

Supporting Information

Base-controlled chemoselectivity: Direct coupