9.02
-8	0	-1	0.15	1.29
-8	0	-1	-1.39	0.90
9	0	1	1.09	1.43
-9	0	-1	0.09	1.20
10	0	1	-0.64	1.20
-10	0	-1	1.30	1.44
11	0	1	0.48	1.86
-11	0	-1	3.25	1.88
11	1	-1	8.58	2.69
-11	1	1	8.02	2.86
-10	1	1	1.17	1.55
-10	1	1	2.34	1.84
10	1	-1	-0.33	1.42
-9	1	1	0.41	1.40
-9	-1	1	-1.78	1.21
-9	1	1	0.83	1.10
9	1	-1	1.49	1.35
-8	1	1	24.88	2.71

-8	-1	1	26.01	2.87
-8	1	1	29.38	2.47
-7	1	1	1.12	0.91
-7	1	1	0.57	1.03
-7	-1	1	4.01	1.44
-6	1	1	90.86	3.73
-6	1	1	93.42	4.24
-6	-1	1	96.94	4.25
-5	-1	1	2.01	0.73
-5	1	1	1.70	0.71
5	-1	-1	1.65	0.54
-5	1	1	0.99	0.63
-5	-1	1	1.03	0.72
-4	1	1	276.45	7.36
4	-1	-1	270.46	6.78
-4	-1	1	281.56	7.90
-3	1	1	59.56	2.22
3	-1	-1	63.28	1.95
-3	-1	1	59.77	2.41
-2	1	1	81.91	2.88
-2	1	1	90.73	2.63
-2	-1	1	90.93	2.72
2	-1	-1	86.73	2.46
-1	1	1	532.71	12.28
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1	1	-1	527.17	12.39
1	-1	-1	538.02	12.40
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-1	1	-1	861.32	20.05
-1	1	-1	876.85	19.62
1	1	1	838.46	20.16
-2	1	-1	79.42	2.79
-2	1	-1	84.82	2.79
2	1	1	76.18	2.75
-3	-1	-1	15.37	1.18
-3	1	-1	16.85	1.28
-3	1	-1	15.49	1.26
3	1	1	16.39	1.27
-4	1	-1	1.94	0.68
-4	1	-1	1.59	0.69
-4	-1	-1	1.28	0.65
4	1	1	1.15	0.66
5	1	1	4.60	1.18
-5	-1	-1	3.02	0.90
-5	1	-1	2.02	0.87
-5	1	-1	2.80	0.97
-6	-1	-1	4.04	1.13
-6	1	-1	4.39	1.27
-6	1	-1	4.26	1.18
6	1	1	4.37	1.29
-7	1	-1	8.25	1.78
-7	1	-1	6.89	1.49
-7	-1	-1	6.35	1.48
7	1	1	7.49	1.63
-8	1	-1	0.56	1.15
8	1	1	-2.10	1.03
-8	1	-1	0.00	1.23
-8	-1	-1	1.23	1.08

-9	-1	-1	1.11	1.32
-9	1	-1	0.08	1.40
9	1	1	-0.24	1.27
-9	1	-1	-0.31	1.07
-10	1	-1	3.07	1.66
-10	1	-1	-0.76	1.32
10	1	1	1.40	1.50
11	1	1	1.49	1.89
-11	1	-1	2.31	2.01
11	-1	1	4.36	2.33
-11	2	1	-0.05	1.84
11	2	-1	5.11	2.27
-10	2	1	-1.04	1.31
10	2	-1	-2.26	1.33
9	2	-1	3.64	1.69
-9	2	1	6.88	1.94
-9	2	1	3.81	1.52
-8	2	1	1.21	1.12
-8	2	1	1.04	1.24
8	2	-1	1.30	1.17
7	2	-1	255.26	8.73
-7	2	1	261.39	8.01
-7	2	1	252.03	8.85
6	2	-1	3.75	1.16
-6	2	1	4.54	1.04
-6	2	1	2.24	1.07
5	-2	-1	194.04	4.98
5	2	-1	194.10	6.38
-5	2	1	192.61	5.93
-5	2	1	186.32	6.50
4	2	-1	69.19	2.98
-4	2	1	69.78	3.14
4	-2	-1	69.03	2.27
-4	2	1	69.96	2.83
3	-2	-1	21.41	1.16
-3	2	1	19.09	1.38
-3	2	1	18.84	1.43
3	2	-1	19.32	1.32
2	2	-1	263.23	6.73
2	-2	-1	258.18	6.46
-2	2	1	255.58	6.71
-2	2	1	251.13	7.01
-1	2	1	86.28	2.97
1	-2	-1	87.97	2.79
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-1	2	1	88.57	2.79
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-2	2	-1	28.89	1.54
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-3	2	-1	439.53	11.35
-3	2	-1	420.20	10.98
3	2	1	431.32	11.34

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-4	2	-1	20.85	1.73
-5	2	-1	334.21	9.54
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-5	2	-1	338.11	10.10
-6	2	-1	28.35	2.50
6	2	1	28.62	2.46
-6	2	-1	30.40	2.33
7	2	1	38.59	3.06
-7	2	-1	30.05	2.81
-7	2	-1	38.54	2.76
-8	2	-1	4.02	1.46
-8	2	-1	2.10	1.45
8	2	1	1.65	1.32
-9	2	-1	18.58	2.49
9	2	1	22.54	2.97
-9	2	-1	20.36	2.85
-10	2	-1	1.96	1.82
-10	2	-1	0.18	1.49
10	2	1	1.18	1.72
11	2	1	0.44	2.14
-11	2	-1	-0.45	1.99
-11	3	1	-0.05	1.82
11	3	-1	-0.18	2.09
10	3	-1	3.24	2.09
-10	3	1	2.36	1.82
-9	3	1	3.65	1.66
9	3	-1	-1.28	1.23
-8	3	1	38.16	2.86
8	3	-1	42.21	3.40
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7	3	-1	42.45	3.21
-7	3	1	43.35	3.26
-7	3	1	46.84	2.96
6	3	-1	216.49	7.41
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-6	3	1	207.85	7.57
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-4	3	1	224.43	7.08
4	-3	-1	215.78	5.71
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-4	3	1	216.44	6.49
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2	3	-1	452.42	11.18
2	-3	-1	442.48	10.88
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-1	3	1	408.98	10.23
1	3	-1	410.76	10.15
0	3	1	160.00	4.76
0	3	-1	159.93	4.51
0	3	-1	161.35	4.94

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-2	3	-1	107.81	3.74
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-3	3	-1	-0.19	0.81
4	3	1	3.41	0.97
-4	3	-1	0.63	0.73
-4	3	-1	3.62	1.13
5	3	1	1.86	1.00
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-5	3	-1	1.07	0.83
-6	3	-1	25.22	2.40
6	3	1	28.02	2.53
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7	3	1	7.50	1.81
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-8	3	-1	3.81	1.48
-8	3	-1	6.41	1.90
8	3	1	8.43	2.06
9	3	1	5.83	2.14
-9	3	-1	5.25	1.86
-9	3	-1	6.81	2.12
10	3	1	8.91	2.44
-10	3	-1	9.10	2.45
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9	4	-1	0.64	1.67
8	4	-1	1.12	1.54
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3	4	-1	3.01	0.91
2	4	-1	67.09	3.11
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-2	4	1	67.60	3.07
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9	5	1	3.20	2.15
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10	5	1	-0.37	2.20
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3	6	-1	24.51	2.28
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0	6	-1	24.04	2.42
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1	6	1	8.26	1.63
1	6	1	9.78	1.66
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2	6	1	0.45	1.10
2	6	1	-0.92	1.12
-2	6	-1	0.26	0.88
3	6	1	32.28	2.74
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4	6	1	0.99	1.46
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-9	6	-1	3.14	2.32
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5	9	1	-0.25	2.10
4	10	-1	0.05	2.40
3	10	-1	2.23	2.60
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3	10	1	3.41	2.69
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-3	0	-2	-0.36	0.41
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9	0	2	-0.36	1.09
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2	5	2	14.43	1.84
3	5	2	2.34	1.04
3	5	2	1.07	1.18
-3	5	-2	0.55	0.76
-3	5	-2	0.85	1.39
-4	5	-2	-0.43	1.44
-4	5	-2	1.20	0.88
4	5	2	-1.18	1.12
-5	5	-2	5.45	1.85
-5	5	-2	6.72	1.52
5	5	2	5.39	1.70
-6	5	-2	51.69	3.57
6	5	2	45.62	3.77
-6	5	-2	48.06	3.52

7	5	2	5.56	1.99
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-7	5	-2	3.61	1.67
8	5	2	0.48	1.76
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-8	5	-2	2.66	1.49
-9	5	-2	-0.21	1.72
9	5	2	0.55	2.04
10	5	2	-0.05	2.20
9	6	-2	7.60	2.79
-9	6	2	3.21	2.28
8	6	-2	0.34	1.69
-8	6	2	-1.68	1.24
-7	6	2	3.47	1.71
7	6	-2	9.96	2.55
-6	6	2	0.78	1.32
6	6	-2	1.63	1.46
-5	6	2	36.92	3.20
5	6	-2	34.50	3.34
-4	6	2	29.15	2.70
4	6	-2	18.38	2.51
-4	6	2	23.91	2.71
-3	6	2	-0.56	1.03
3	6	-2	0.30	1.10
-3	6	2	1.08	1.54
-2	6	2	27.66	2.43
2	6	-2	28.55	2.65
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0	6	2	45.70	2.95
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3	6	2	35.51	2.93
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-4	6	-2	11.94	2.51
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4	6	2	12.13	2.22
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7	6	2	7.11	2.34

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7	7	-2	-1.08	1.59
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3	7	-2	12.43	2.34
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3	7	2	0.29	1.50
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7	8	2	7.73	3.06
6	9	-2	-0.85	1.96
5	9	-2	0.06	2.03
4	9	-2	-0.98	2.08
3	9	-2	1.83	2.13
2	9	-2	2.89	2.15
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5	9	2	-0.73	2.24
3	10	-2	0.59	2.20
2	10	-2	-1.16	2.36
1	10	-2	12.14	3.19
0	10	2	1.72	2.45
0	10	-2	-0.51	2.57
1	10	2	-3.74	2.02
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-2	10	-2	-0.56	2.99
2	10	2	-0.30	2.15
3	10	2	-0.60	2.59
-3	10	-2	-2.08	2.86
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-10	0	3	-0.24	1.45
-9	0	3	15.17	2.49
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-7	0	3	67.33	3.73
-7	0	3	70.43	3.33
-6	0	3	-0.13	0.72

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-5	0	3	567.76	14.95
5	0	-3	573.35	14.07
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-4	0	3	0.24	0.58
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0	0	-3	-0.12	0.33
0	0	3	0.60	0.34
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-1	0	-3	0.75	0.53
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-2	0	-3	-0.38	0.41
2	0	3	0.40	0.44
3	0	3	12.91	1.25
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6	0	3	-0.95	0.73
7	0	3	18.07	2.19
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8	0	3	0.19	0.96
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-9	1	3	9.16	2.19
-8	1	3	7.66	1.82
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-7	1	3	2.90	1.02
-7	1	3	4.21	1.38
6	-1	-3	45.78	2.23
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9	1	3	10.37	2.32
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3	4	3	206.55	7.07
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0	3	-4	1.34	0.92
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-2	3	-4	2.50	1.21
-2	-3	-4	3.55	0.98
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-1	-2	-5	4.61	0.74
-2	-2	-5	65.65	2.91
-2	2	-5	77.99	2.55
2	-2	5	72.94	2.81
-2	2	-5	72.21	3.34
-2	-2	-5	68.51	2.89
-3	-2	-5	273.04	7.95
-3	2	-5	270.17	6.82
-3	2	-5	266.65	7.96
3	2	5	287.23	8.62
-3	-2	-5	283.66	7.85
-4	2	-5	0.79	0.85
-4	-2	-5	1.60	0.89
4	2	5	2.23	1.04
-4	2	-5	2.38	1.22
5	2	5	0.73	1.07
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-5	-2	-5	1.64	0.94
-5	2	-5	2.23	1.10

-6	2	-5	24.00	2.44
-6	2	-5	22.73	2.41
6	-2	5	21.75	2.37
6	2	5	23.75	2.58
-6	-2	-5	24.13	2.45
-7	2	-5	61.62	3.80
7	-2	5	63.22	4.00
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8	-2	5	4.00	1.65
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9	-2	5	14.59	2.71
9	2	5	14.95	2.82
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9	3	-5	0.41	1.48
8	3	-5	12.69	2.10
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-7	3	5	1.77	1.30
7	3	-5	3.43	1.41
6	-3	-5	60.23	3.09
6	3	-5	56.09	3.13
6	3	-5	59.93	3.44
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5	-3	-5	-0.06	1.02
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4	-3	-5	22.04	1.85
4	3	-5	24.58	1.51
4	3	-5	23.74	1.83
-4	3	5	19.08	1.96
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3	-3	-5	135.11	4.74
3	3	-5	133.55	3.79
3	3	-5	128.10	4.69
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-6	1	-6	4.67	1.42

-7	1	-6	8.05	1.90
7	-1	6	12.01	2.16
7	1	6	13.20	2.31
-8	-1	-6	1.86	1.47
8	-1	6	-1.37	1.17
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8	1	6	-0.81	1.17
-9	-1	-6	10.66	2.35
-9	1	-6	4.37	1.87
9	1	6	0.56	1.37
9	-1	6	5.48	2.32
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-10	1	-6	-0.58	1.68
10	-1	6	-1.30	1.86
-10	-1	-6	0.92	1.90
11	2	-6	4.67	2.61
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-8	-2	6	7.36	1.76
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6	2	-6	53.49	2.95
5	2	-6	5.75	1.10
-5	2	6	4.04	1.21
5	-2	-6	5.47	1.26
-5	-2	6	4.32	0.94
4	-2	-6	30.65	2.16
-4	2	6	31.61	2.25
-4	-2	6	33.06	1.76
4	2	-6	29.23	1.89
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-7	-2	-6	1.02	1.26
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-8	-2	-6	2.42	1.32
8	-2	6	5.61	2.01
9	2	6	-1.71	1.57
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8	3	6	23.82	3.36
8	-3	6	22.06	2.93
9	-3	6	2.14	2.01
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-3	-1	7	0.72	0.55
3	-1	-7	2.40	0.95
3	1	-7	0.43	0.62
2	-1	-7	209.15	6.48
2	1	-7	203.92	5.85
-2	-1	7	210.71	5.68
2	1	-7	209.29	6.38
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1	1	-7	27.77	1.87
1	-1	-7	29.65	2.06
1	-1	-7	29.65	1.25
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1	1	-7	28.12	1.83
-1	-1	7	27.52	1.56
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0	-1	-7	1.55	0.57
0	-1	7	1.39	0.68
0	1	-7	0.16	0.62
0	1	7	2.73	0.95
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1	1	7	13.43	1.54
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2	1	7	193.53	6.50
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4	1	7	1.26	1.05

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8	1	7	16.82	2.90
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8	-1	7	14.51	2.71
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0	-2	-7	0.98	0.92
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-9	-1	8	3.80	1.83
8	1	-8	5.86	1.58
-8	-1	8	4.02	1.57
-8	1	8	4.11	1.61
7	-1	-8	0.95	1.30
-7	1	8	0.36	1.11
7	1	-8	0.05	0.90
-7	-1	8	-0.65	0.87
6	1	-8	0.19	0.93
6	1	-8	0.30	0.73
6	-1	-8	2.96	1.29
-6	1	8	1.18	1.07
-6	-1	8	1.05	0.91
5	1	-8	49.67	2.39
5	-1	-8	51.88	3.03
-5	1	8	46.62	2.83
5	1	-8	42.39	2.79
-5	-1	8	47.87	2.59
4	1	-8	113.54	3.81
4	1	-8	111.01	4.55
-4	-1	8	116.05	4.02
4	-1	-8	114.61	4.60
-4	1	8	110.83	4.46
-3	1	8	497.21	13.02
3	-1	-8	478.67	13.03
-3	-1	8	489.45	12.28

3	1	-8	494.89	12.05
3	1	-8	465.02	12.93
2	-1	-8	7.56	0.59
2	1	-8	5.09	1.09
2	-1	-8	7.07	1.34
-2	1	8	6.81	1.19
2	1	-8	5.32	1.09
-2	-1	8	6.32	0.90
1	1	-8	71.84	3.26
-1	-1	8	78.39	2.91
1	1	-8	75.14	3.15
1	-1	-8	76.24	2.43
-1	1	8	77.51	3.48
1	-1	-8	70.45	3.36
0	-1	-8	232.26	6.34
0	-1	-8	242.00	7.34
0	-1	8	236.75	6.81
0	1	8	235.46	7.40
0	1	-8	242.92	7.10
0	1	-8	236.47	7.01
-1	-1	-8	5.56	1.20
-1	1	-8	6.95	1.33
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-1	1	-8	7.65	1.18
1	1	8	9.92	1.51
-1	-1	-8	7.74	1.04
-2	-1	-8	174.89	5.71
2	1	8	179.76	6.20
-2	1	-8	174.70	6.01
-2	-1	-8	179.17	5.69
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-2	1	-8	168.19	5.43
-3	1	-8	2.40	1.02
-3	-1	-8	1.33	1.02
3	1	8	-0.35	0.99
-3	1	-8	1.60	1.06
-3	-1	-8	1.25	0.84
3	-1	8	1.56	0.90
-4	-1	-8	59.38	3.28
4	1	8	56.69	3.49
4	-1	8	50.17	3.12
-4	-1	-8	52.21	2.86
-4	1	-8	46.41	3.14
5	1	8	5.34	1.66
5	-1	8	3.51	1.53
-5	1	-8	4.93	1.43
-5	-1	-8	6.88	1.53
-6	-1	-8	-0.15	0.97
6	-1	8	0.80	1.42
6	1	8	1.86	1.46
-6	1	-8	1.58	1.20
-7	-1	-8	-0.52	1.06
7	1	8	1.27	1.57
-7	1	-8	1.18	1.02
7	-1	8	0.00	1.39
-8	1	-8	1.06	1.59
8	1	8	0.63	1.53
8	-1	8	-0.64	1.70
-8	-1	-8	-0.69	1.29
9	1	8	1.87	2.02

9	-1	8	-2.99	1.63
-9	-1	-8	0.01	1.39
-9	1	-8	3.77	1.84
-11	2	8	-3.18	1.70
-10	2	8	0.88	2.00
-9	2	8	7.68	2.19
9	2	-8	9.50	1.96
-8	2	8	4.35	1.70
8	2	-8	4.99	1.47
7	2	-8	-0.24	0.93
7	-2	-8	-1.58	1.18
7	2	-8	1.24	1.11
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-6	-2	8	139.17	5.06
6	2	-8	135.63	5.77
6	-2	-8	139.17	5.52
-6	2	8	137.59	5.77
6	2	-8	138.01	5.04
5	2	-8	2.86	0.97
5	-2	-8	2.65	1.27
-5	-2	8	3.63	0.92
-5	2	8	-0.06	0.95
-4	-2	8	35.49	1.88
4	2	-8	36.23	2.22
4	2	-8	38.28	2.66
4	-2	-8	39.48	2.63
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3	2	-8	18.68	1.69
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2	2	-8	62.82	3.13
-2	2	8	68.22	3.31
2	-2	-8	62.11	3.29
-2	-2	8	65.08	2.20
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-2	-2	-8	16.90	1.82
2	-2	8	18.61	1.55
-2	2	-8	14.20	1.87
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-3	-2	-8	1.29	1.05

-3	2	-8	1.60	1.24
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3	-2	8	0.12	0.80
-4	-2	-8	8.07	1.51
4	-2	8	7.94	1.54
-4	-2	-8	5.72	1.41
4	2	8	9.39	1.92
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5	-2	8	53.60	3.54
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-5	-2	-8	52.99	3.40
-6	-2	-8	159.12	6.61
6	-2	8	155.11	6.70
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-7	-2	-8	1.44	1.44
7	-2	8	1.28	1.55
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8	2	8	-1.49	1.32
8	-2	8	1.72	1.92
-8	2	-8	1.26	1.54
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9	2	8	-1.20	1.57
9	-2	8	0.52	2.13
-9	-2	-8	1.31	1.74
-9	2	-8	2.24	1.84
-11	3	8	7.11	2.92
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5	-3	-8	44.01	3.06
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4	-3	-8	2.12	1.36
4	3	-8	1.67	1.05
3	-3	-8	8.94	1.65
3	3	-8	8.97	1.52
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2	-3	-8	151.76	5.74
-2	3	8	156.39	5.84
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2	-3	8	39.10	2.23
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3	-3	8	20.07	1.88
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-3	-3	-8	20.88	2.11
-4	-3	-8	-0.24	1.08
-4	3	-8	-0.22	1.20
4	-3	8	-0.18	0.84
4	3	8	-0.51	1.19
5	3	8	3.59	1.73
5	-3	8	5.31	1.57
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6	3	8	17.94	2.84
6	-3	8	16.80	2.39
7	3	8	12.87	2.75
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8	-3	8	4.11	2.23
8	3	8	2.65	2.17
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5	4	-8	3.01	1.43
4	-4	-8	11.20	1.96
4	4	-8	15.67	1.96
-4	4	8	5.27	1.32
3	-4	-8	12.52	2.02
3	4	-8	14.26	1.99
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3	5	-8	4.58	1.62
3	-5	-8	1.72	1.42
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-2	5	8	12.29	2.10
2	-5	-8	10.06	1.89
2	5	-8	8.23	1.96
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4	5	8	3.69	2.01
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5	6	-8	13.25	2.57
5	-6	-8	12.57	2.14
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4	-6	-8	6.06	1.76
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-7	0	-9	1.64	1.42
7	0	9	1.70	1.50
8	0	9	-1.45	1.67
-8	0	-9	0.47	1.37
-9	0	-9	0.57	1.51
9	0	9	6.61	2.57
-11	1	9	2.19	2.21
-10	1	9	4.33	2.10
-9	-1	9	5.44	1.86
-9	1	9	4.23	1.83
9	1	-9	4.16	1.72
-8	1	9	5.11	1.85

-8	-1	9	4.99	1.62
8	-1	-9	6.46	1.80
8	1	-9	3.26	1.48
-7	-1	9	1.35	1.08
7	1	-9	1.77	1.25
-7	1	9	0.35	1.13
7	-1	-9	-0.15	1.21
-6	1	9	10.41	1.81
6	-1	-9	9.43	1.80
-6	-1	9	8.72	1.51
6	1	-9	9.75	1.77
5	-1	-9	158.42	6.00
5	1	-9	151.86	4.94
5	1	-9	157.47	6.04
-5	1	9	156.76	5.78
-5	-1	9	164.98	5.47
4	-1	-9	94.20	4.22
-4	-1	9	90.43	3.50
4	1	-9	84.80	4.06
-4	1	9	87.68	3.91
4	1	-9	91.69	3.40
3	-1	-9	45.19	2.85
3	1	-9	47.00	2.45
-3	-1	9	47.97	2.27
-3	1	9	45.47	2.68
-2	-1	9	17.74	1.39
2	1	-9	17.57	1.64
2	1	-9	19.05	1.85
2	-1	-9	18.97	0.98
-2	1	9	19.22	1.81
1	1	-9	51.28	2.75
-1	-1	9	48.15	2.41
-1	1	9	46.61	2.77
1	1	-9	49.58	2.69
1	-1	-9	52.92	3.01
1	-1	-9	48.36	1.96
0	-1	9	191.19	5.91
0	1	-9	191.54	6.22
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-1	1	-9	61.97	2.97
1	-1	9	55.91	2.79
-1	-1	-9	55.88	2.64
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-2	1	-9	8.13	1.40
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2	1	9	6.63	1.54
2	-1	9	7.85	1.39
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-3	-1	-9	7.77	1.48
-3	1	-9	9.31	1.69
3	-1	9	9.14	1.53
3	1	9	10.15	1.81
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4	-1	9	5.30	1.46

-4	-1	-9	5.56	1.37
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5	1	9	50.75	3.60
5	-1	9	44.88	3.27
-5	-1	-9	52.37	3.38
-5	1	-9	43.87	3.24
-6	-1	-9	0.33	1.20
-6	1	-9	5.24	1.64
6	1	9	2.56	1.67
6	-1	9	2.33	1.62
7	1	9	12.45	2.68
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7	-1	9	13.45	2.56
-8	-1	-9	-2.72	1.22
8	-1	9	-0.77	1.51
8	1	9	-1.35	1.64
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9	-1	9	12.07	2.88
9	1	9	6.20	2.77
-9	1	-9	15.55	2.85
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9	2	-9	0.22	1.32
-8	2	9	0.58	1.48
8	2	-9	1.40	1.33
7	2	-9	10.03	1.85
-7	2	9	6.43	1.87
7	-2	-9	5.76	1.75
6	2	-9	75.92	4.06
6	-2	-9	76.35	4.11
-6	2	9	73.60	4.13
6	2	-9	77.47	3.58
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4	2	-9	3.85	1.18
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5	2	9	39.77	3.60
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6	2	9	0.84	1.54
6	-2	9	0.53	1.52
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-7	-2	-9	0.48	1.34
7	2	9	0.64	1.76
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8	2	9	3.34	2.09
8	-2	9	7.50	2.57
-8	2	-9	11.50	2.39
9	-2	9	-3.30	1.95
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9	3	-9	1.75	1.50
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7	-3	-9	7.83	1.92
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4	3	-9	0.48	1.07

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-7	3	-9	17.11	2.73
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6	0	10	10.01	2.32
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2	-1	11	2.23	1.27
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-3	-1	-11	101.92	4.56
3	-1	11	93.10	4.67
-3	-1	-11	94.73	4.40
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4	-1	11	0.82	1.31
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5	-1	11	7.51	2.07
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6	-1	11	-0.07	1.72
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7	-1	11	20.28	3.29
-7	-1	-11	22.24	3.07
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10	2	-11	0.84	1.92

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8	2	-11	0.78	1.29
8	-2	-11	0.26	1.61
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7	2	-11	24.03	2.72
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-6	2	11	28.54	2.81
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3	2	-11	3.58	1.38
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7	-2	11	2.11	1.82
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9	0	-12	-1.21	1.57
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7	0	-12	0.04	1.43
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6	0	-12	1.09	1.37
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2	4	-13	2.83	1.60
2	-4	-13	4.22	1.87
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1	-4	-13	2.15	1.61
1	4	-13	3.94	1.68
0	-4	-13	0.57	1.60

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0	4	-13	0.83	1.79
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7	5	-13	2.44	1.71
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8	0	-15	-0.88	1.79
7	0	-15	4.17	2.12
6	0	-15	-2.15	1.45
5	0	-15	59.80	4.43
4	0	-15	-0.41	1.04
3	0	-15	6.13	2.08
2	0	-15	0.88	1.59
1	0	-15	-0.26	1.31
0	0	-15	-0.17	1.63
1	0	15	-2.75	1.75
-1	0	-15	-0.91	1.53
2	0	15	2.29	2.07
-2	0	-15	-0.08	1.29
-3	0	-15	25.05	2.83
3	0	15	16.18	3.04
-4	0	-15	-0.29	1.40
4	0	15	1.50	2.33
-5	0	-15	5.16	2.07
5	0	15	2.63	2.55
-6	0	-15	-0.16	1.65
6	0	15	-4.61	2.40
10	1	-15	0.96	2.36
9	1	-15	-0.22	2.04
9	-1	-15	-2.11	2.06
8	-1	-15	4.92	2.51
8	1	-15	6.36	2.35
7	-1	-15	2.91	2.15
7	1	-15	-1.49	1.36
6	1	-15	1.68	1.68
6	-1	-15	-1.08	1.52
5	1	-15	1.49	1.52
5	-1	-15	1.44	1.66
-5	1	15	2.67	1.66
4	-1	-15	22.21	3.06
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4	1	-15	19.94	2.78
3	-1	-15	4.41	1.81
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3	1	-15	3.33	1.75
-2	1	15	1.38	1.60
2	1	-15	1.85	1.59
2	-1	-15	1.89	1.77
1	-1	-15	4.54	1.93
1	1	-15	3.68	1.74
-1	1	15	3.65	2.09
0	-1	-15	11.18	2.37
0	1	15	8.57	2.31
0	1	-15	7.60	1.99
-1	-1	-15	1.72	1.79
-1	1	-15	0.84	1.23
1	1	15	0.20	1.85
-1	1	-15	3.76	1.87
-2	-1	-15	-1.43	1.61
-2	1	-15	0.69	1.34
2	1	15	2.15	2.15
-3	-1	-15	1.68	1.37
3	1	15	0.27	1.99

-3	1	-15	0.73	1.42
-4	-1	-15	-0.27	1.40
-4	1	-15	0.68	1.67
4	1	15	1.69	2.44
-5	-1	-15	-0.69	1.67
5	1	15	3.27	2.48
-5	1	-15	2.04	1.71
6	1	15	-0.83	2.49
-6	1	-15	-1.02	1.69
-6	-1	-15	2.95	1.88
10	2	-15	-0.49	2.23
9	2	-15	5.47	2.35
9	-2	-15	7.79	2.98
-9	2	15	6.53	2.67
8	2	-15	4.50	2.20
-8	2	15	6.72	2.47
8	-2	-15	2.05	2.36
-7	2	15	-0.85	1.65
7	2	-15	-0.46	1.41
7	-2	-15	-2.06	1.89
6	-2	-15	-0.18	1.67
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5	-2	-15	3.78	1.98
-4	2	15	28.30	2.98
4	-2	-15	24.24	3.21
4	2	-15	22.86	2.92
3	-2	-15	-0.83	0.98
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-2	2	15	1.11	1.61
2	2	-15	0.02	1.34
2	-2	-15	0.51	1.66
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-1	2	15	3.68	2.01
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0	2	-15	5.11	1.91
0	2	15	3.44	2.18
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0	2	-15	8.87	1.96
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4	2	15	1.21	2.46
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-9	3	15	1.40	2.37
9	3	-15	3.83	2.41
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8	-3	-15	1.67	2.39
8	3	-15	3.69	2.22
7	-3	-15	-2.52	1.87
7	3	-15	-2.04	1.42
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3	3	-15	1.05	1.65
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1	3	-15	2.90	1.83
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-9	4	15	1.29	2.48
8	-4	-15	0.66	2.42
-8	4	15	0.10	2.24
8	4	-15	3.56	2.19
-7	4	15	-1.95	1.88
7	-4	-15	-0.57	2.06
7	4	-15	0.15	1.77
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6	5	-15	1.05	1.93
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3	5	-15	2.48	1.89
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6	6	-15	1.62	2.12
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10	0	-16	1.24	2.86
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7	0	-16	-0.34	1.66
6	0	-16	0.40	1.69
5	0	-16	1.43	1.68
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-1	0	-16	-1.48	1.74
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5	6	-16	0.53	1.96
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4	6	-16	-0.25	1.90
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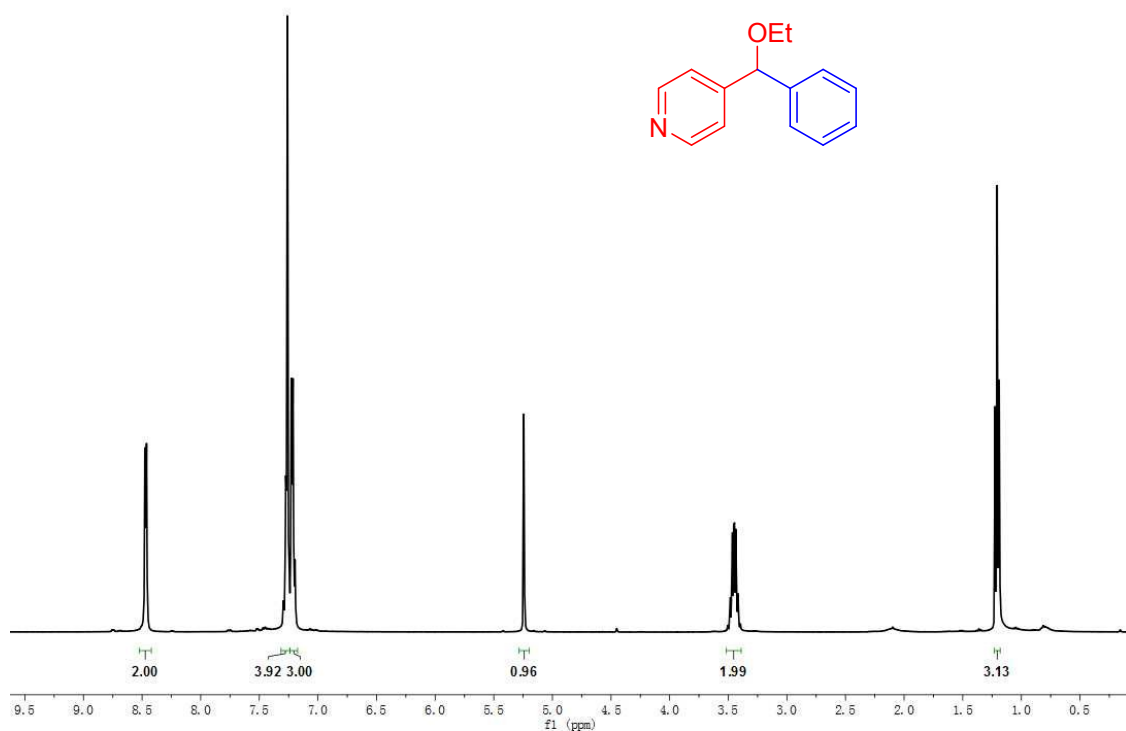
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9. References.

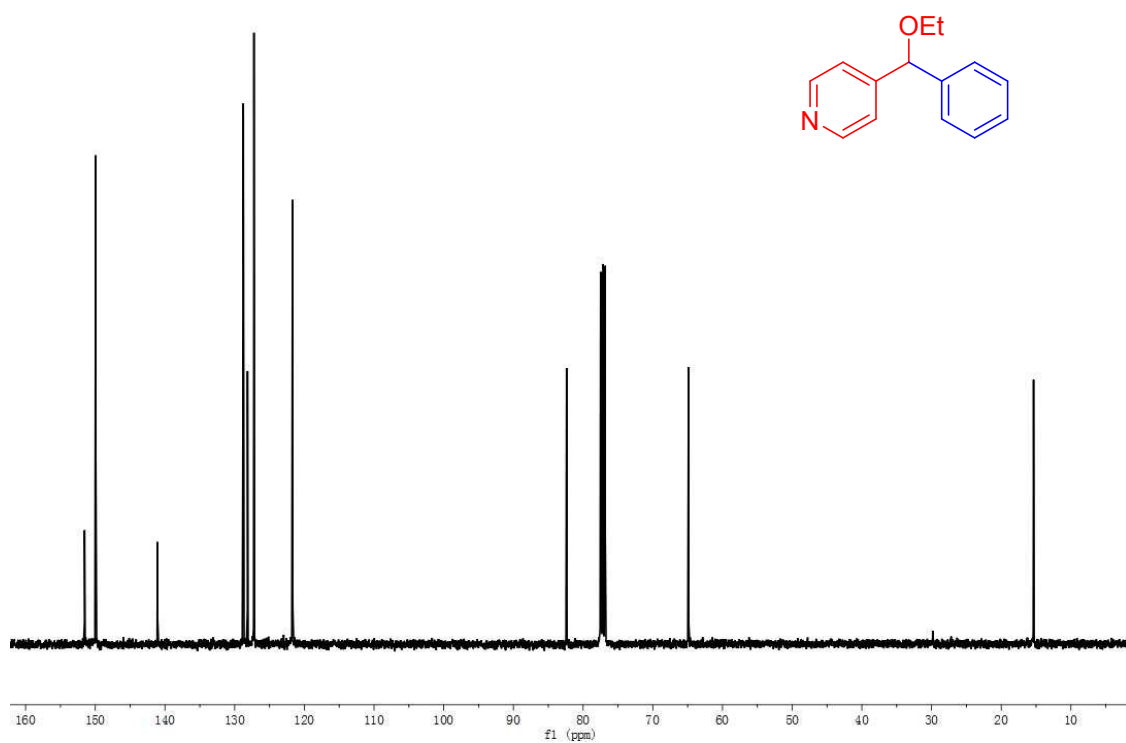
1. F. Gao, B.-S. Kim, P. J. Walsh, *Chem. Sci.*, **2016**, 7, 976-983.
2. J. Yang , G. B. Dudley, *J. Org. Chem.*, **2009**, 74, 7998.
3. A. R. Rivero, B.-S. Kim, P. J. Walsh, *Org. Lett.*, **2016**, 18, 1590-1593.

10. NMR Spectra.

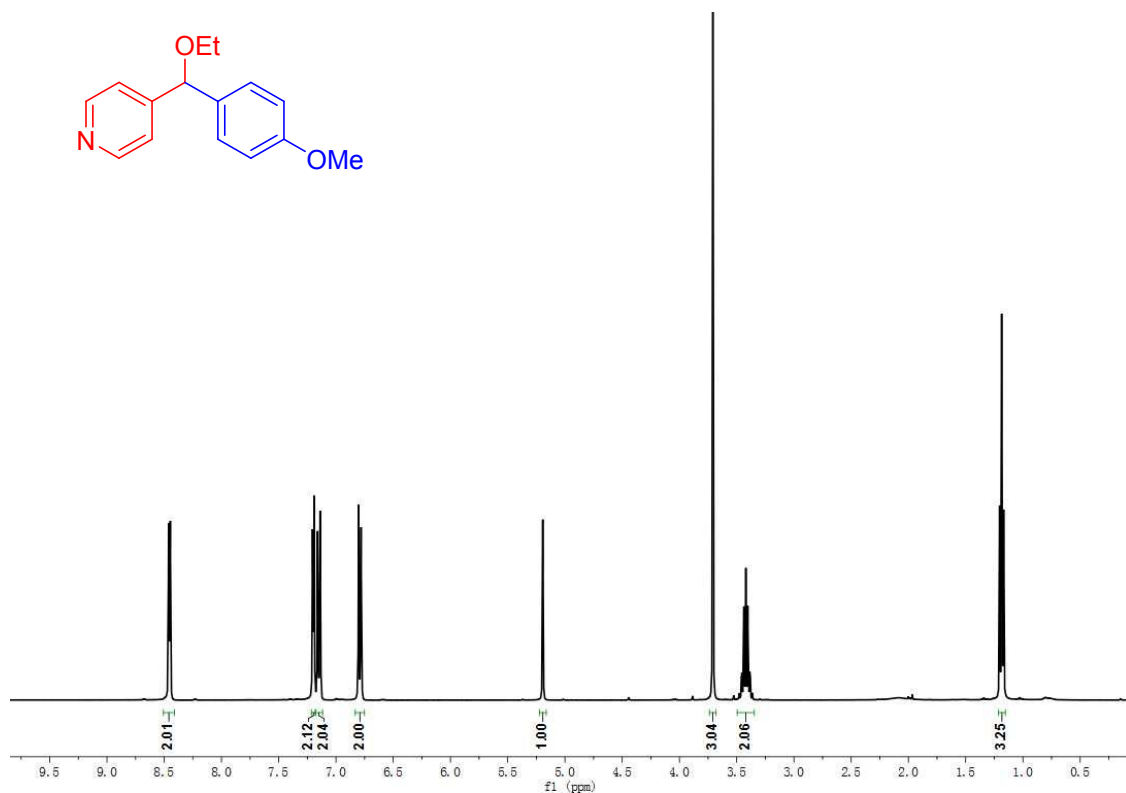
¹H NMR spectra (400 MHz, Chloroform-*d*) of 4-(Ethoxy(phenyl)methyl)pyridine (3aa)



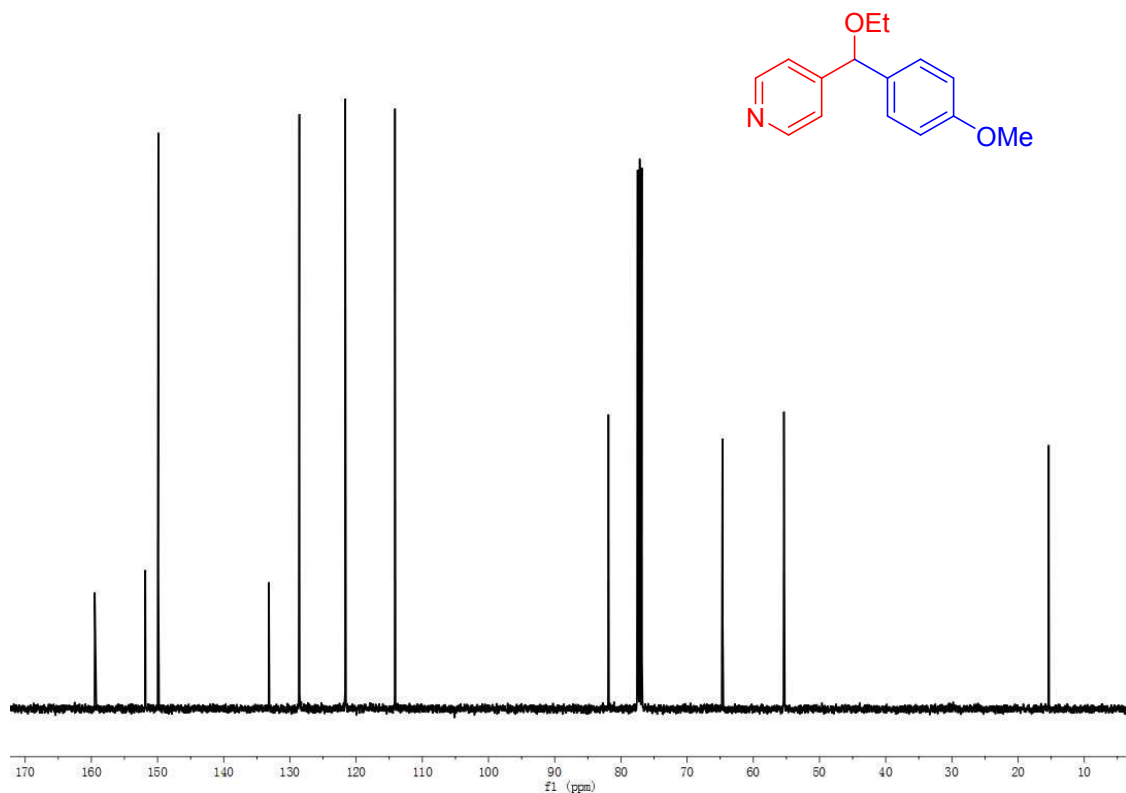
¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 4-(Ethoxy(phenyl)methyl)pyridine (3aa)



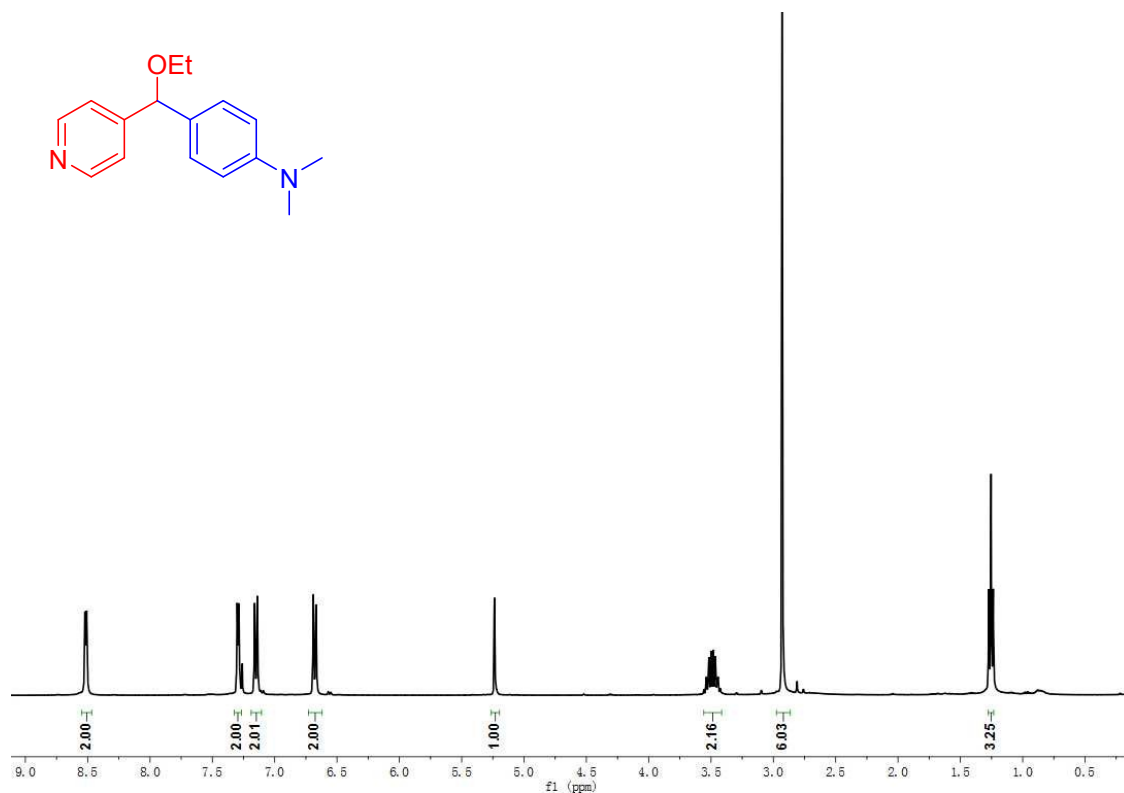
¹H NMR spectra (400 MHz, Chloroform-*d*) of 4-(Ethoxy(4-methoxyphenyl)methyl)pyridine (3ab)



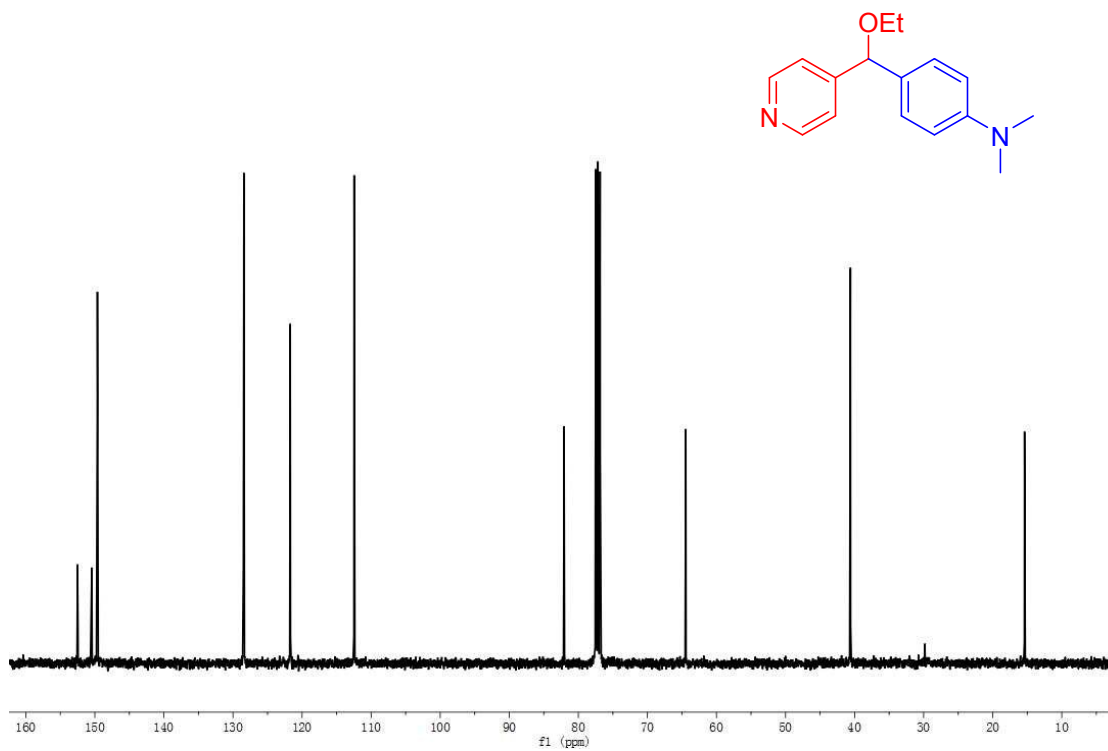
¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 4-(Ethoxy(4-methoxyphenyl)methyl)pyridine (3ab)



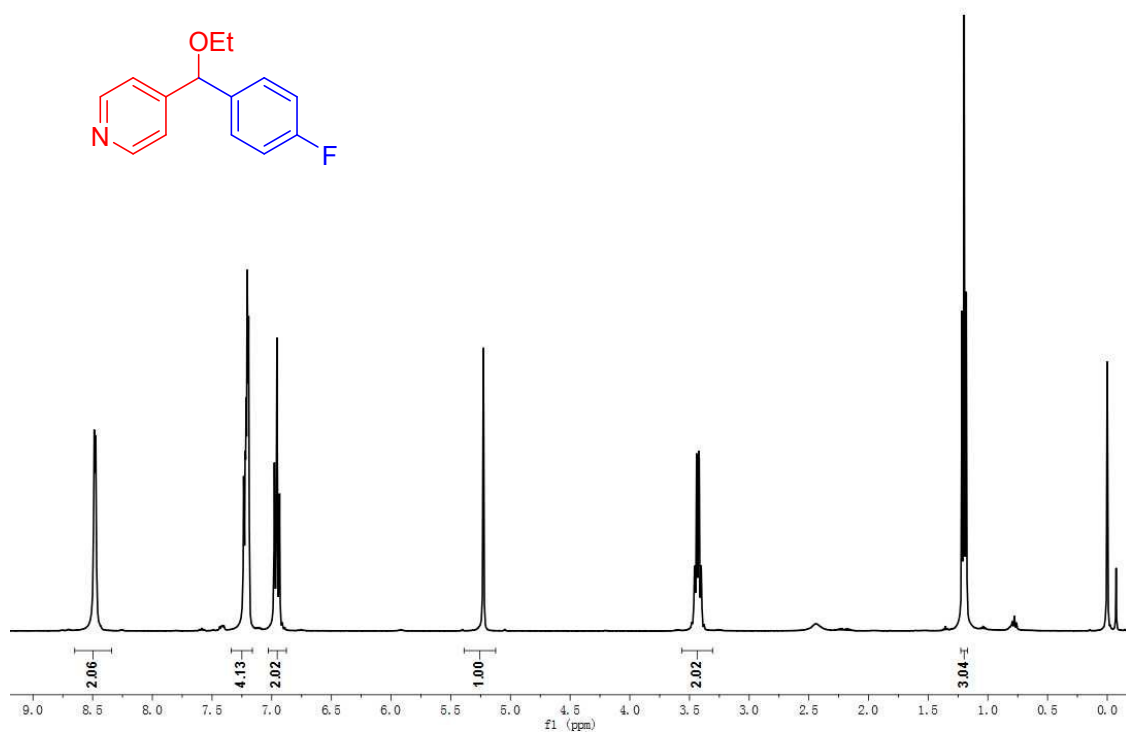
¹H NMR spectra (400 MHz, Chloroform-*d*) of 4-(Ethoxy(pyridin-4-yl)methyl)-N,N-dimethylaniline (3ac)



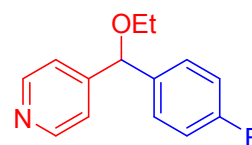
¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 4-(Ethoxy(pyridin-4-yl)methyl)-N,N-dimethylaniline (3ac)

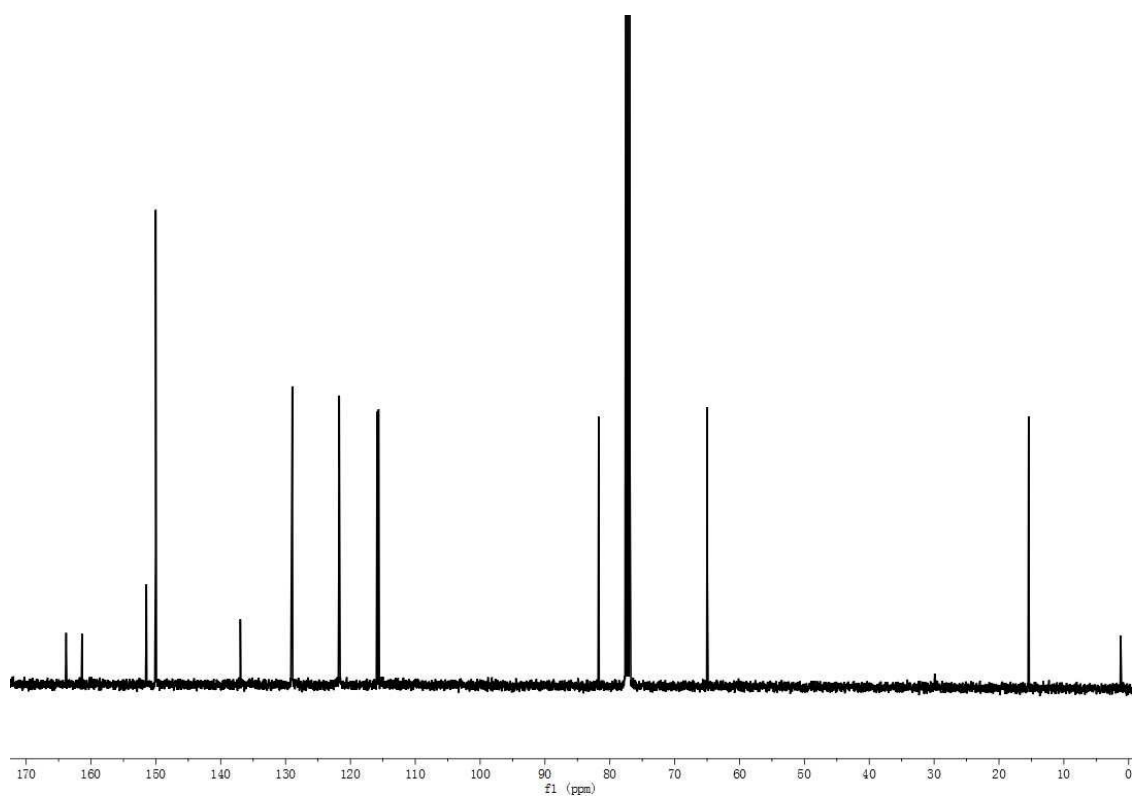


^1H NMR spectra (400 MHz, Chloroform-*d*) of 4-(Ethoxy(4-fluorophenyl)methyl)pyridine (3ad)

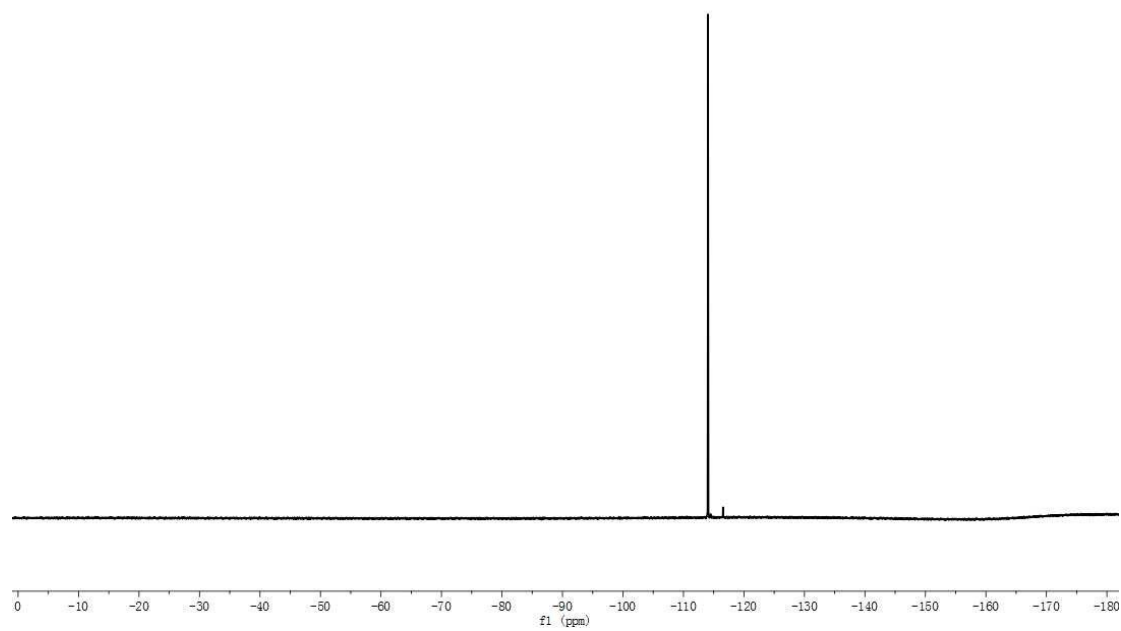


$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 4-(Ethoxy(4-fluorophenyl)methyl)pyridine (3ad)

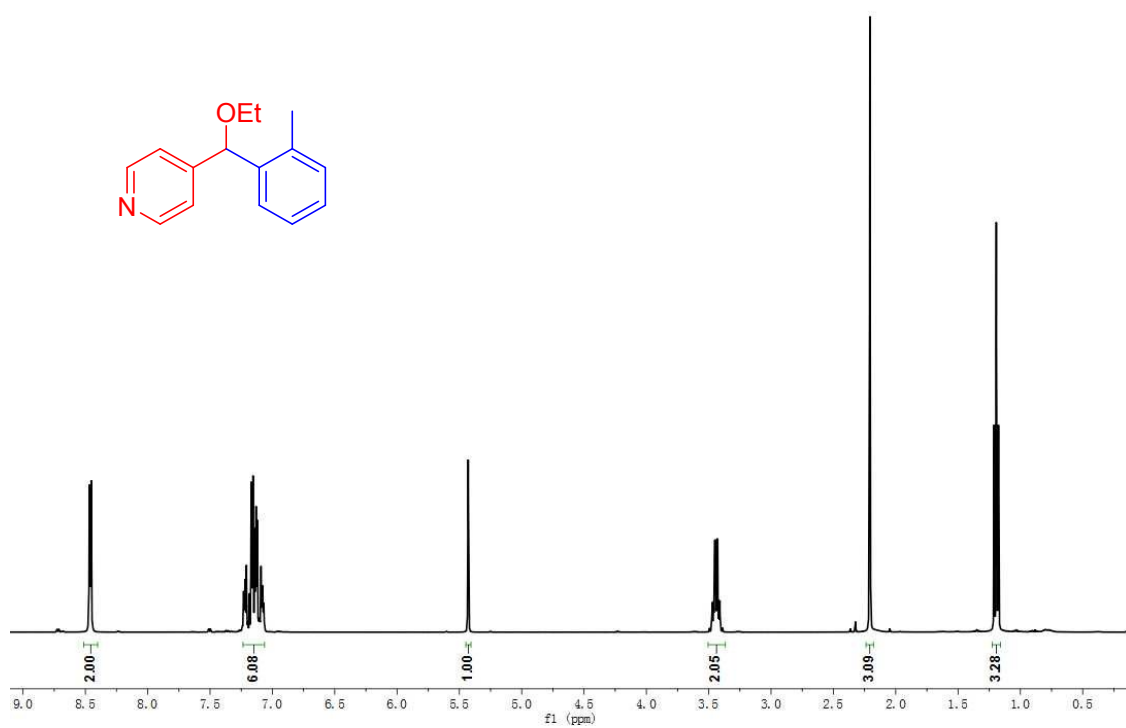




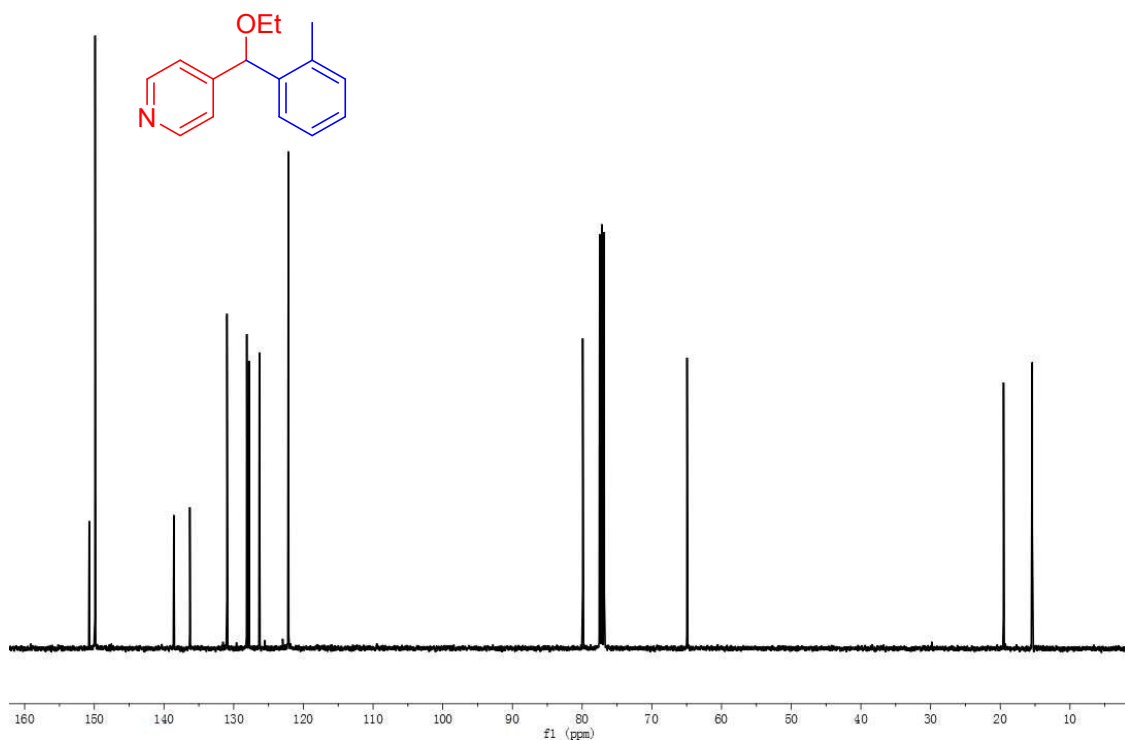
^{19}F NMR spectra (376 MHz, Chloroform-*d*) of 4-(Ethoxy(4-fluorophenyl)methyl)pyridine (3ad)



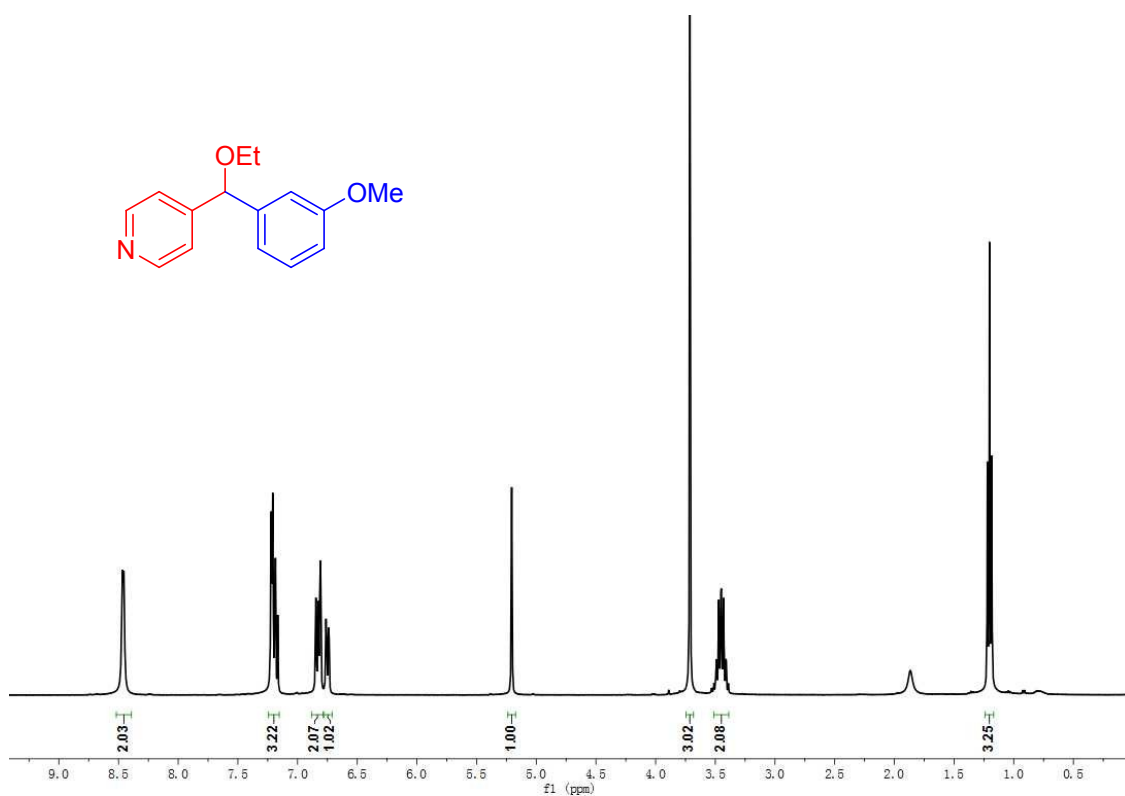
¹H NMR spectra (400 MHz, Chloroform-*d*) of 4-(Ethoxy(*o*-tolyl)methyl)pyridine (3ae)



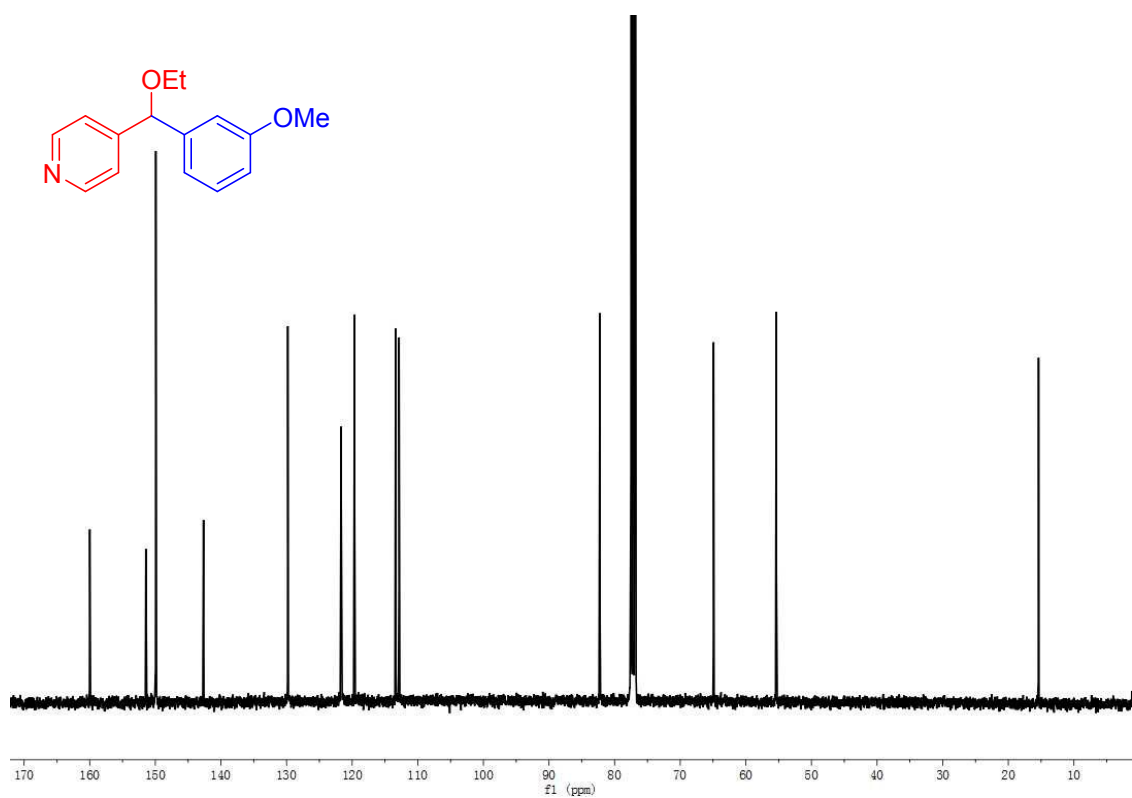
¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 4-(Ethoxy(*o*-tolyl)methyl)pyridine (3ae)



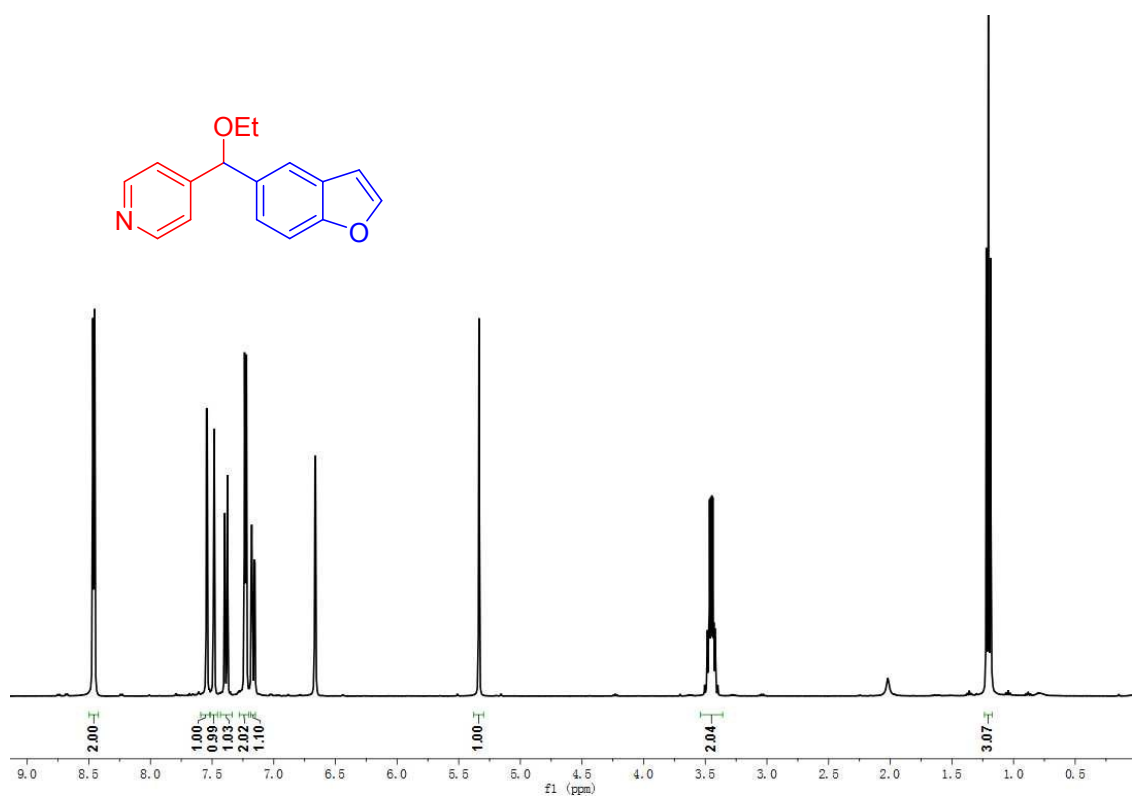
¹H NMR spectra (400 MHz, Chloroform-*d*) of 4-(Ethoxy(3-methoxyphenyl)methyl)pyridine (3af)



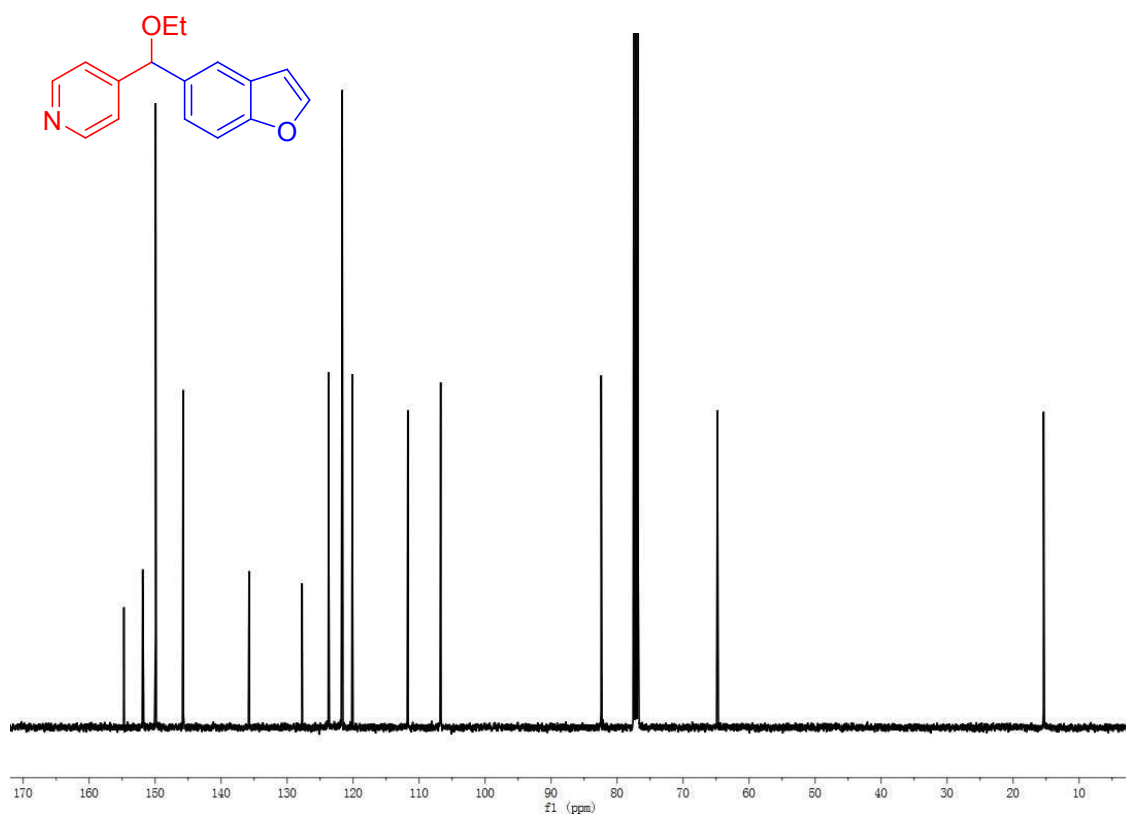
¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 4-(Ethoxy(3-methoxyphenyl)methyl)pyridine (3af)



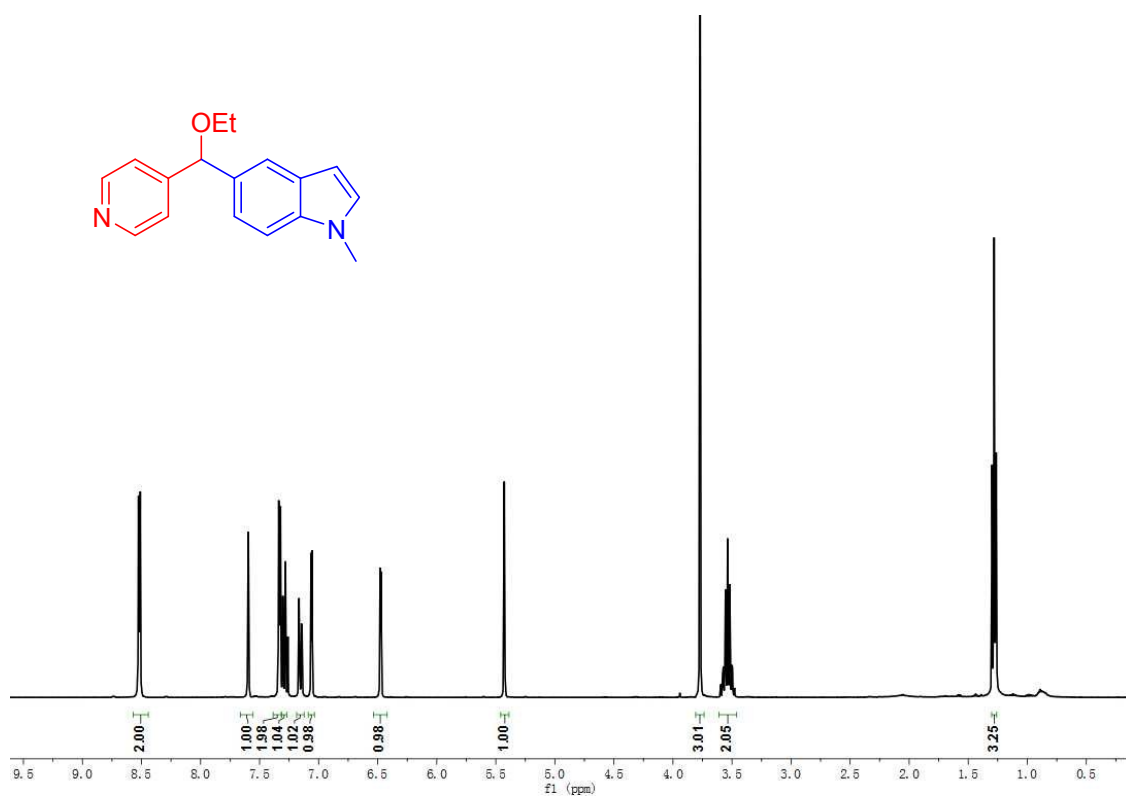
¹H NMR spectra (400 MHz, Chloroform-*d*) of 4-(Benzofuran-5-yl(ethoxy)methyl)pyridine (3ag)



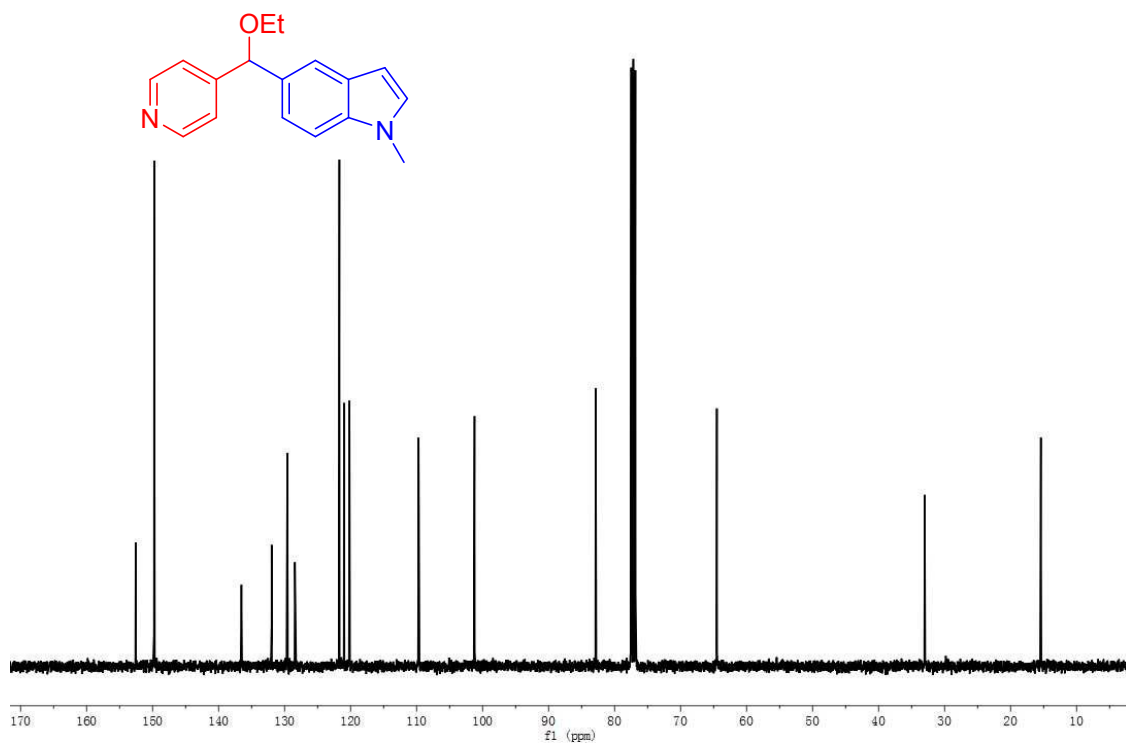
¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 4-(Benzofuran-5-yl(ethoxy)methyl)pyridine (3ag)



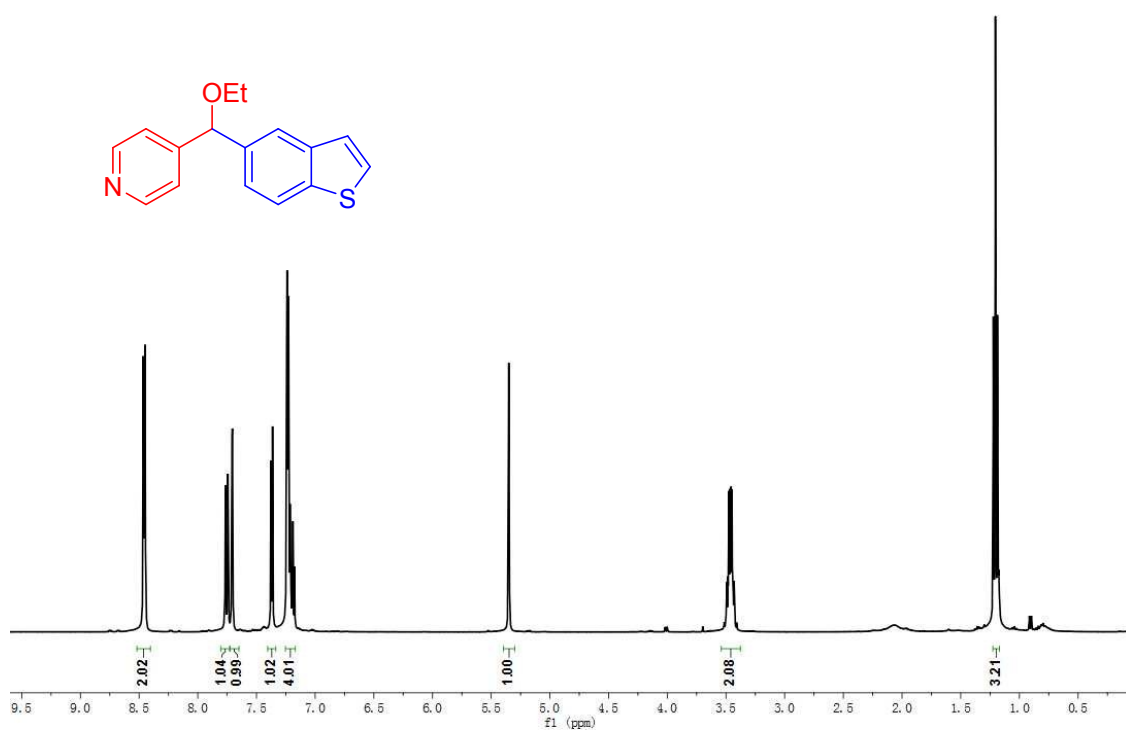
¹H NMR spectra (400 MHz, Chloroform-*d*) of 5-(Ethoxy(pyridin-4-yl)methyl)-1-methyl-1H-indole (3ah)



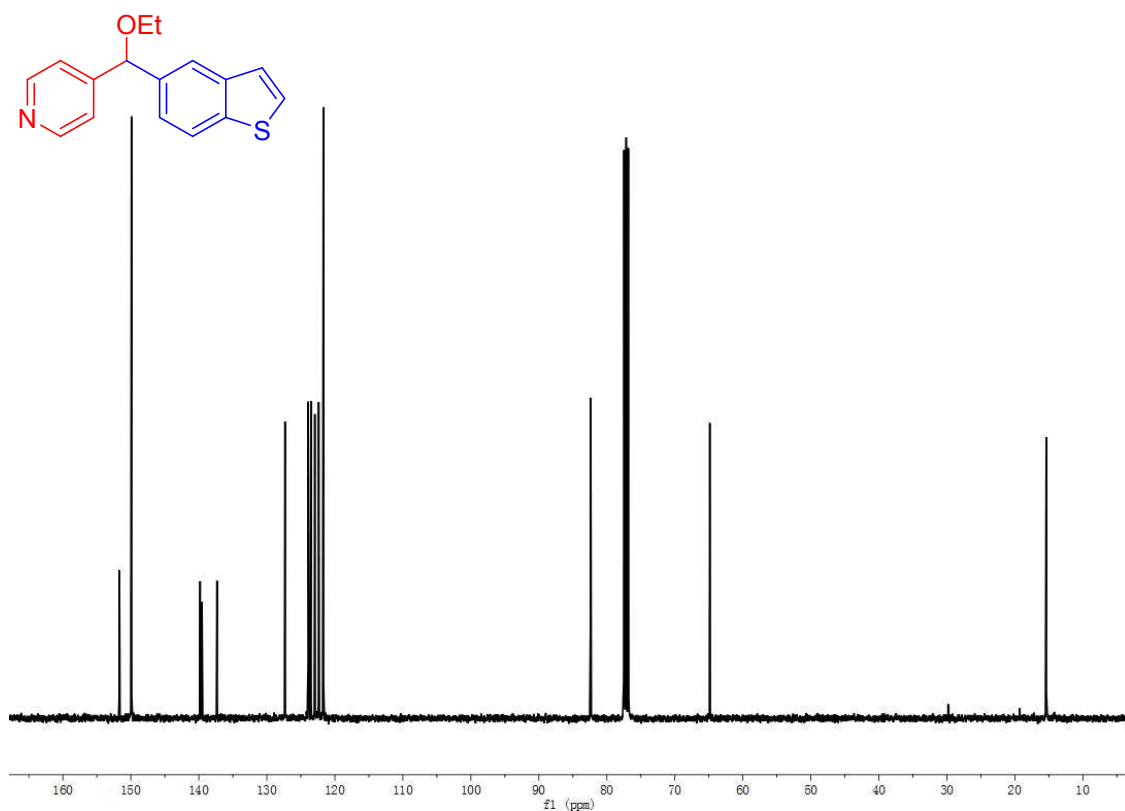
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 5-(Ethoxy(pyridin-4-yl)methyl)-1-methyl-1H-indole (3ah)



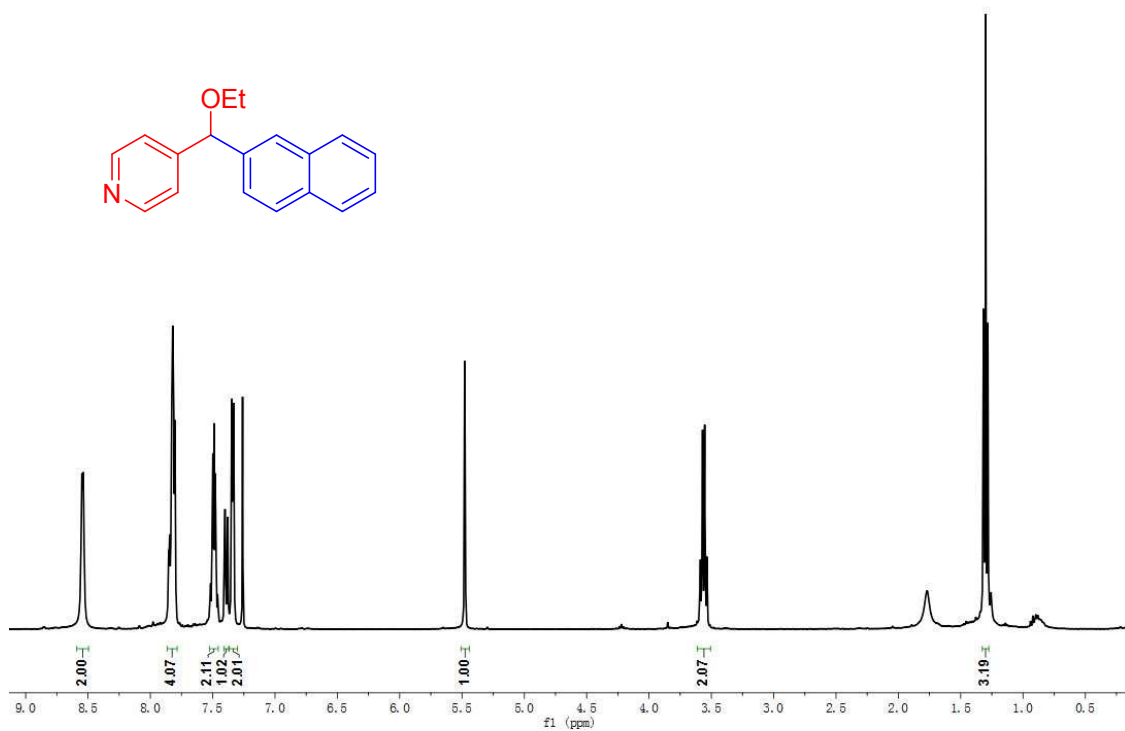
^1H NMR spectra (400 MHz, Chloroform-*d*) of 4-(Benzo[b]thiophen-5-yl(ethoxy)methyl)pyridine (3ai)



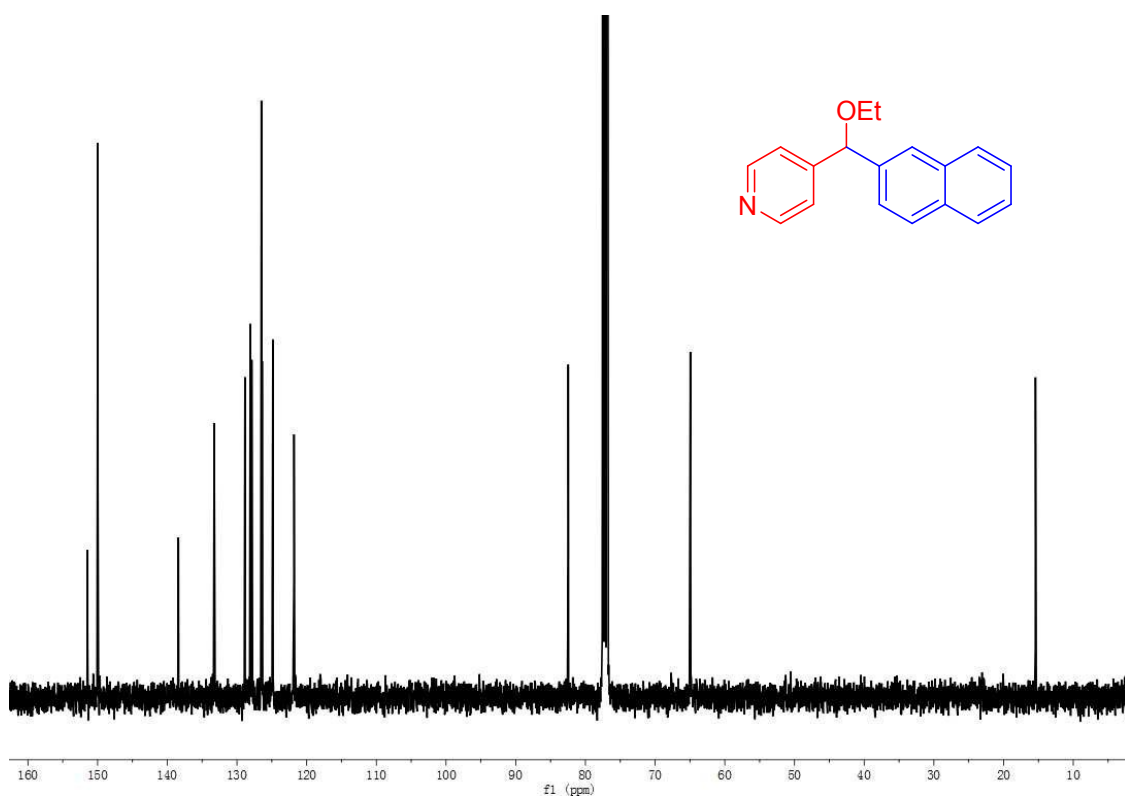
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 4-(Benzo[b]thiophen-5-yl(ethoxy)methyl)pyridine (3ai)



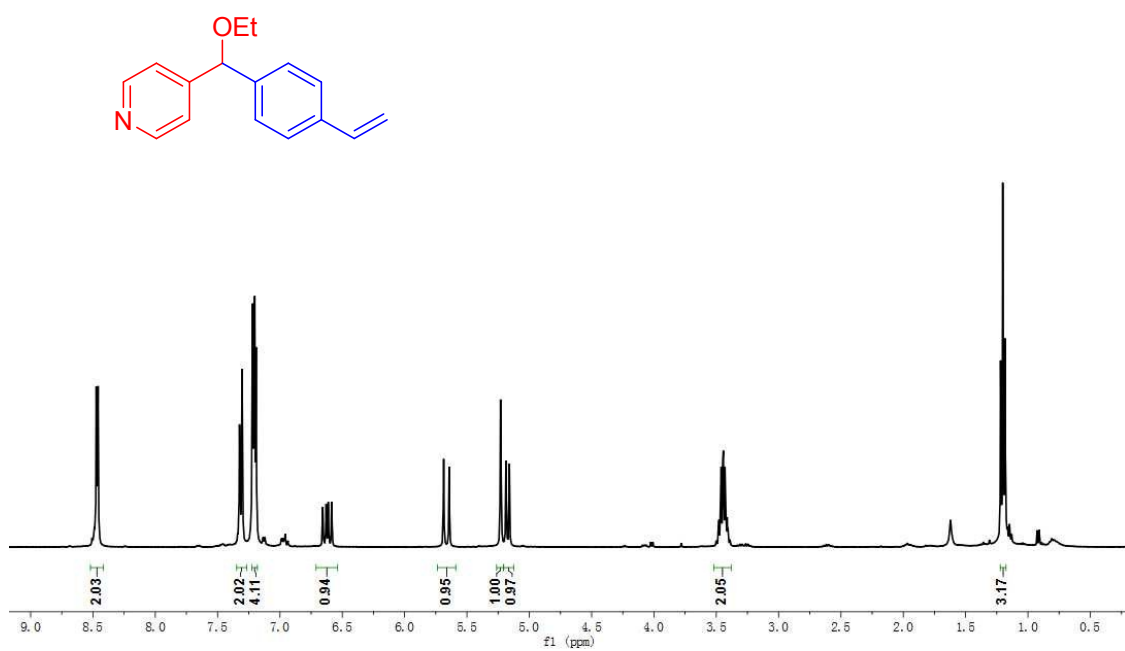
^1H NMR spectra (400 MHz, Chloroform-*d*) of 4-(Ethoxy(naphthalen-2-yl)methyl)pyridine (3aj)



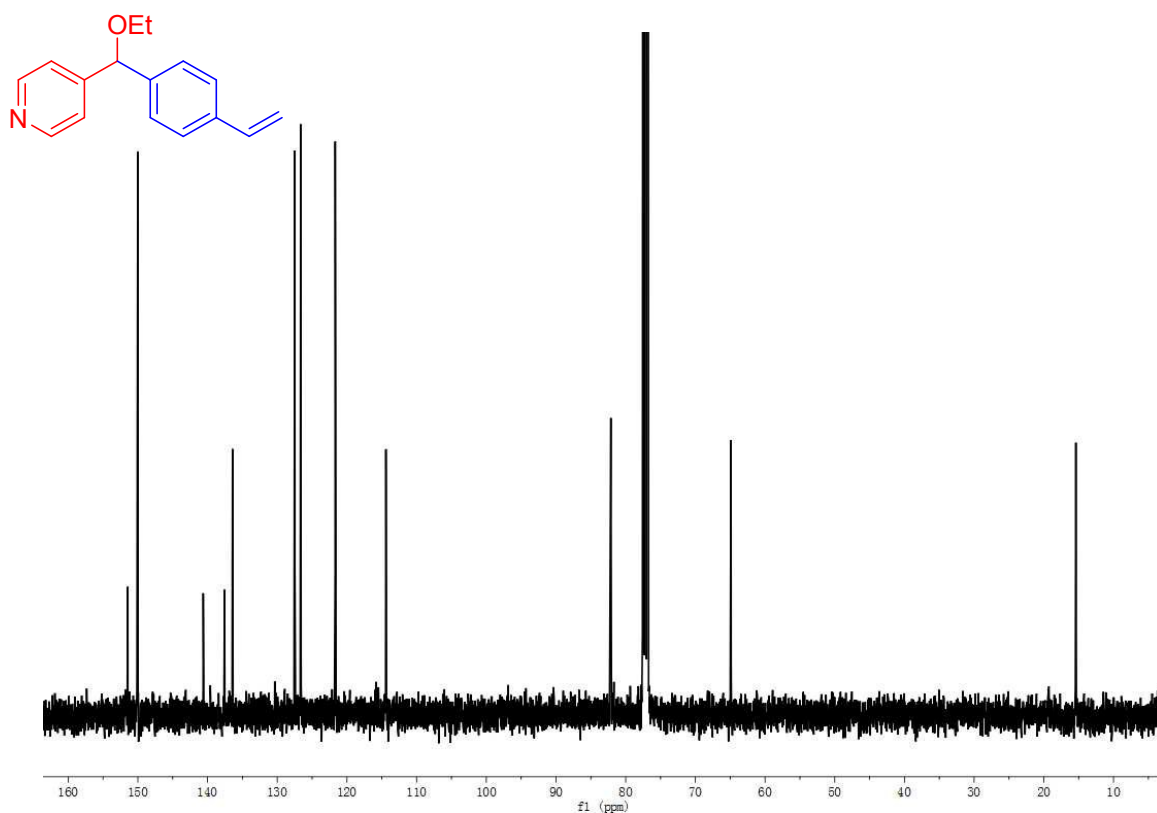
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 4-(Ethoxy(naphthalen-2-yl)methyl)pyridine (3aj)



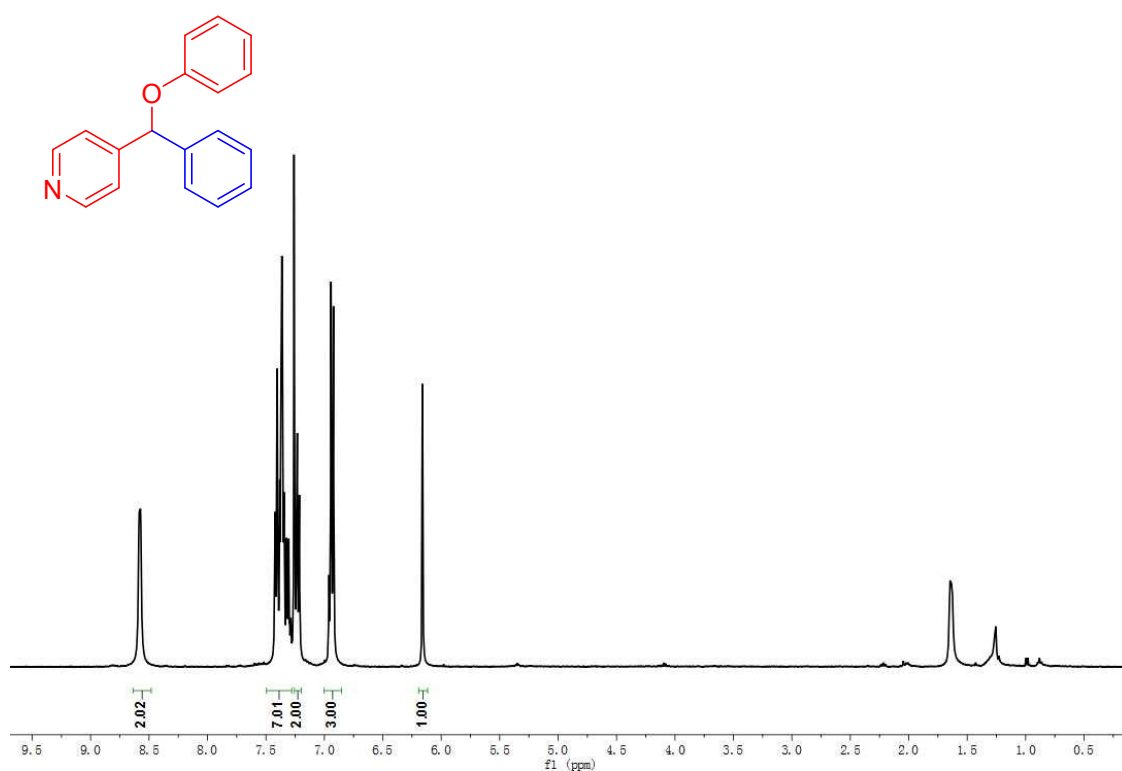
^1H NMR spectra (400 MHz, Chloroform-*d*) of 4-(Ethoxy(4-vinylphenyl)methyl)pyridine (3ak)



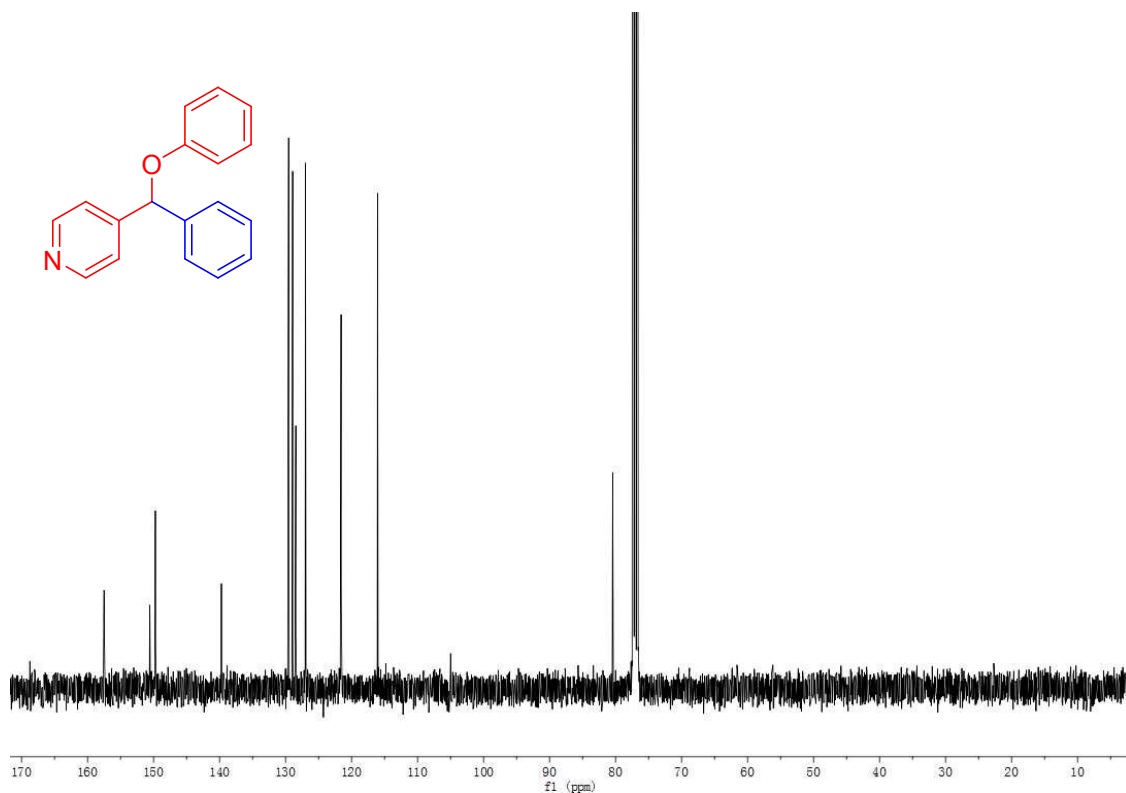
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 4-(Ethoxy(4-vinylphenyl)methyl)pyridine (3ak)



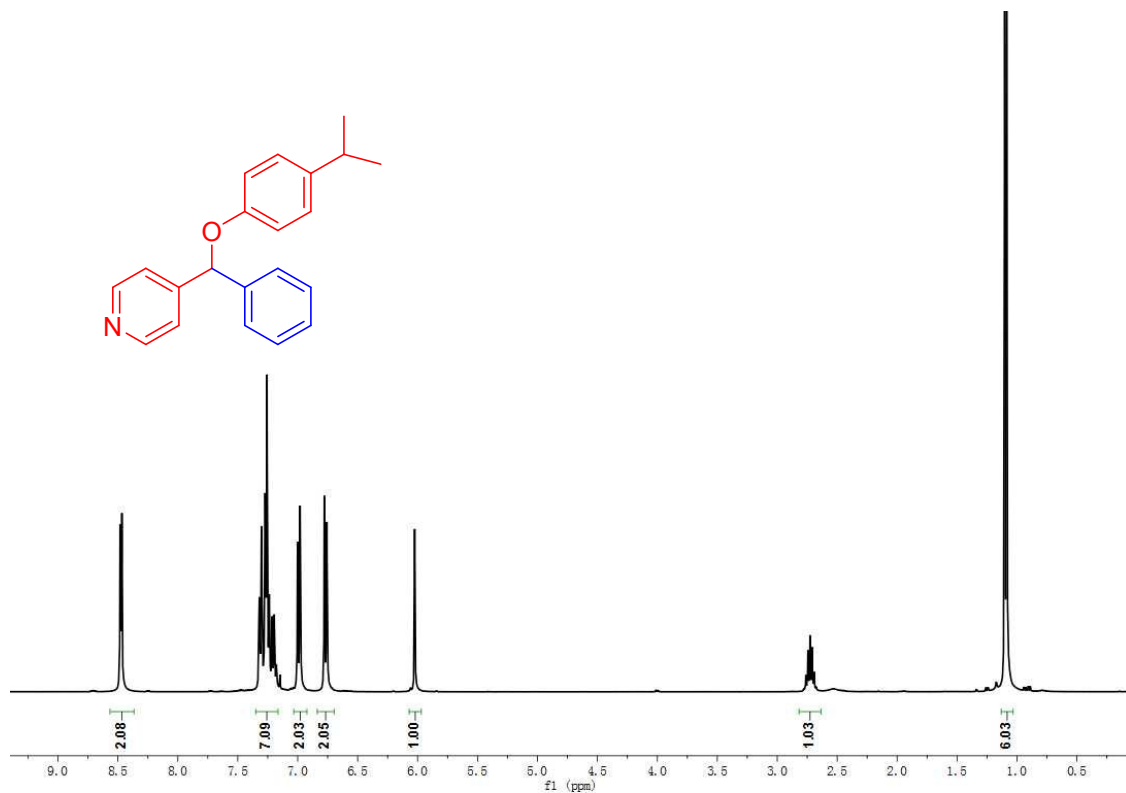
^1H NMR spectra (400 MHz, Chloroform-*d*) of 4-(Phenoxy(phenyl)methyl)pyridine (3ba)



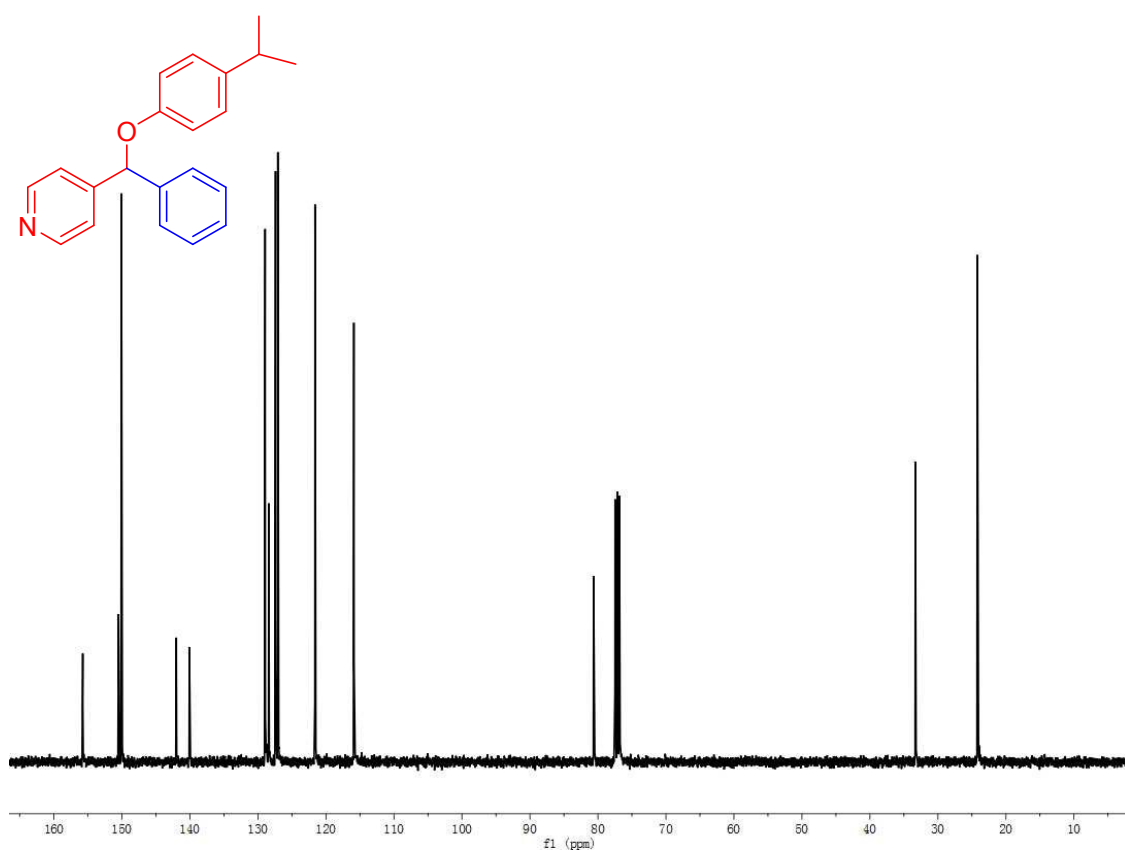
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 4-(Phenoxy(phenyl)methyl)pyridine (3ba)



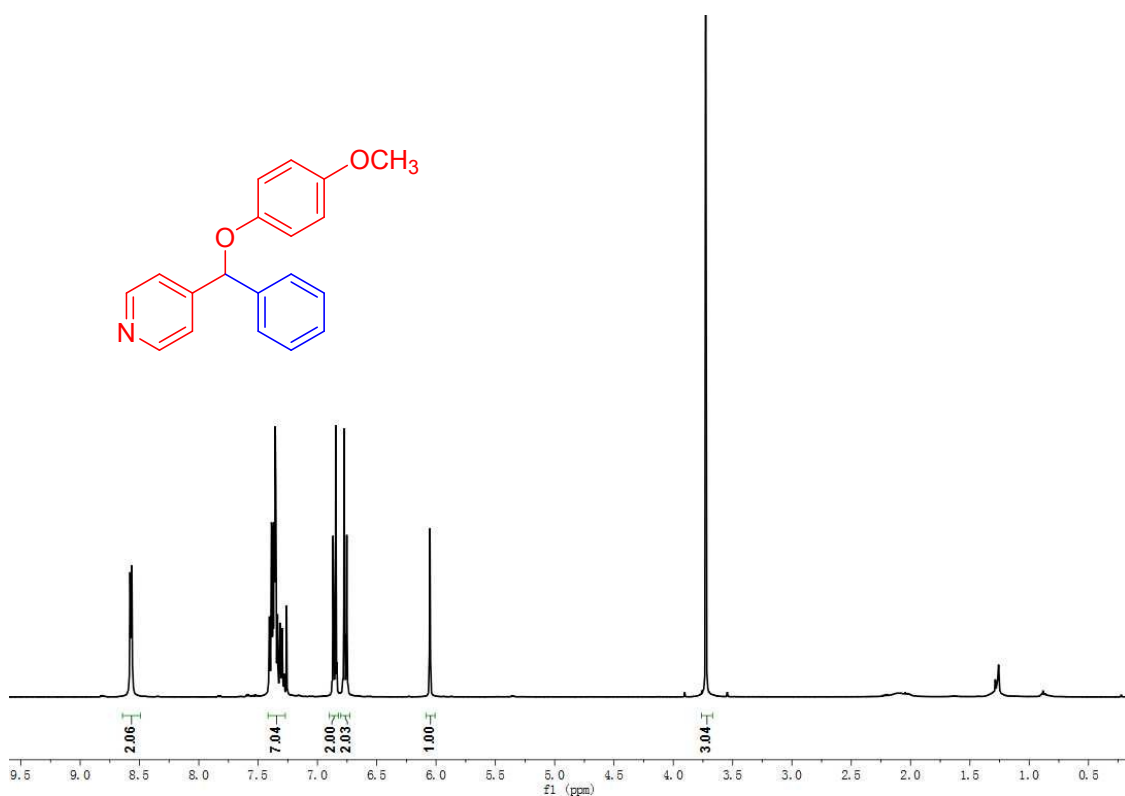
^1H NMR spectra (400 MHz, Chloroform-*d*) of 4-((4-Isopropylphenoxy)(phenyl)methyl)pyridine (3ca)



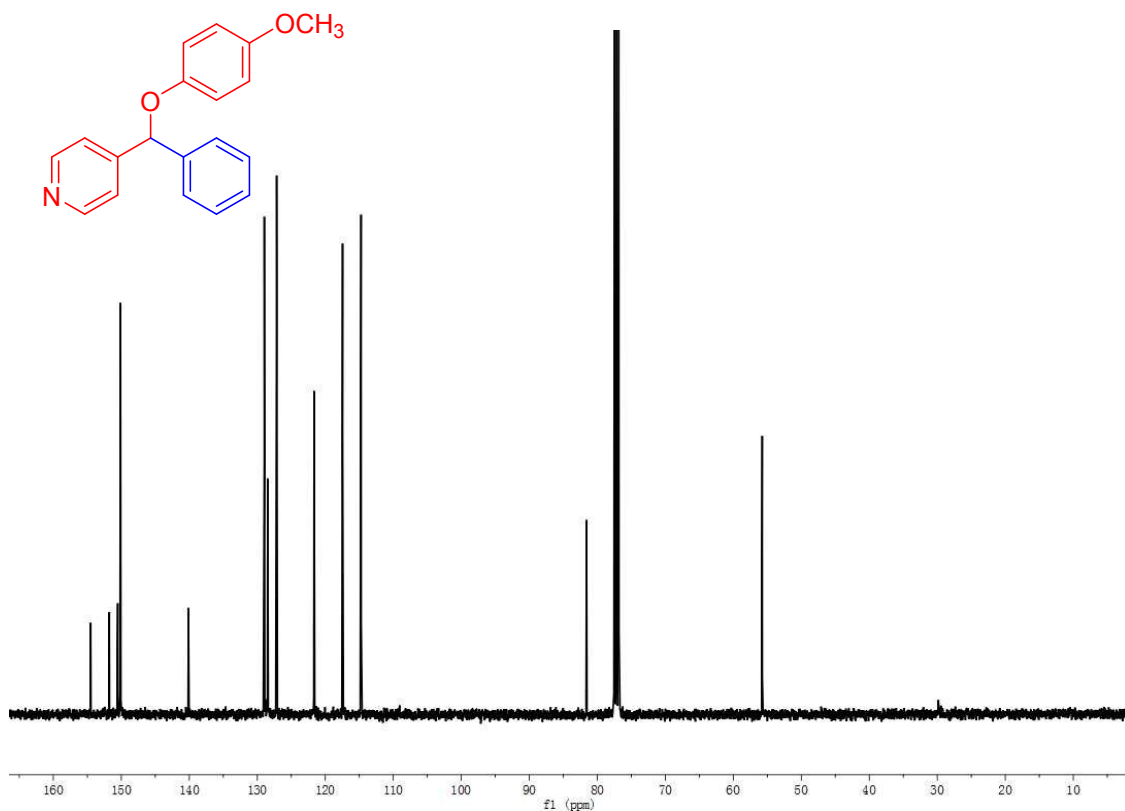
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 4-((4-Isopropylphenoxy)(phenyl)methyl)pyridine (3ca)



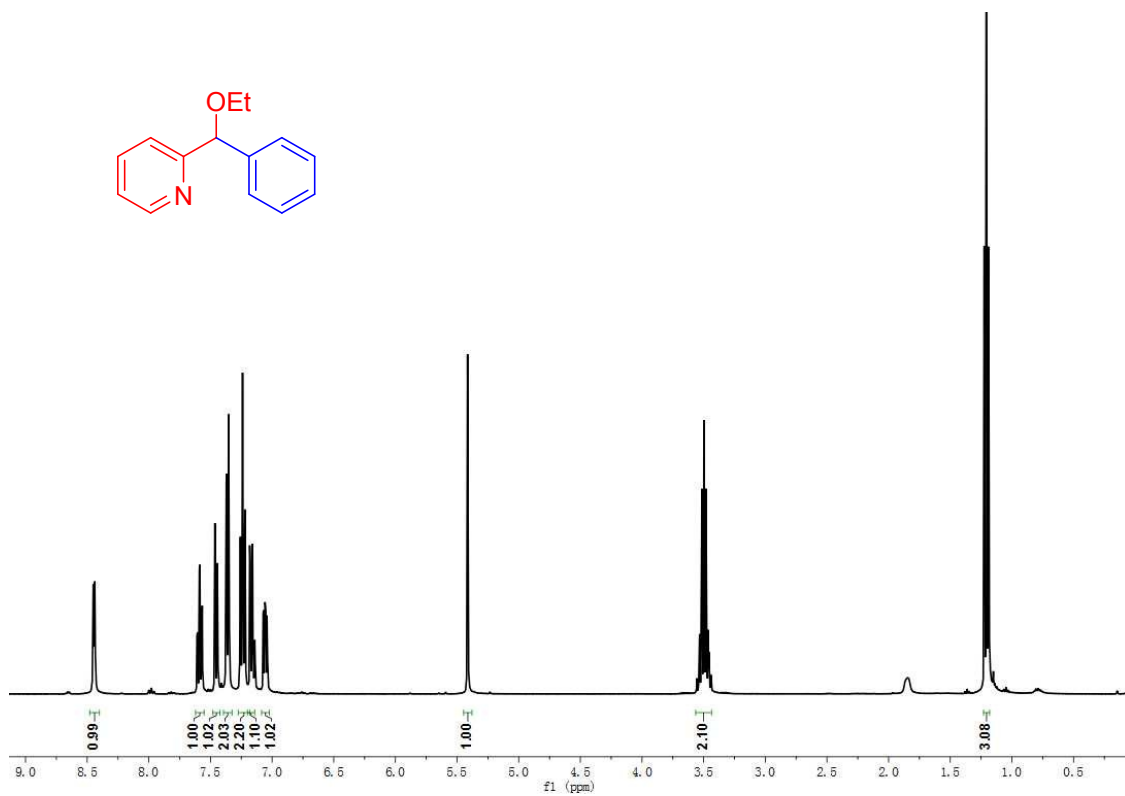
^1H NMR spectra (400 MHz, Chloroform-*d*) of 4-((4-Methoxyphenoxy)(phenyl)methyl)pyridine (3da)



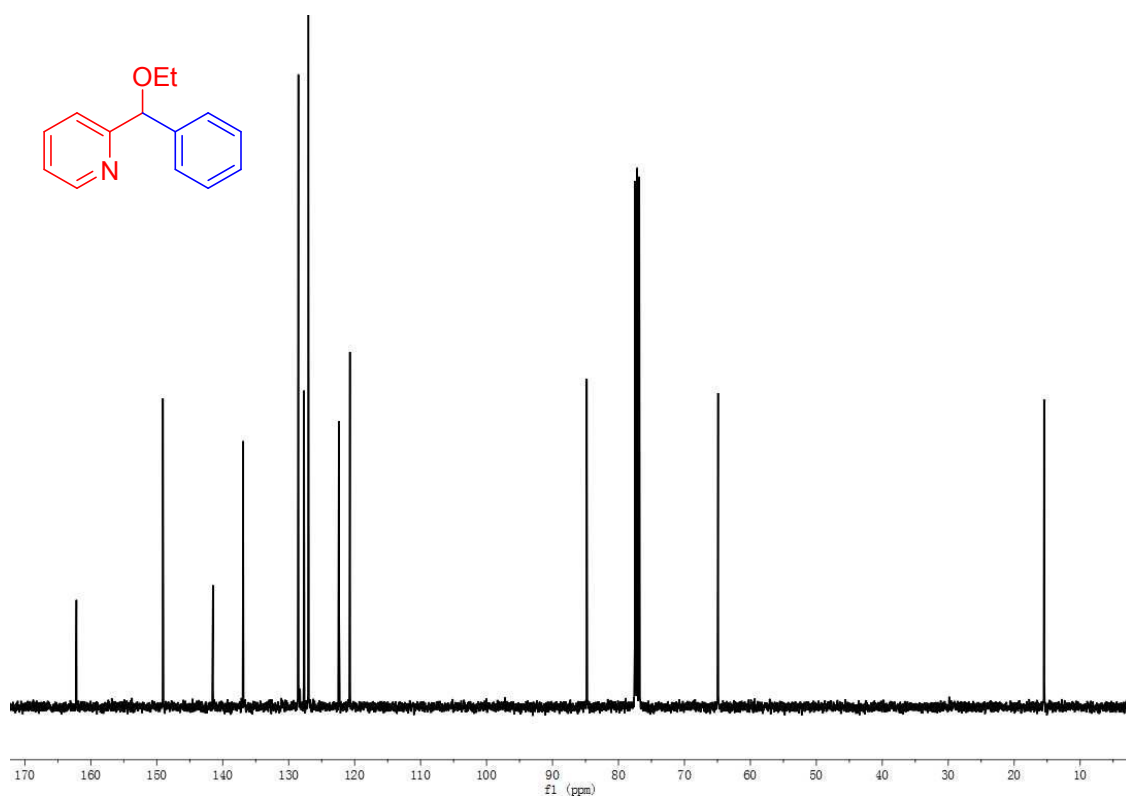
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 4-((4-Methoxyphenoxy)(phenyl)methyl)pyridine (3da)



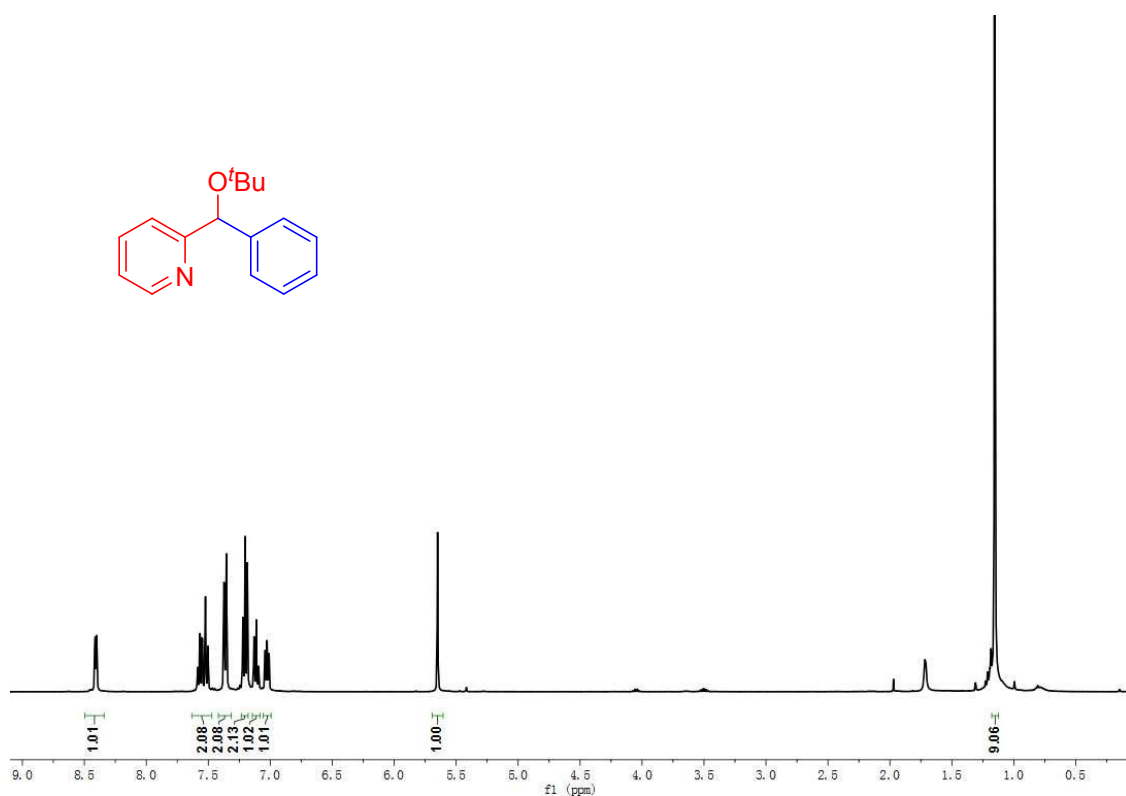
^1H NMR spectra (400 MHz, Chloroform-*d*) of 2-(Ethoxy(phenyl)methyl)pyridine (3ea)



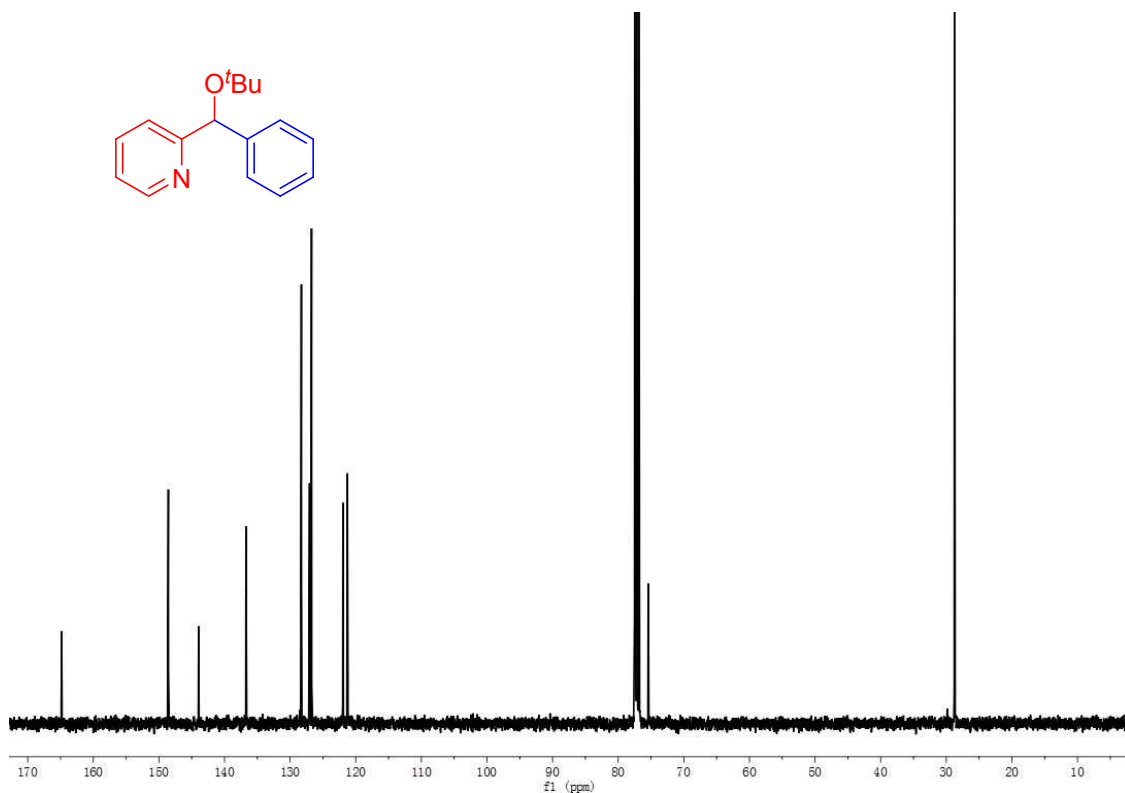
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 2-(Ethoxy(phenyl)methyl)pyridine (3ea)



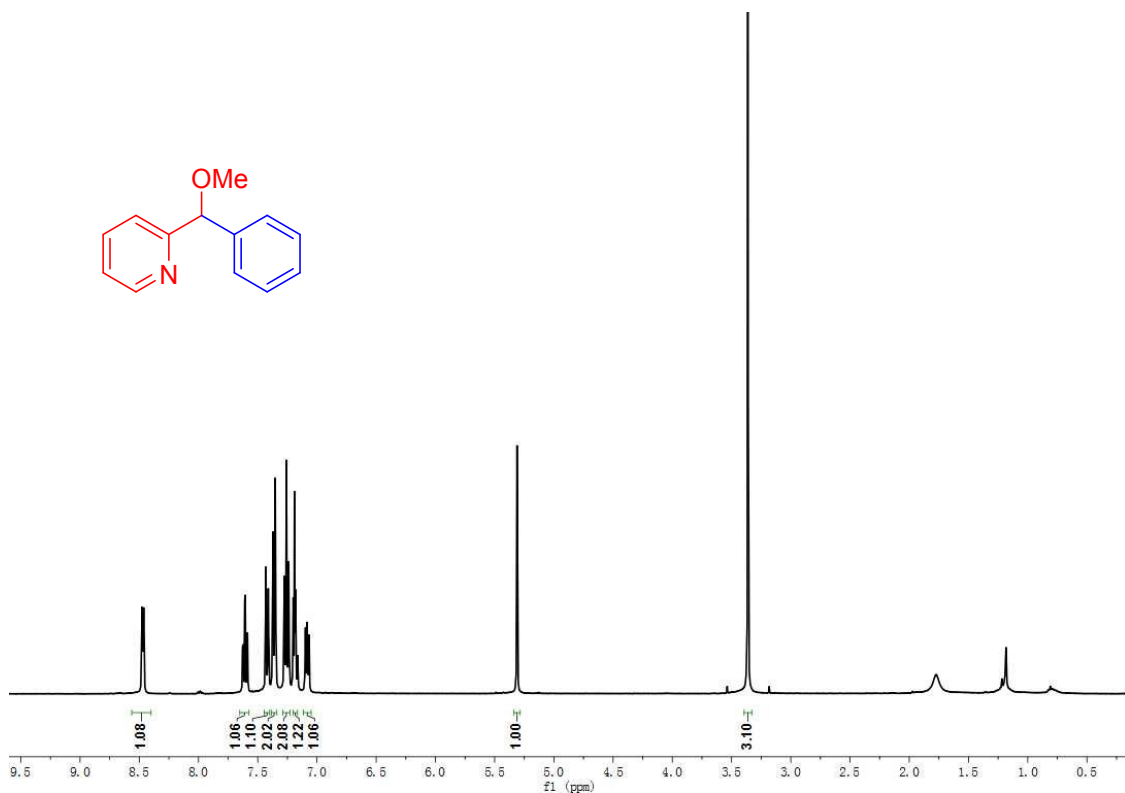
^1H NMR spectra (400 MHz, Chloroform-*d*) of 2-(*tert*-Butoxy(phenyl)methyl)pyridine (3fa)



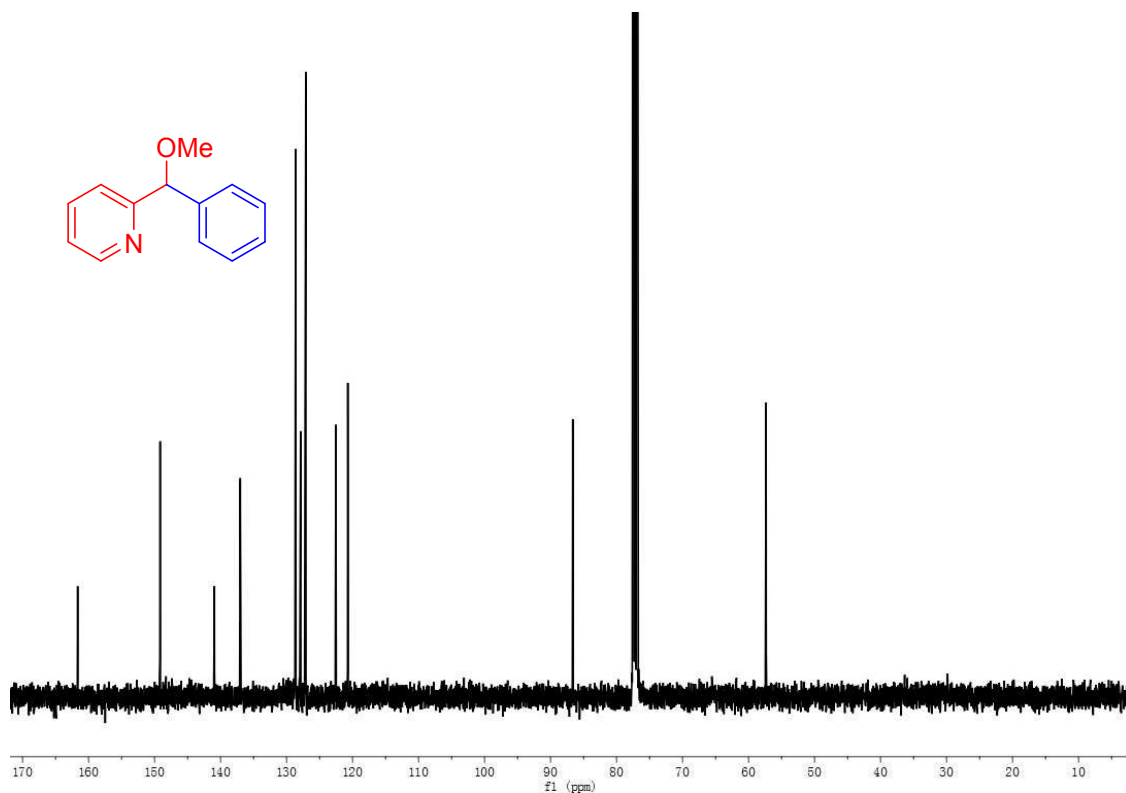
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 2-(*tert*-Butoxy(phenyl)methyl)pyridine (3fa)



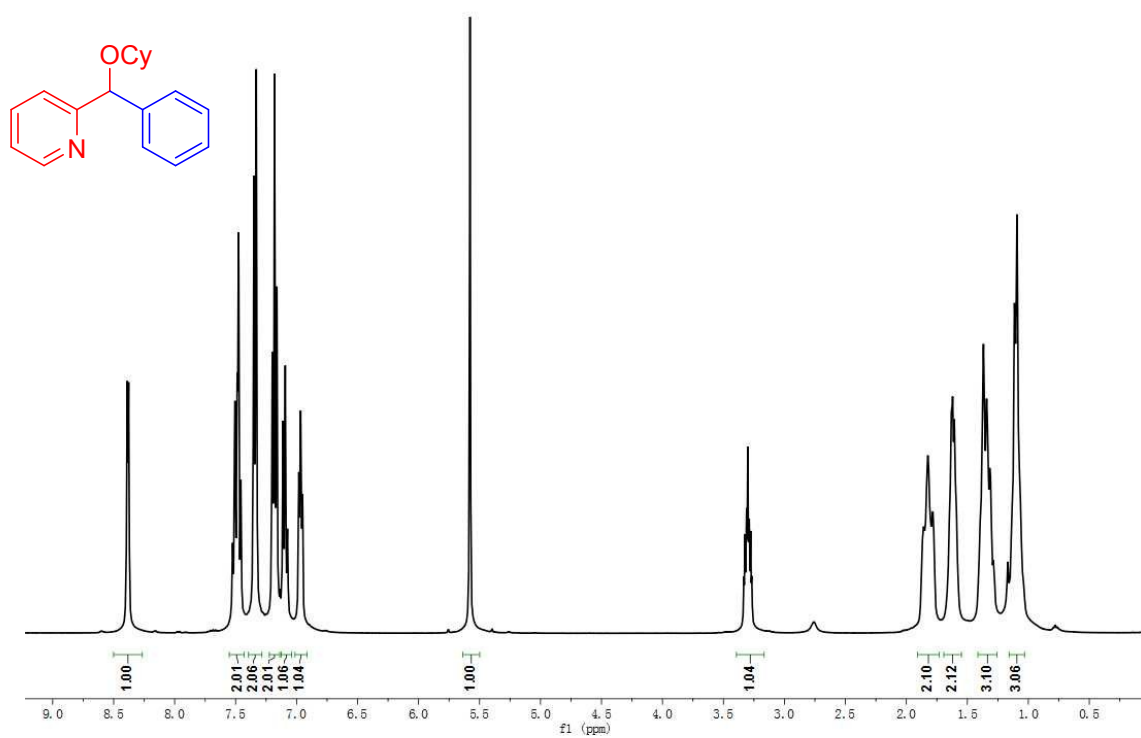
^1H NMR spectra (400 MHz, Chloroform-*d*) of 2-(Methoxy(phenyl)methyl)pyridine (3ga)



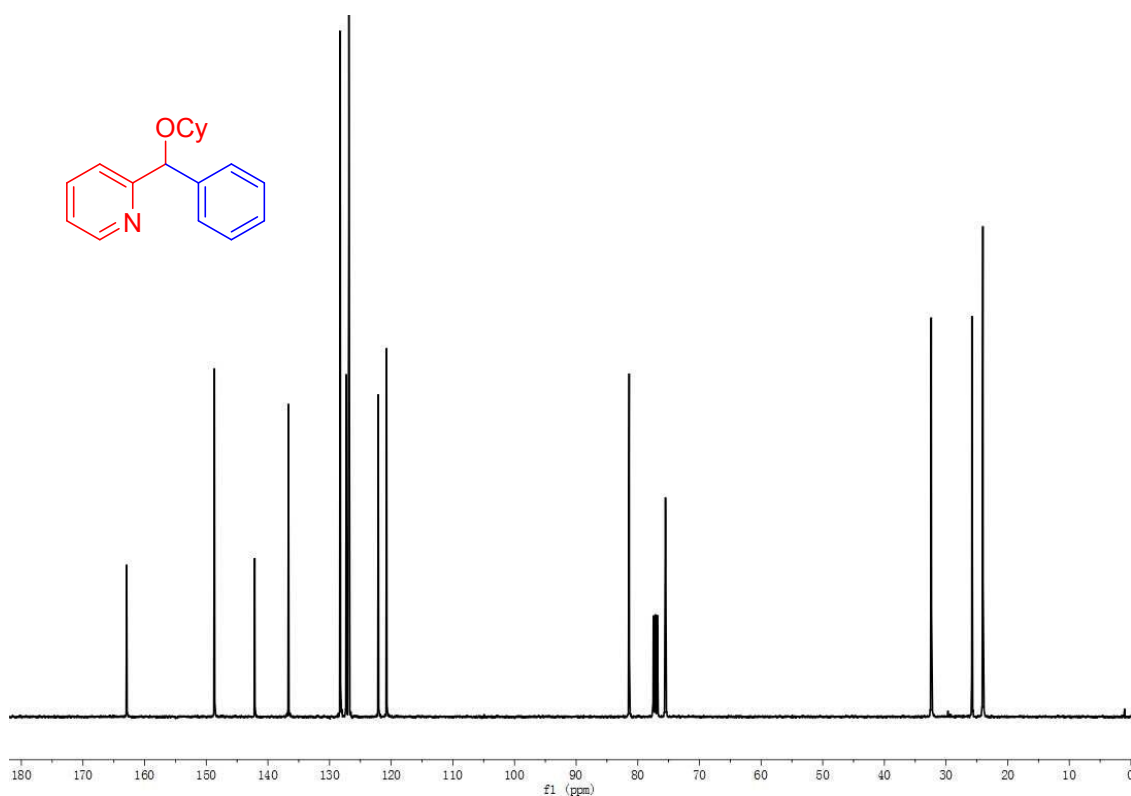
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 2-(Methoxy(phenyl)methyl)pyridine (3ga)



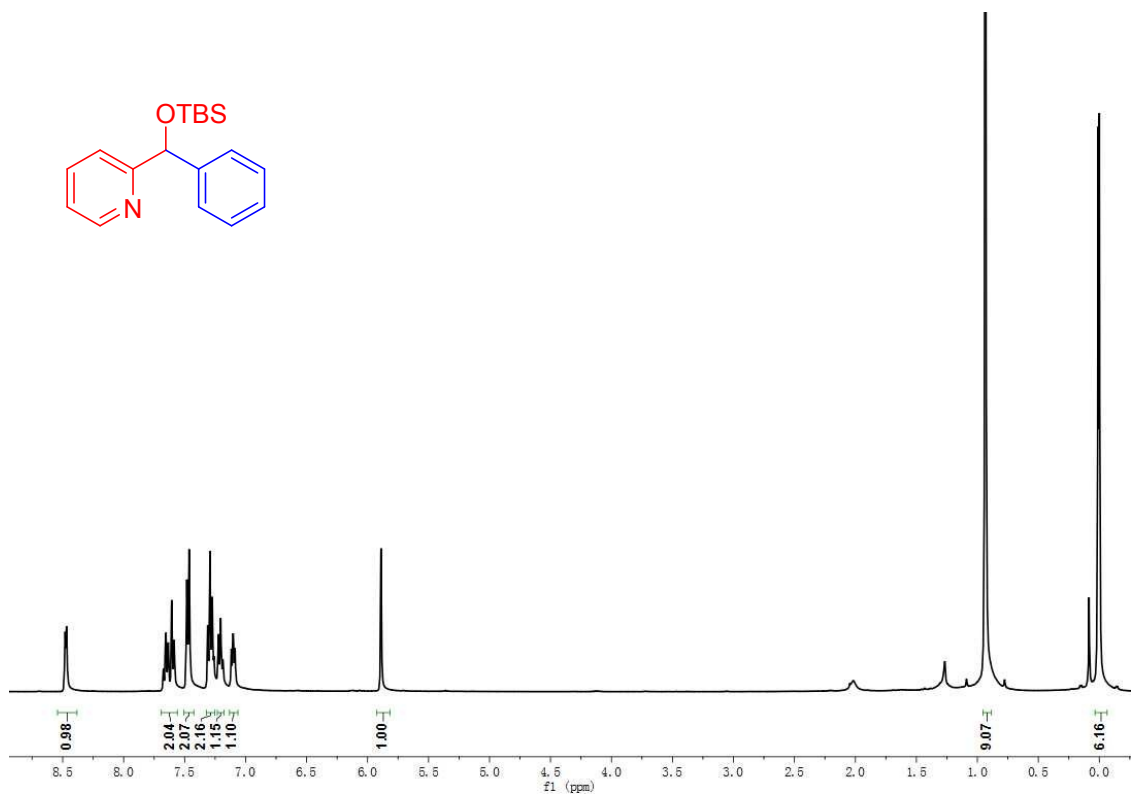
^1H NMR spectra (400 MHz, Chloroform-*d*) of 2-((Cyclohexyloxy)(phenyl)methyl)pyridine (3ha)



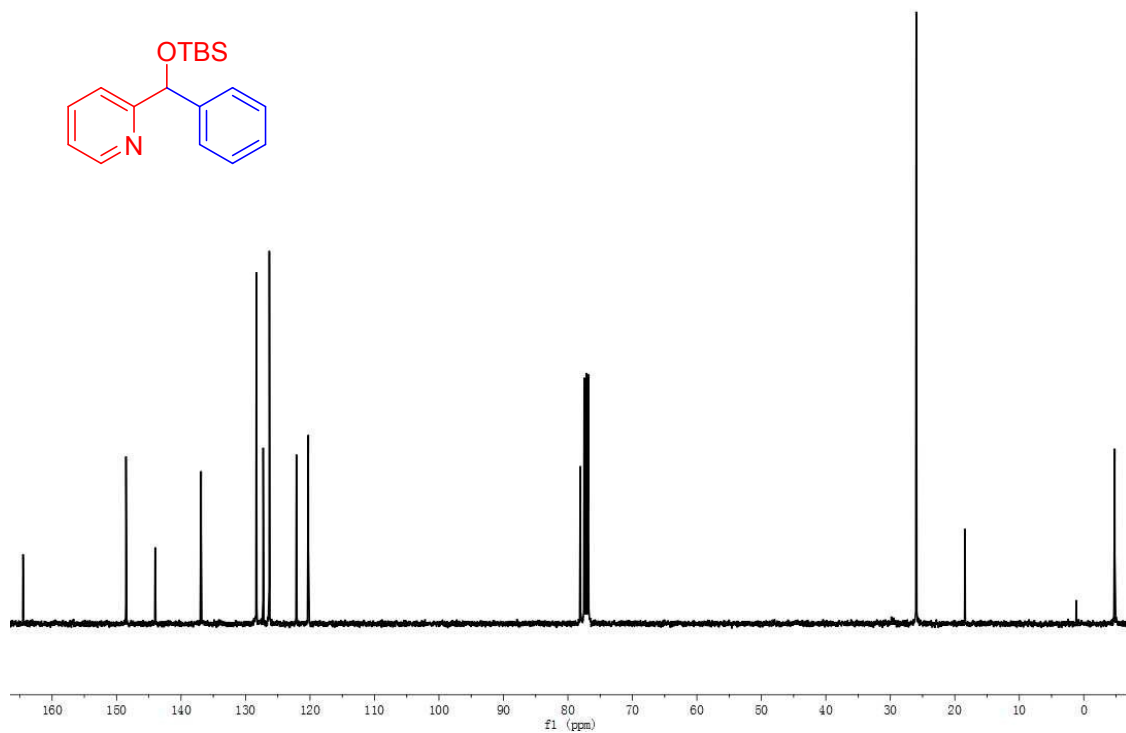
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 2-((Cyclohexyloxy)(phenyl)methyl)pyridine (3ha)



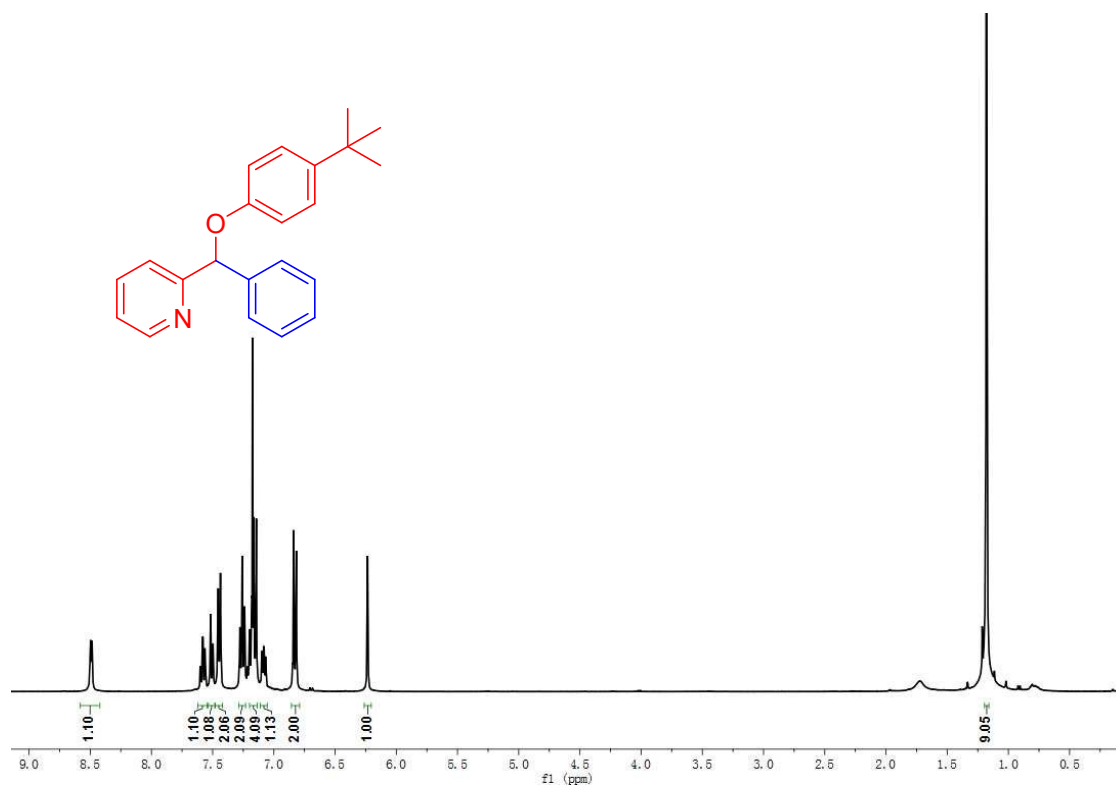
^1H NMR spectra (400 MHz, Chloroform-*d*) of 2-(((*tert*-Butyldimethylsilyl)oxy)(phenyl)methyl)pyridine (3ia)



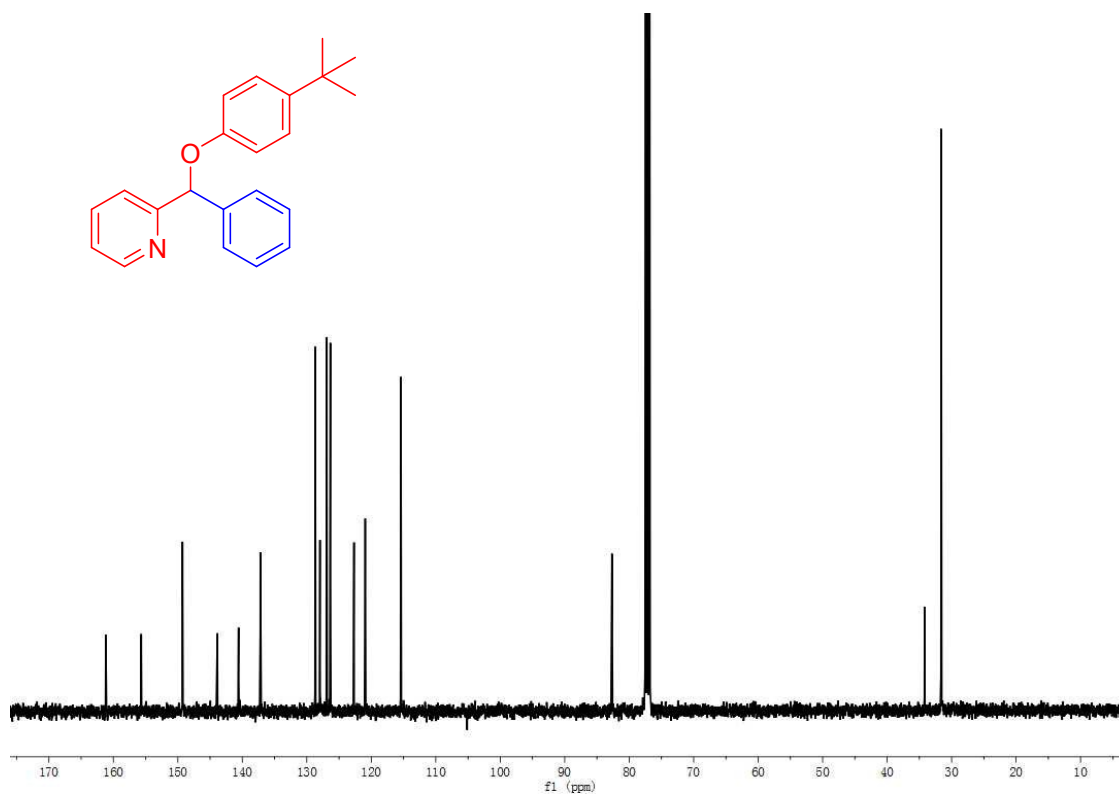
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 2-(((*tert*-Butyldimethylsilyl)oxy)(phenyl)methyl)pyridine (3ia)



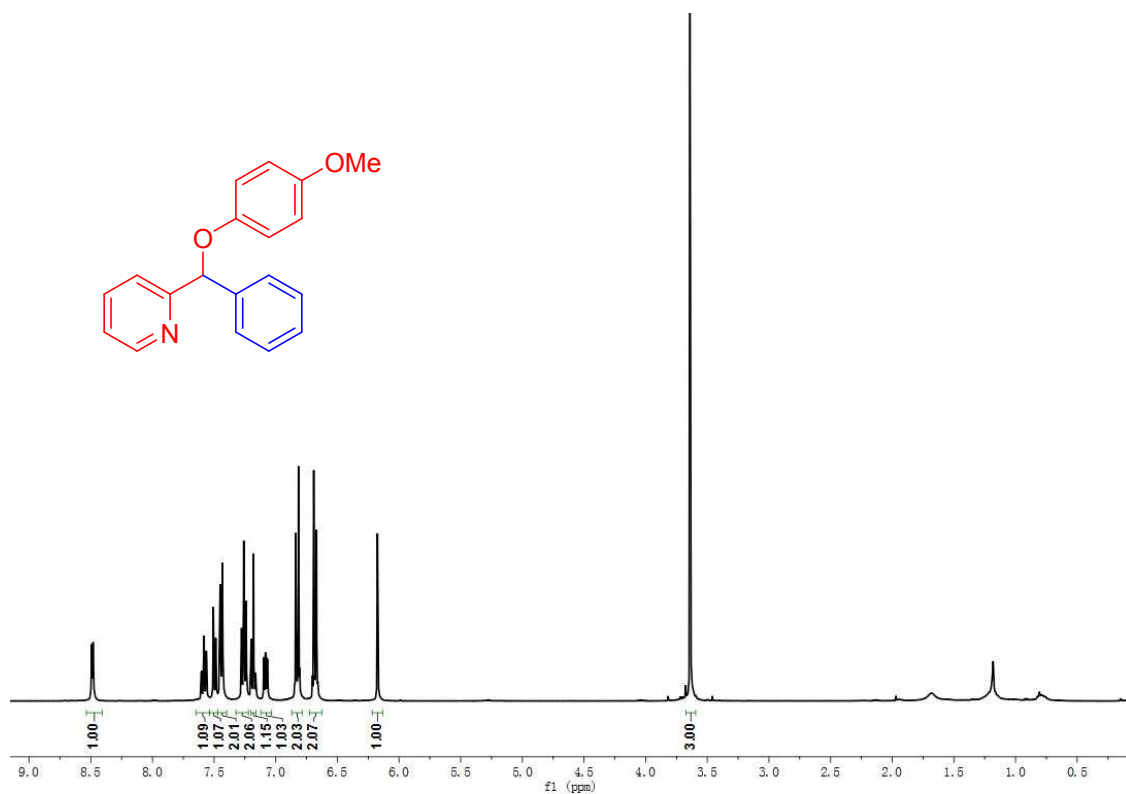
^1H NMR spectra (400 MHz, Chloroform-*d*) of 2-((4-(*tert*-Butyl)phenoxy)(phenyl)methyl)pyridine (3ja)



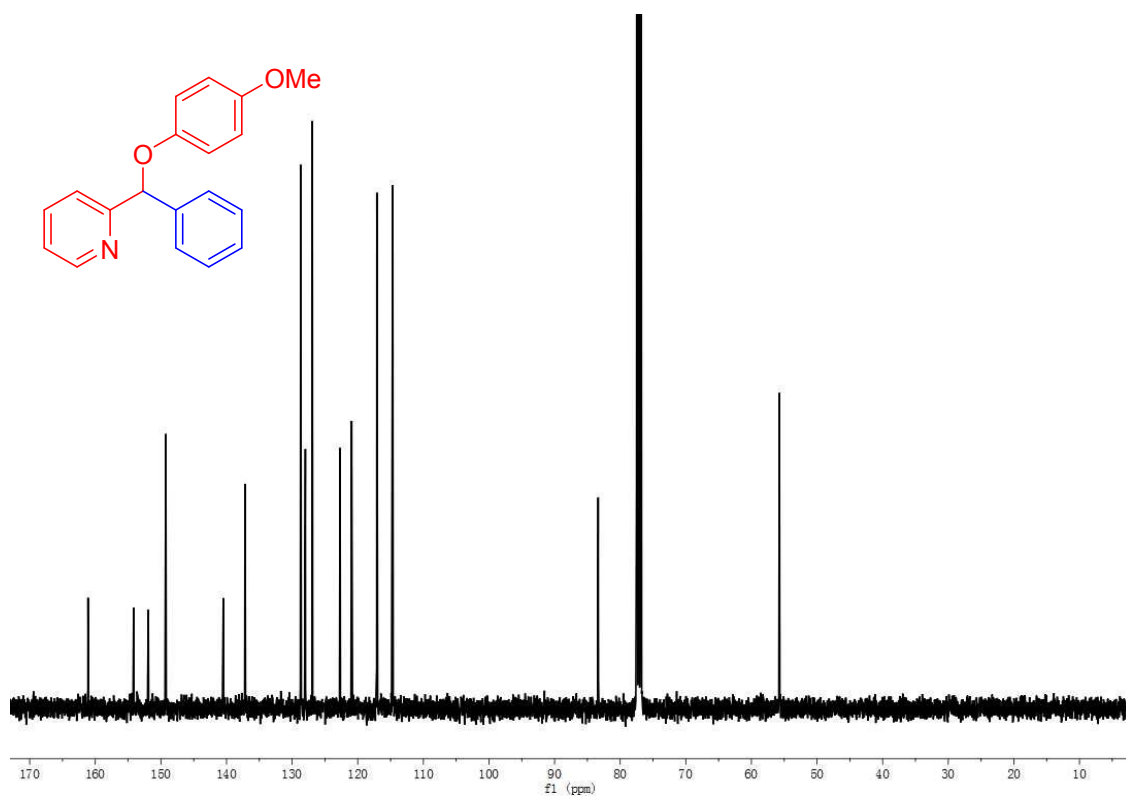
¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 2-((4-*tert*-Butylphenoxy)(phenyl)methyl)pyridine (3ja)



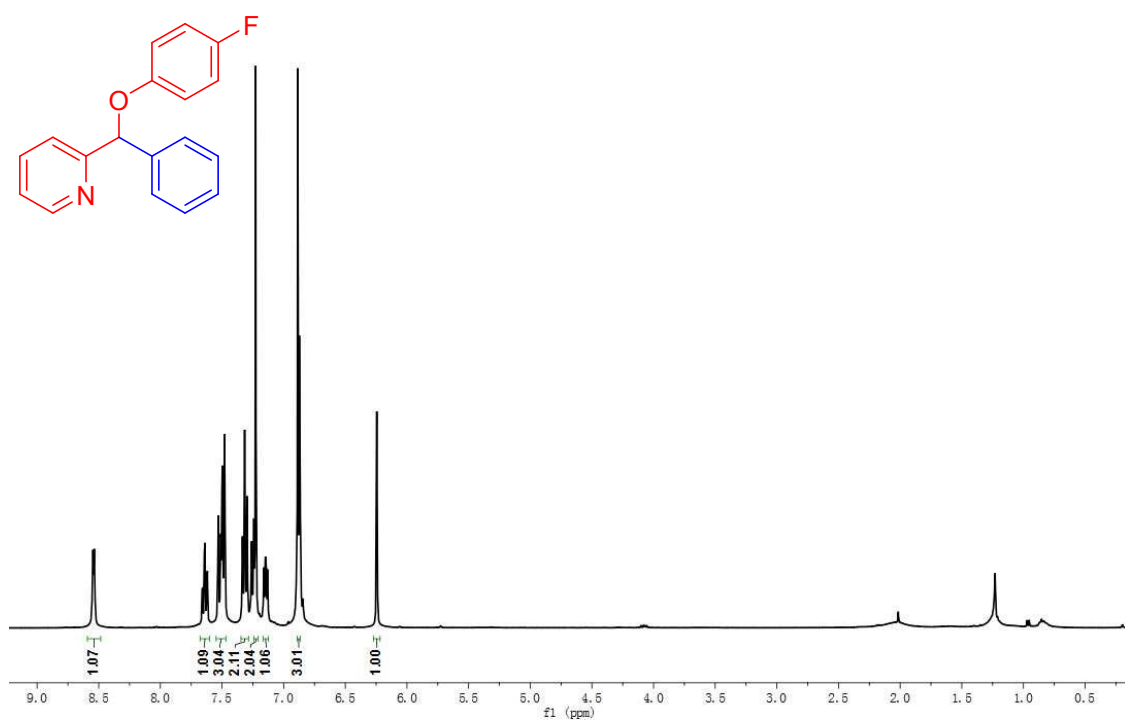
¹H NMR spectra (400 MHz, Chloroform-*d*) of 2-((4-Methoxyphenoxy)(phenyl)methyl)pyridine (3ka)



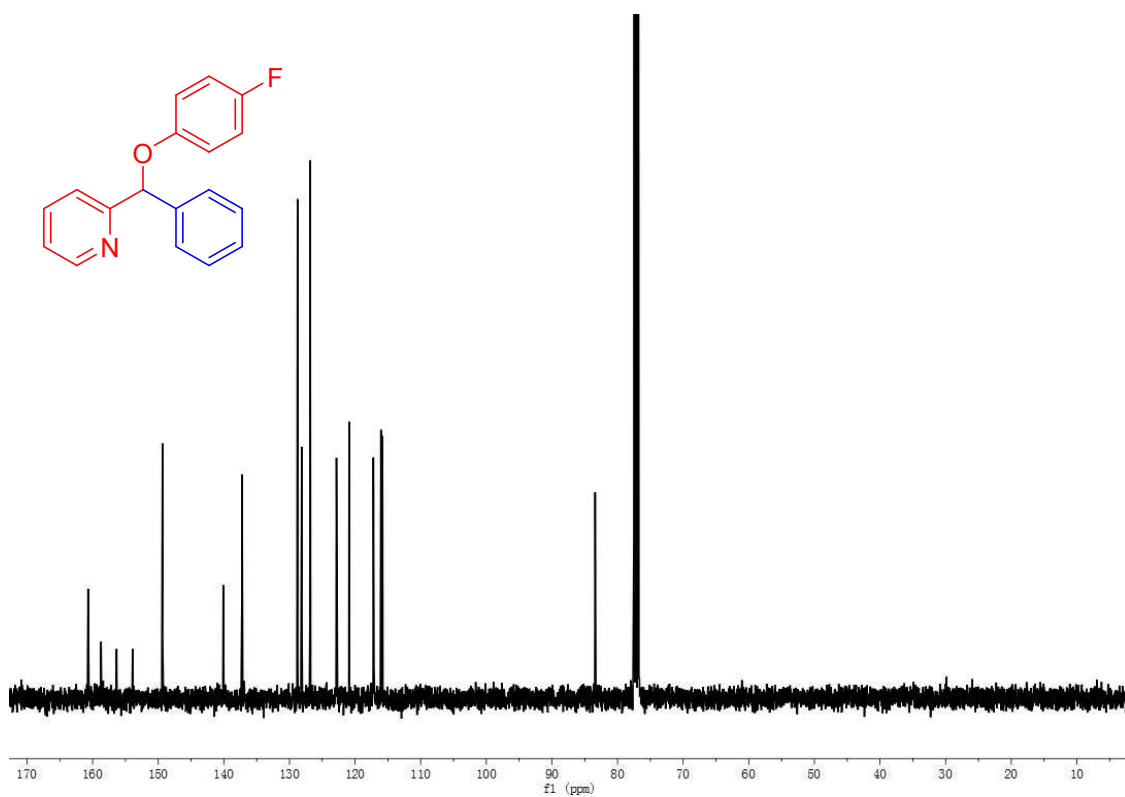
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 2-((4-Methoxyphenoxy)(phenyl)methyl)pyridine (3ka)



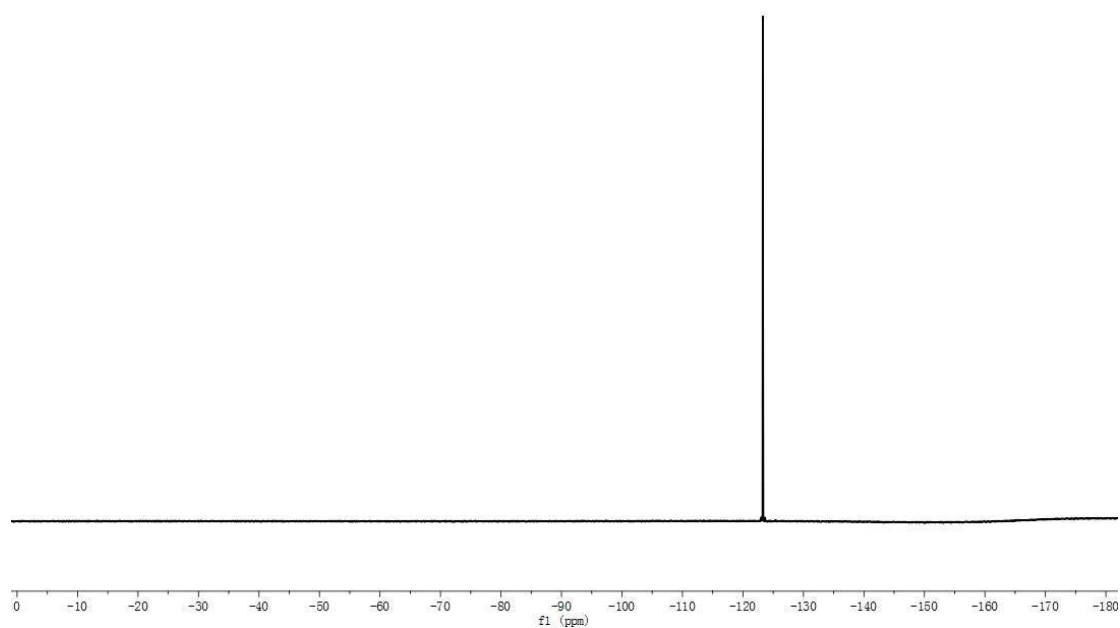
^1H NMR spectra (400 MHz, Chloroform-*d*) of 2-((4-Fluorophenoxy)(phenyl)methyl)pyridine (3la)



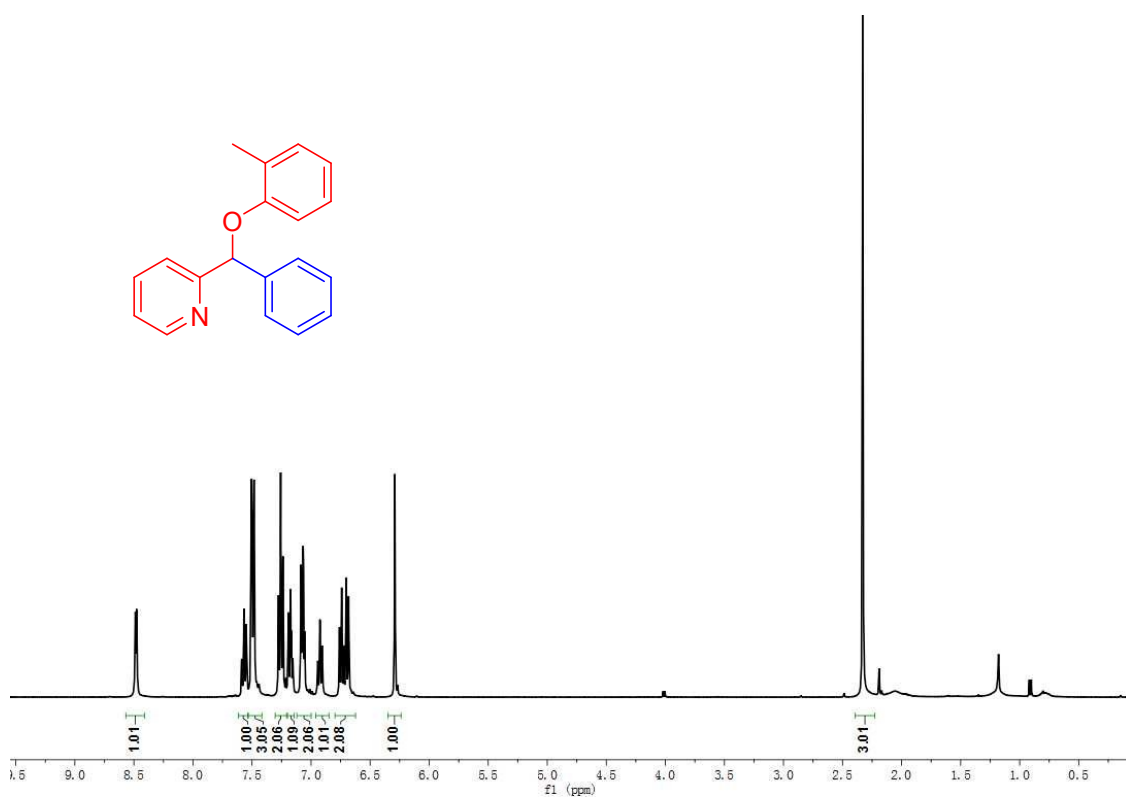
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 2-((4-Fluorophenoxy)(phenyl)methyl)pyridine (3la)



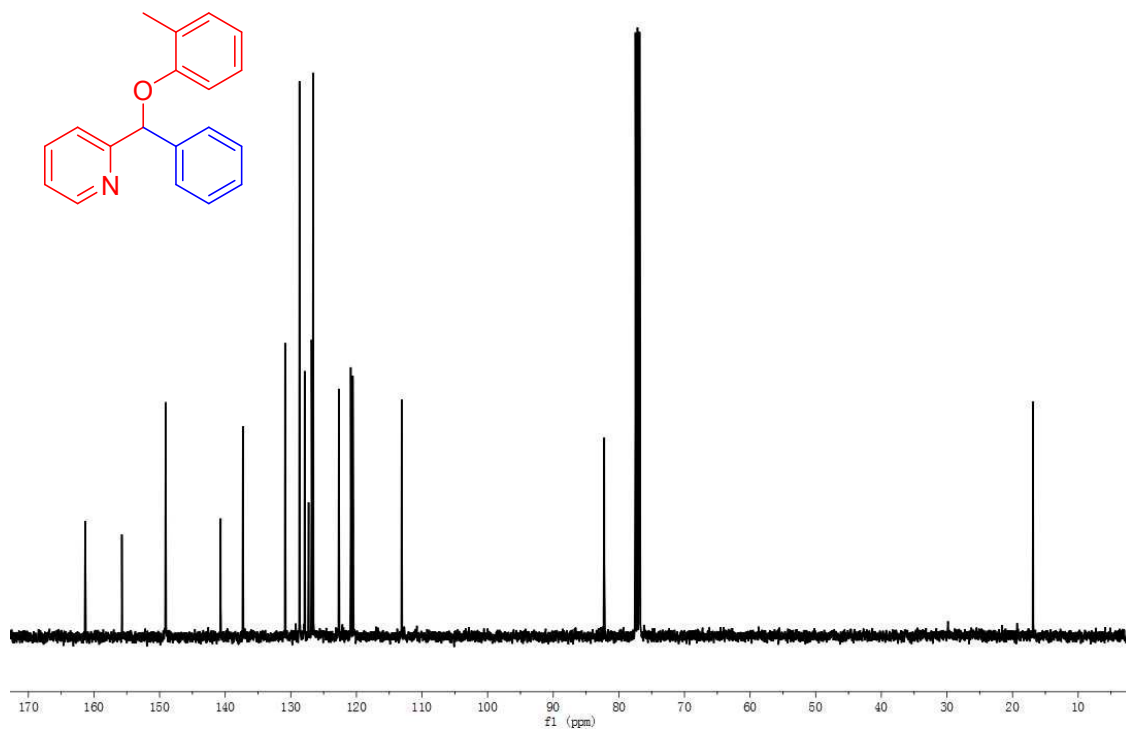
^{19}F NMR spectra (376 MHz, Chloroform-*d*) of 2-((4-Fluorophenoxy)(phenyl)methyl)pyridine (3la)



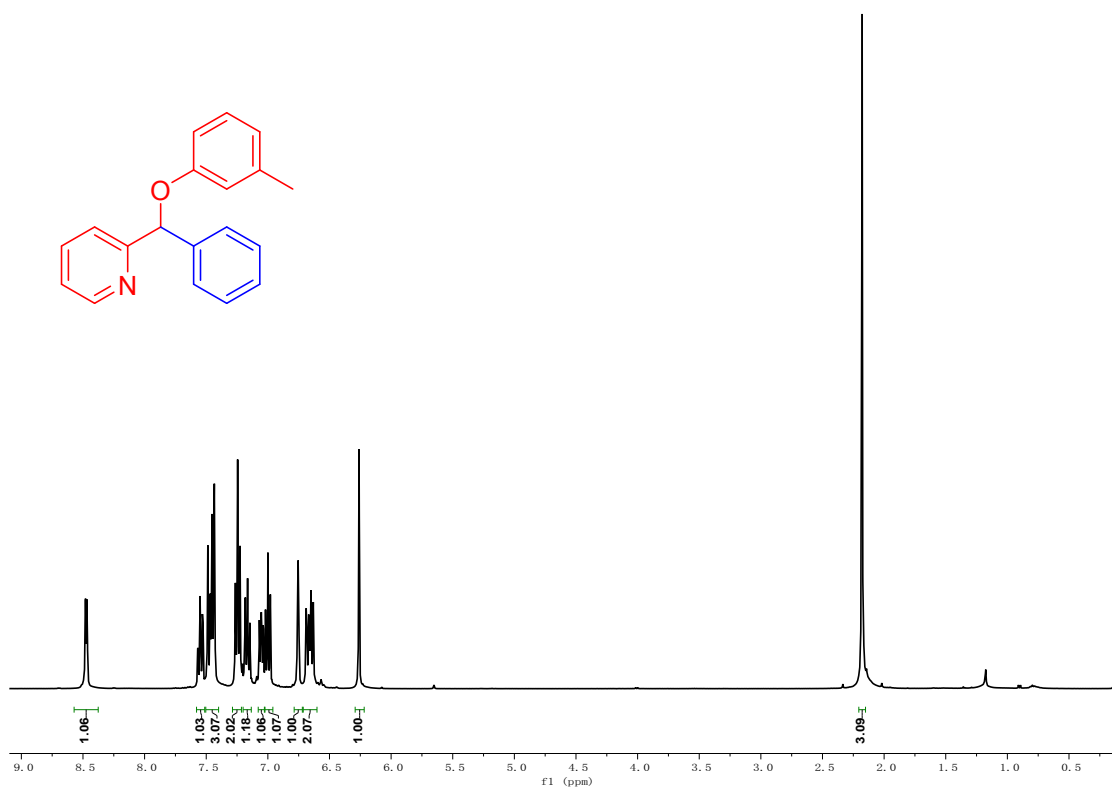
^1H NMR spectra (400 MHz, Chloroform-*d*) of 2-(Phenyl(*o*-tolylloxy)methyl)pyridine (3ma)



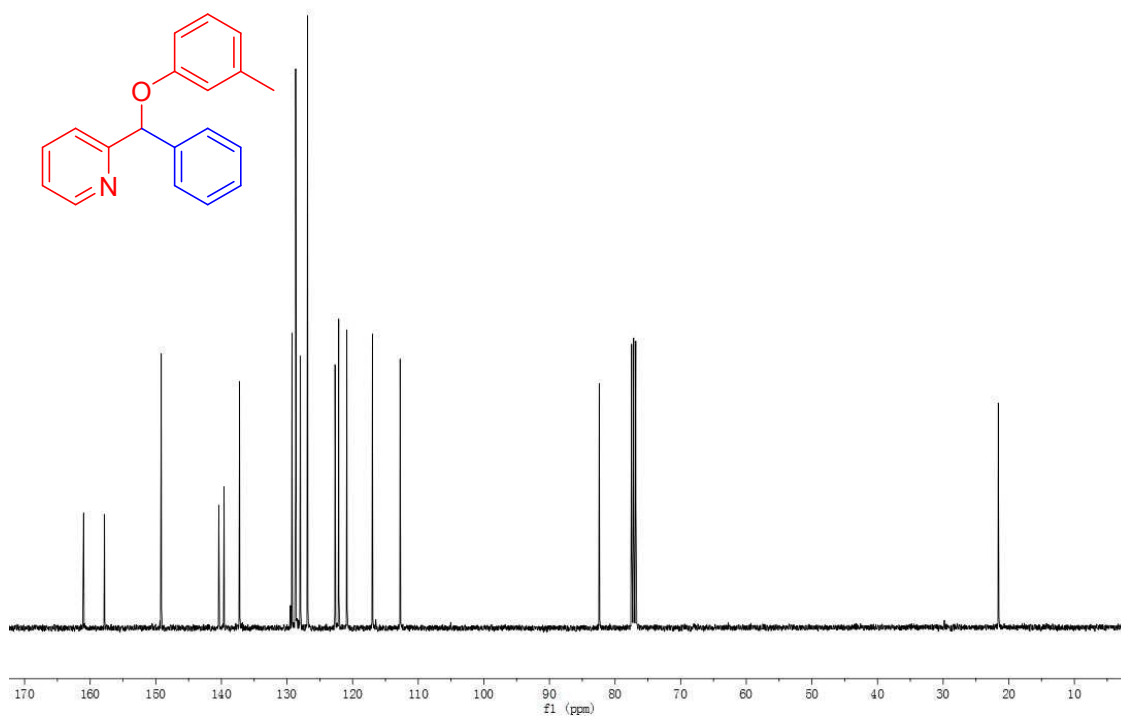
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 2-(Phenyl(o-tolyloxy)methyl)pyridine (3ma)



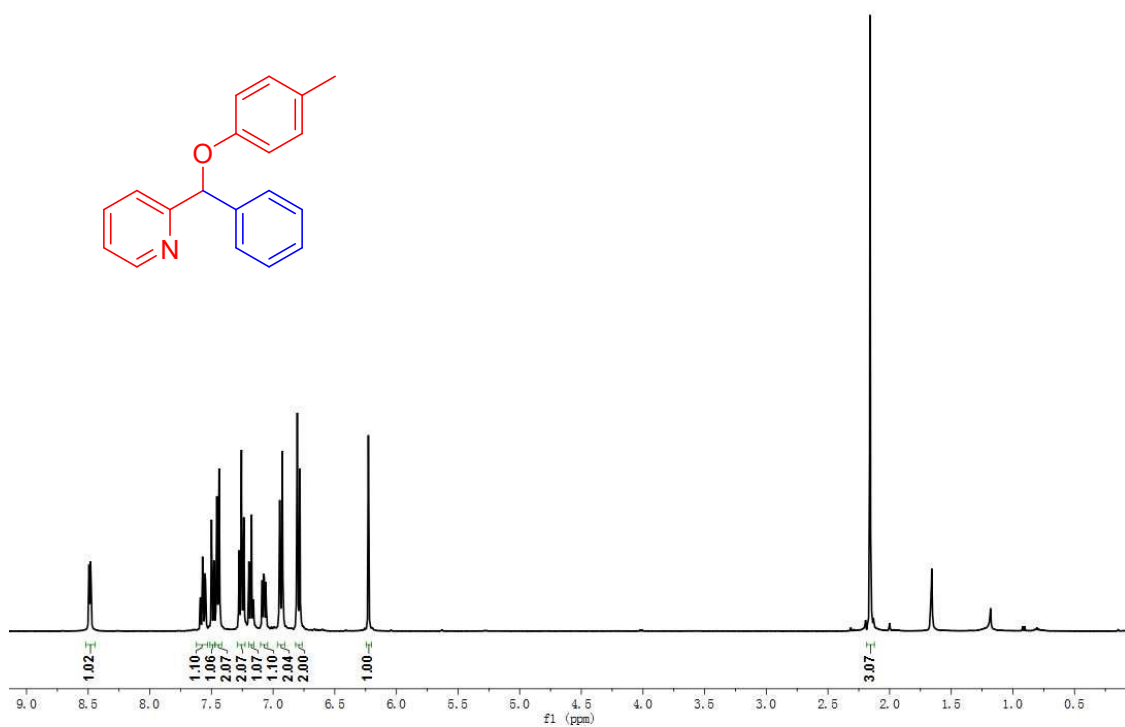
^1H NMR spectra (400 MHz, Chloroform-*d*) of 2-(Phenyl(m-tolyloxy)methyl)pyridine (3na)



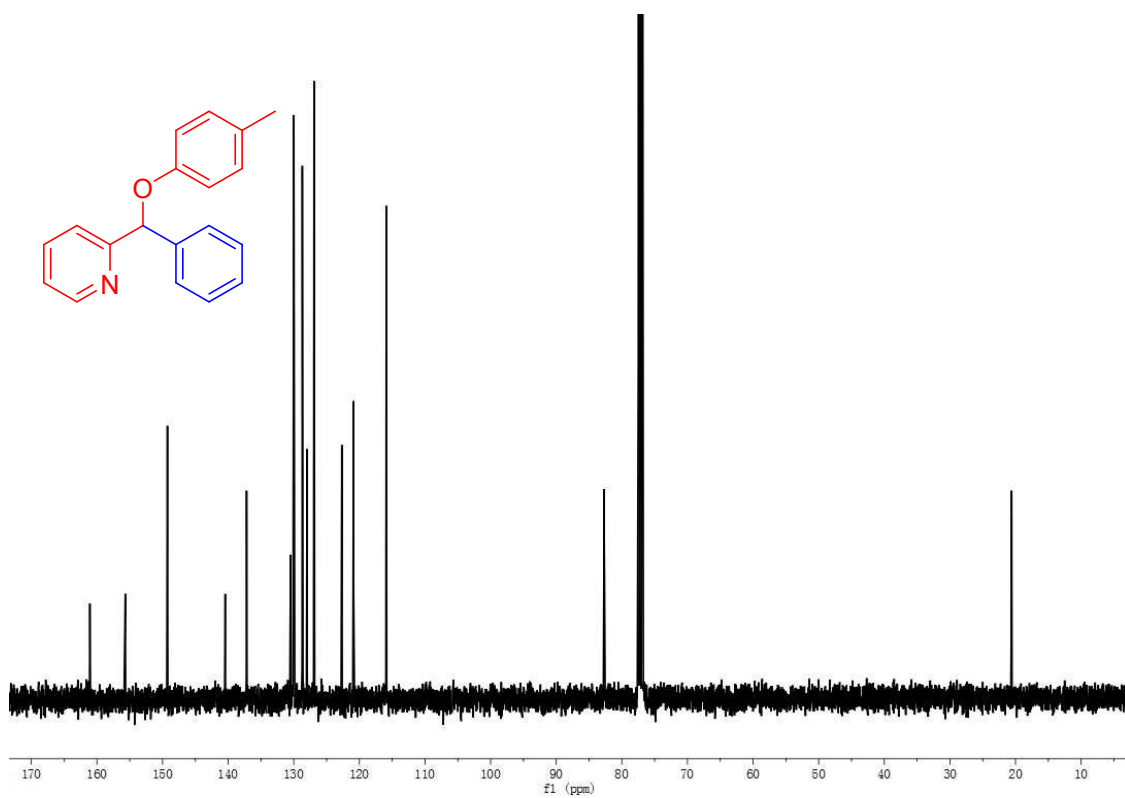
¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 2-(Phenyl(m-tolyloxy)methyl)pyridine (3na)



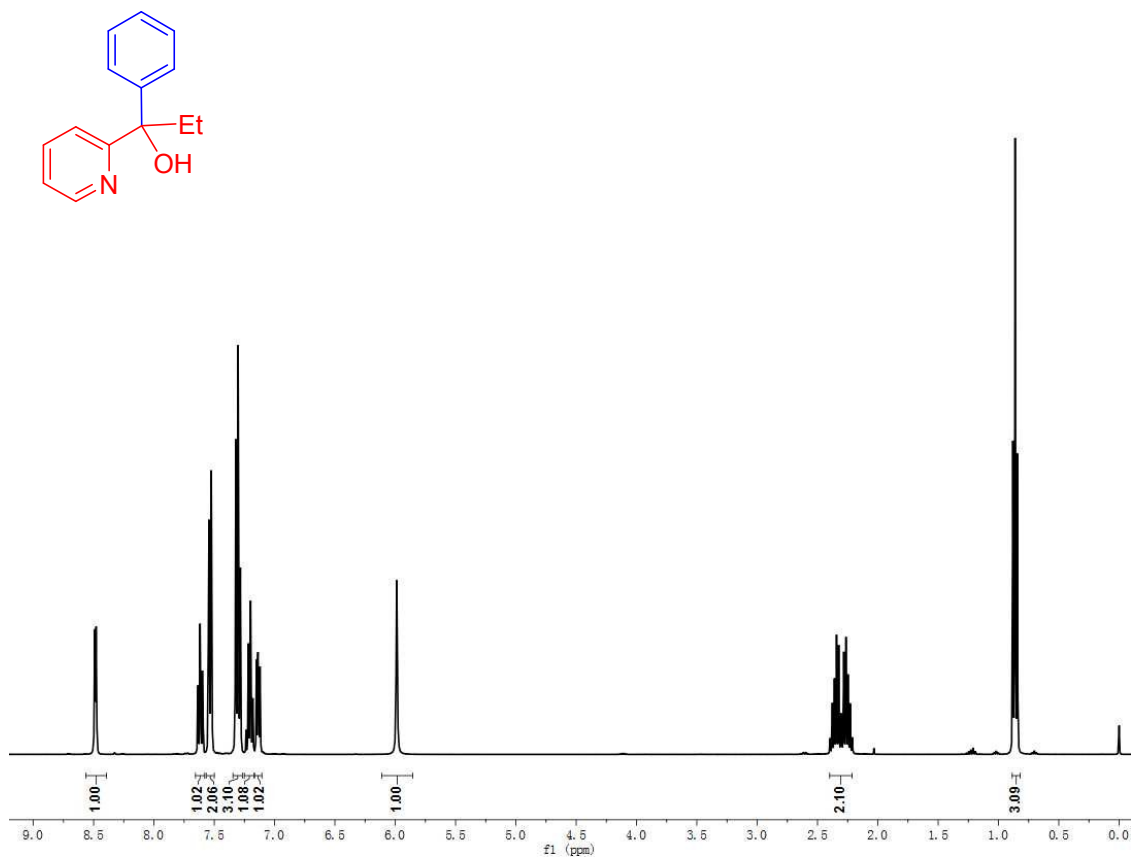
¹H NMR spectra (400 MHz, Chloroform-*d*) of 2-(Phenyl(p-tolyloxy)methyl)pyridine (3oa)



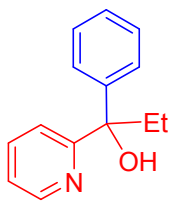
¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 2-(Phenyl(p-tolyloxy)methyl)pyridine (30a)

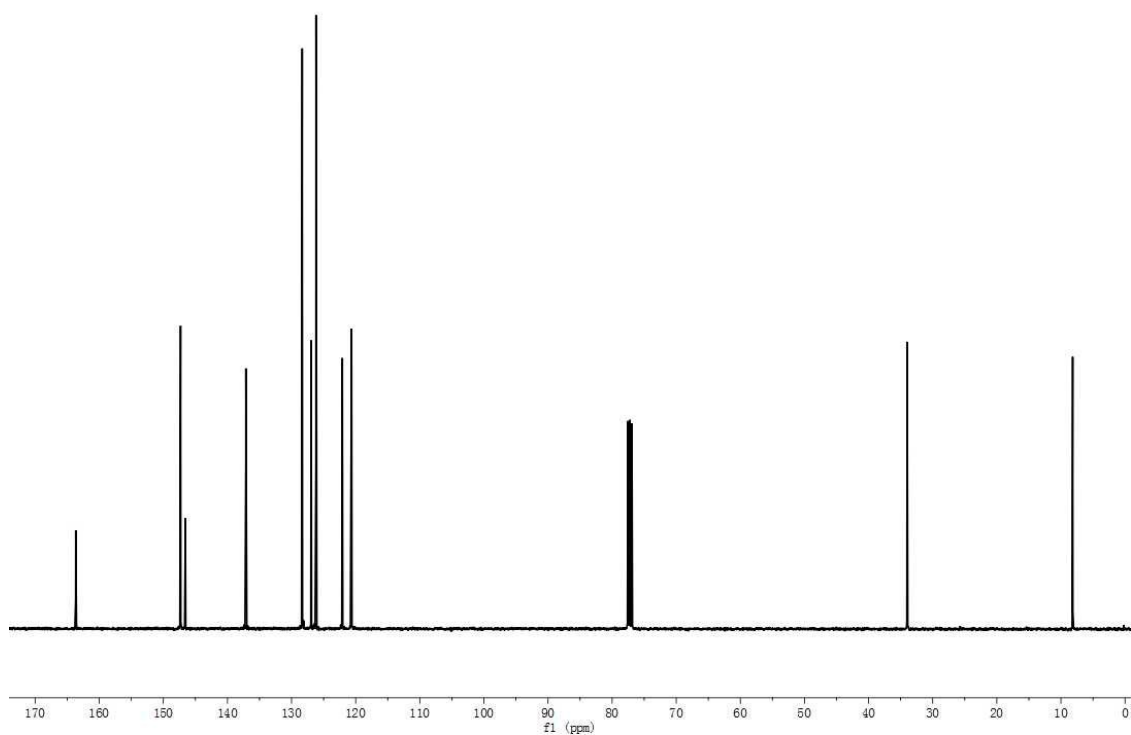


¹H NMR spectra (400 MHz, Chloroform-*d*) of 1-Phenyl-1-(pyridin-2-yl)propan-1-ol (4ea)

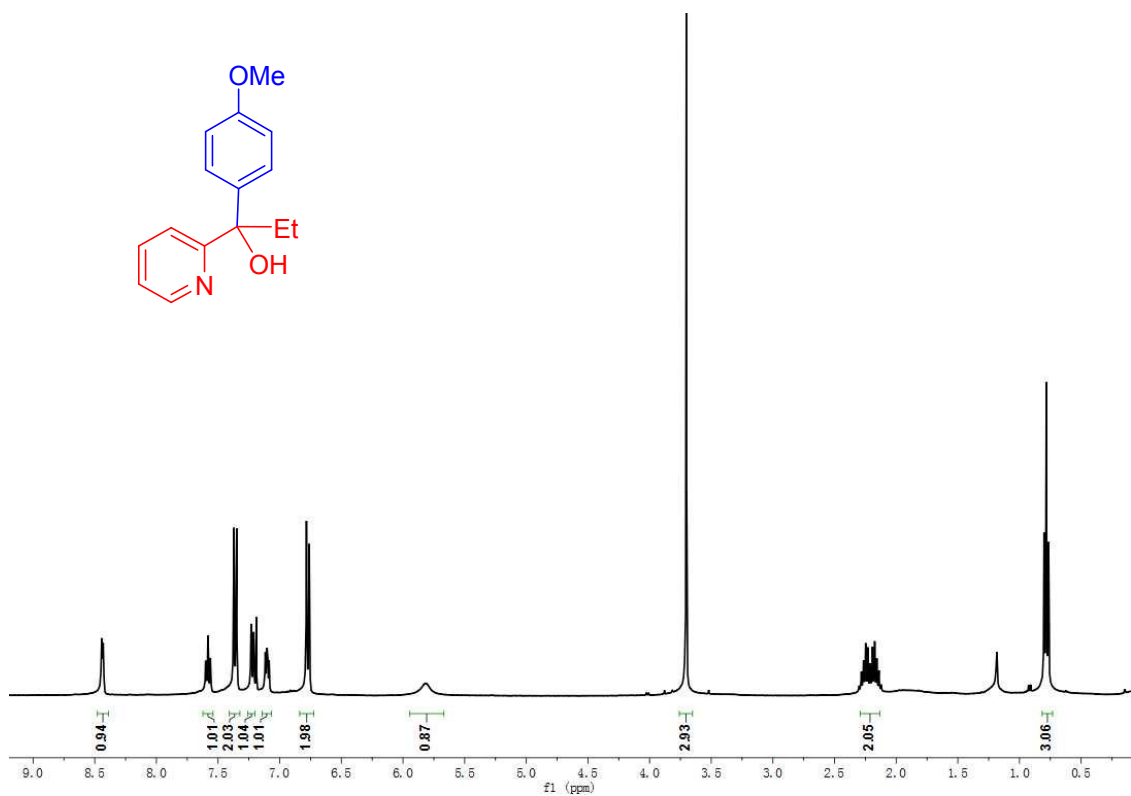


¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 1-Phenyl-1-(pyridin-2-yl)propan-1-ol (4ea)

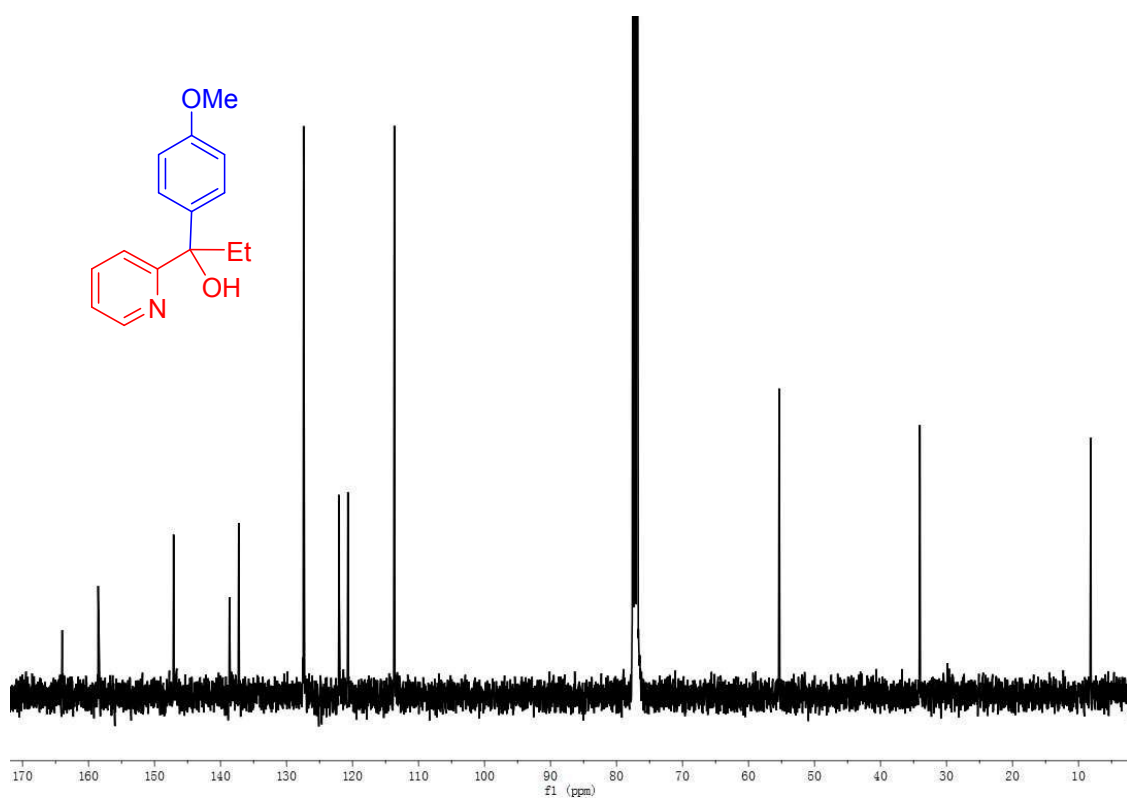




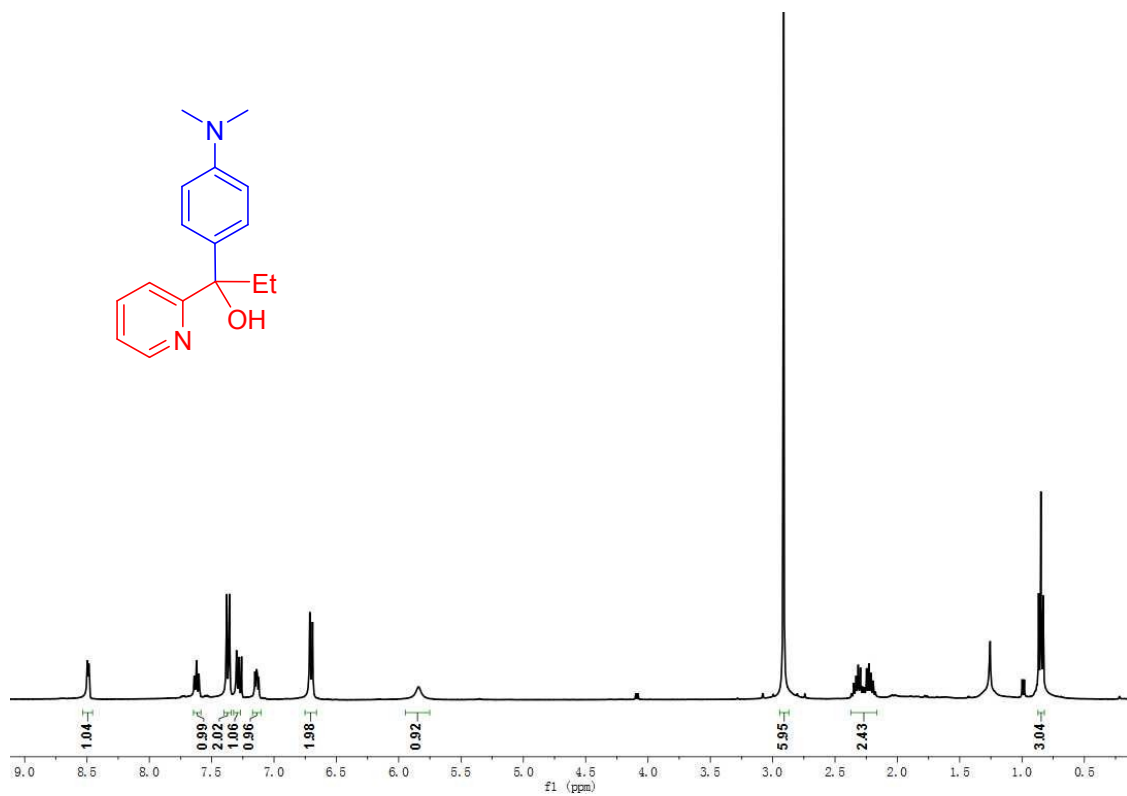
¹H NMR spectra (400 MHz, Chloroform-*d*) of 1-(4-Methoxyphenyl)-1-(pyridin-2-yl)propan-1-ol (4eb)



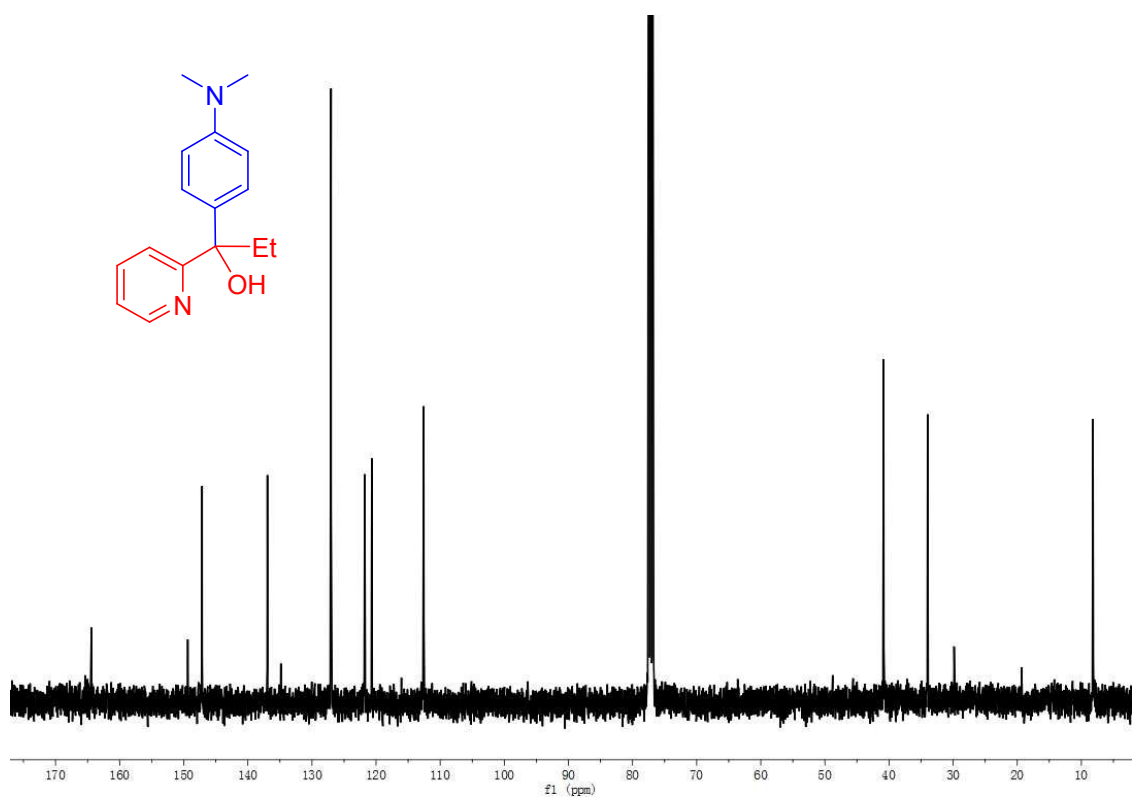
¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 1-(4-Methoxyphenyl)-1-(pyridin-2-yl)propan-1-ol (4eb)



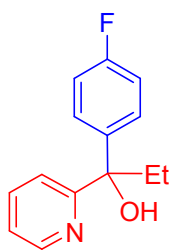
¹H NMR spectra (400 MHz, Chloroform-*d*) of 1-(4-(Dimethylamino)phenyl)-1-(pyridin-2-yl)propan-1-ol (4ec)

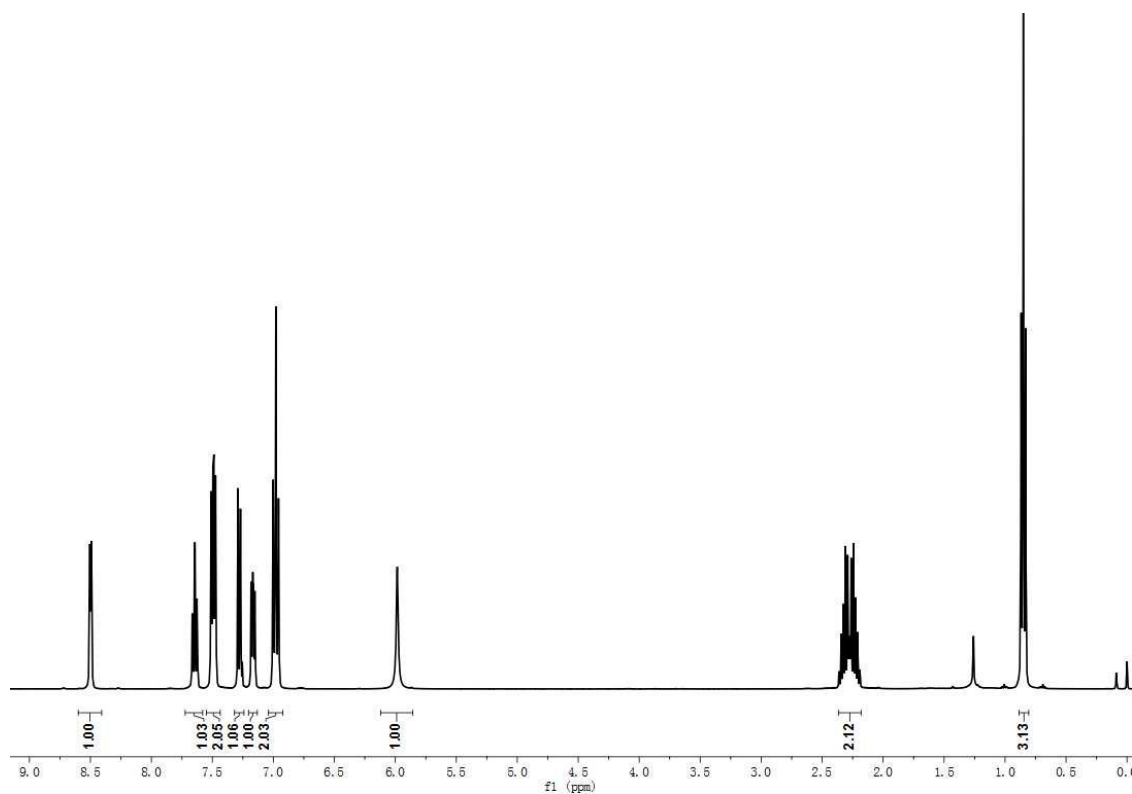


¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 1-(4-(Dimethylamino)phenyl)-1-(pyridin-2-yl)propan-1-ol (4ec)

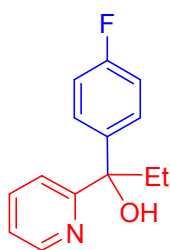


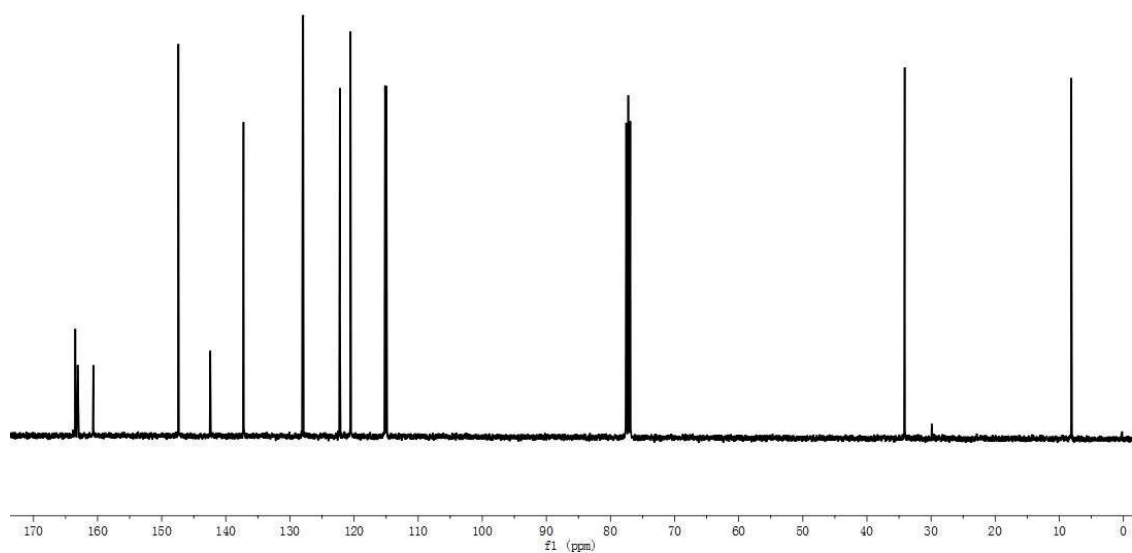
¹H NMR spectra (400 MHz, Chloroform-*d*) of 1-(4-Fluorophenyl)-1-(pyridin-2-yl)propan-1-ol (4ed)



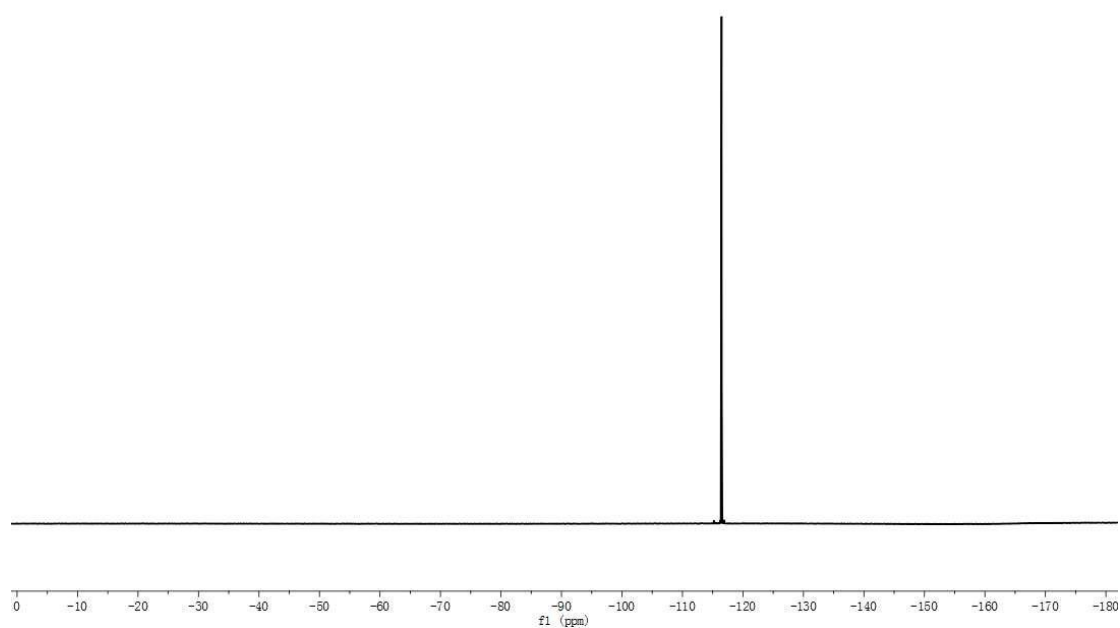


$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, CDCl_3) of 1-(4-Fluorophenyl)-1-(pyridin-2-yl)propan-1-ol (4ed)

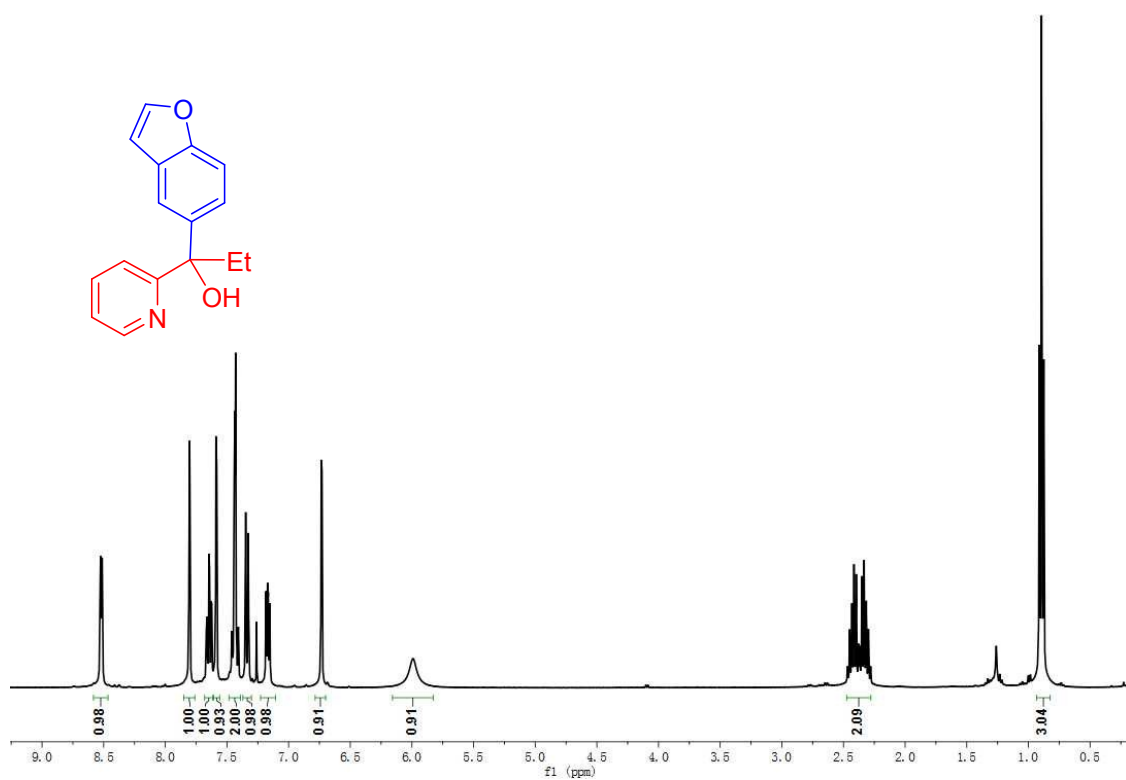




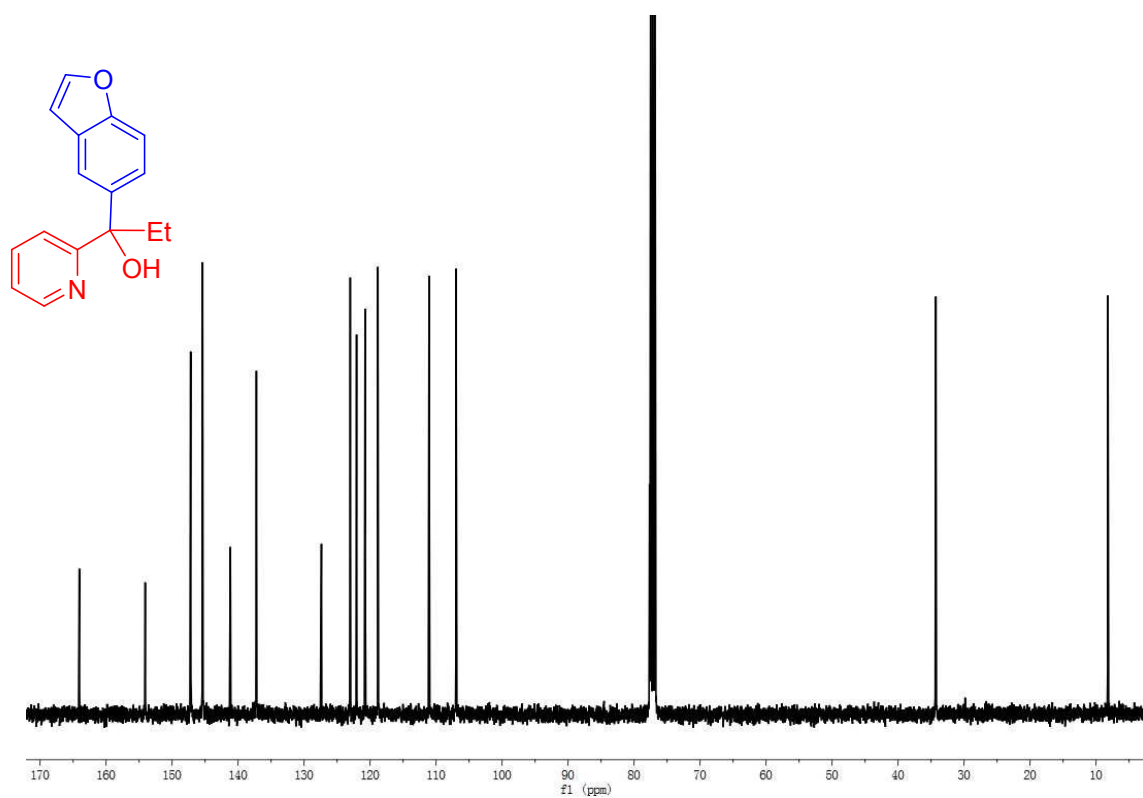
^{19}F NMR spectra (376 MHz, Chloroform-*d*) of 1-(4-Fluorophenyl)-1-(pyridin-2-yl)propan-1-ol (4ed)



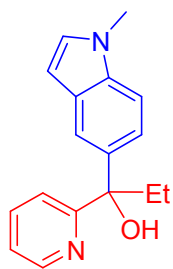
¹H NMR spectra (400 MHz, Chloroform-*d*) of 1-(Benzofuran-5-yl)-1-(pyridin-2-yl)propan-1-ol (4eg)

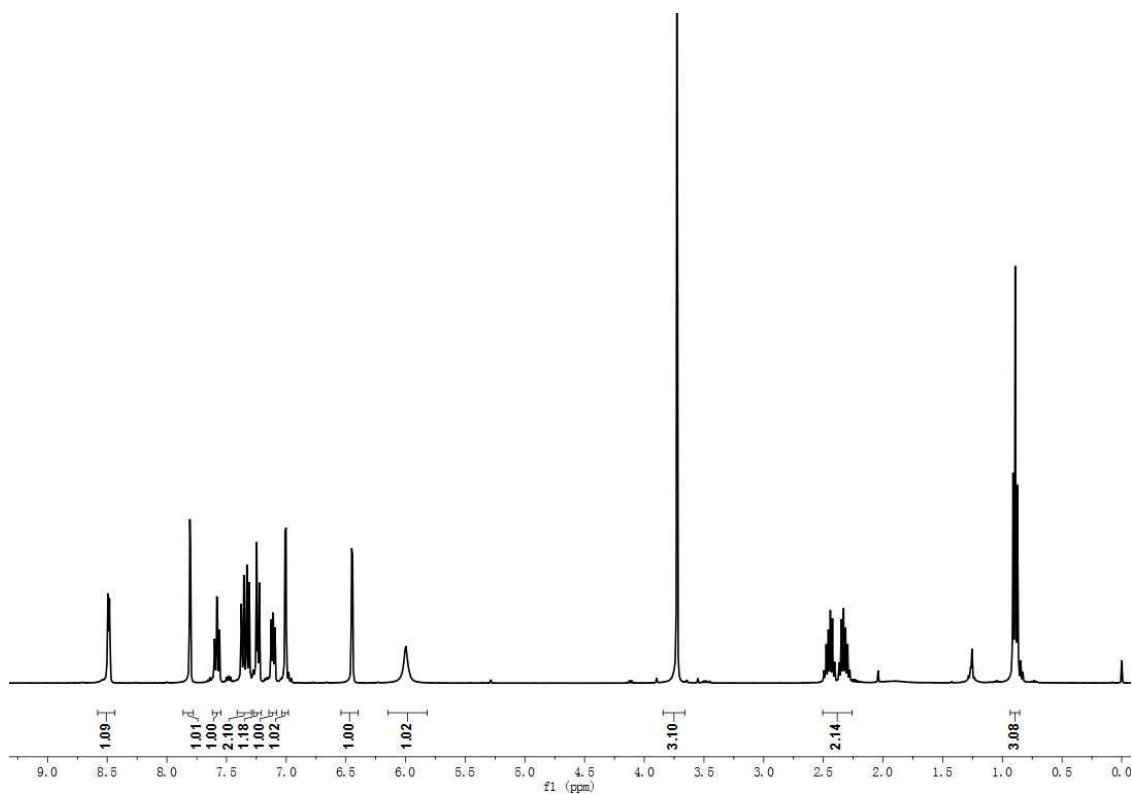


¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 1-(Benzofuran-5-yl)-1-(pyridin-2-yl)propan-1-ol (4eg)

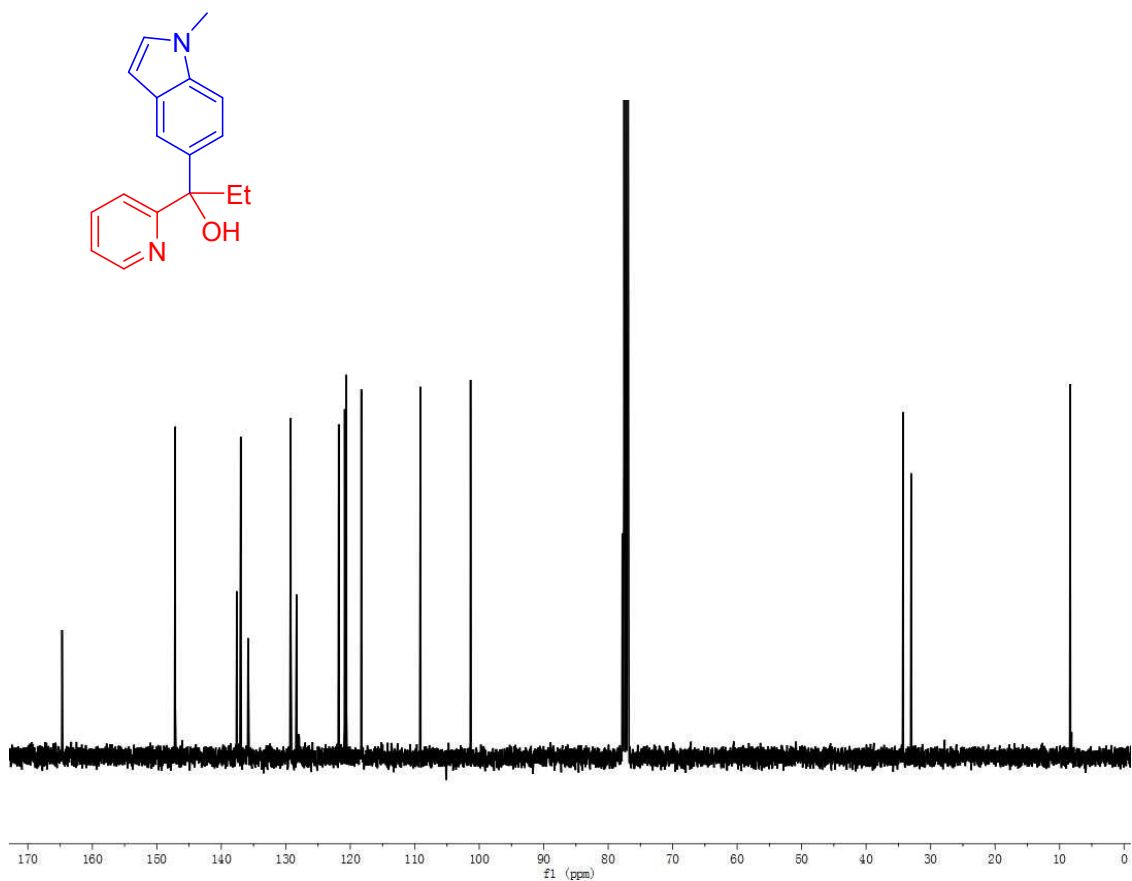


¹H NMR spectra (400 MHz, Chloroform-*d*) of 1-(1-Methyl-1H-indol-5-yl)-1-(pyridin-2-yl)propan-1-ol (4eh)

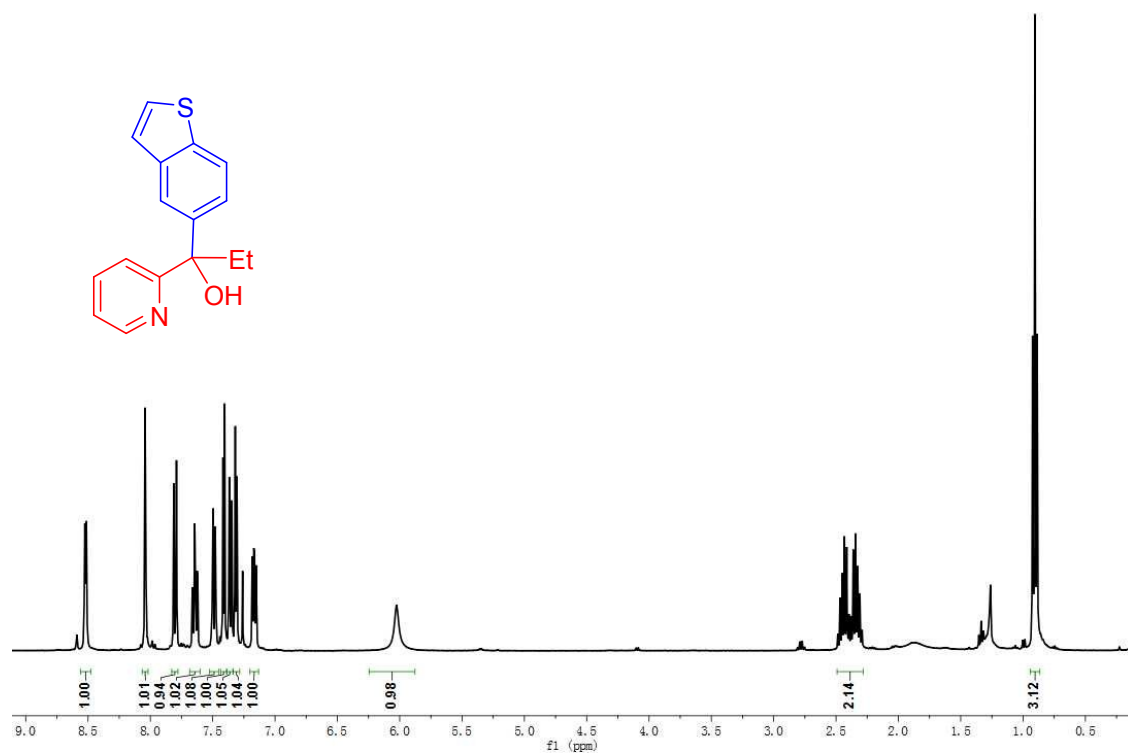




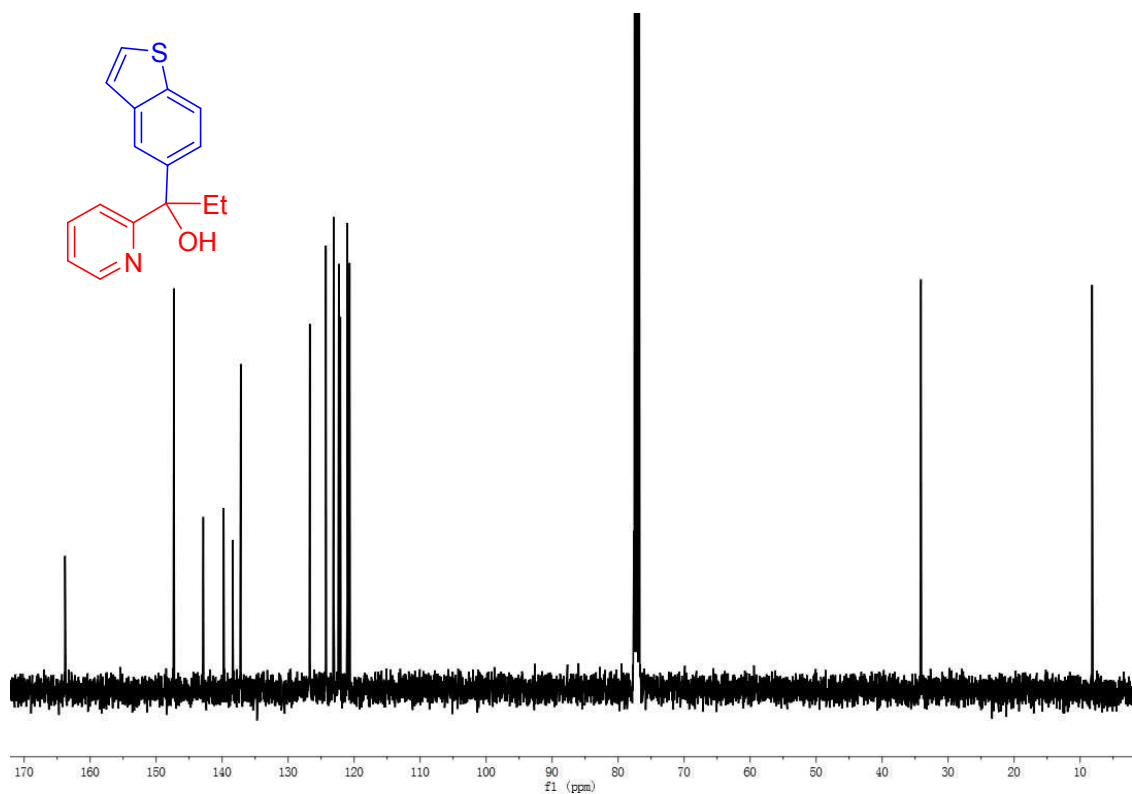
$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform- d) of 1-(1-Methyl-1H-indol-5-yl)-1-(pyridin-2-yl)propan-1-ol (4eh)



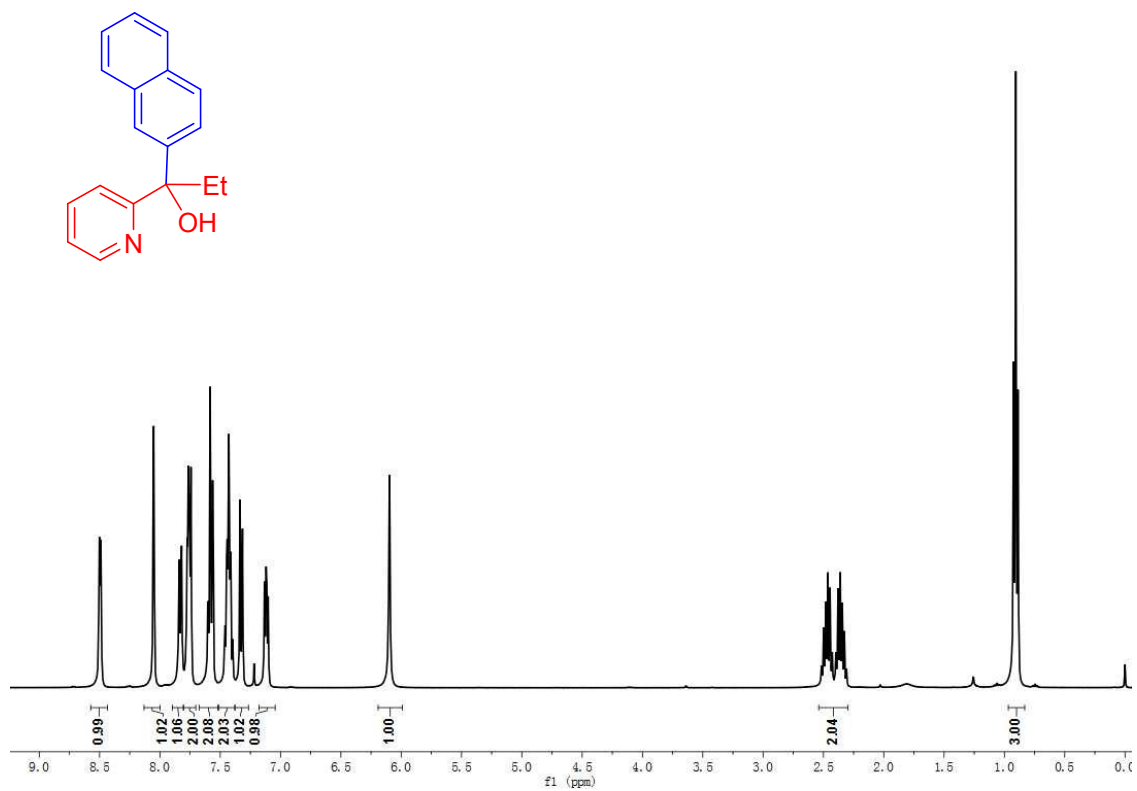
¹H NMR spectra (400 MHz, Chloroform-*d*) of 1-(Benzo[*b*]thiophen-5-yl)-1-(pyridin-2-yl)propan-1-ol (4ei)



¹³C{¹H} NMR spectra (100 MHz, Chloroform-*d*) of 1-(Benzo[*b*]thiophen-5-yl)-1-(pyridin-2-yl)propan-1-ol (4ei)



¹H NMR spectra (400 MHz, Chloroform-*d*) of 1-(Naphthalen-2-yl)-1-(pyridin-2-yl)propan-1-ol (4ej)



$^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100 MHz, Chloroform-*d*) of 1-(Naphthalen-2-yl)-1-(pyridin-2-yl)propan-1-ol (4ej)

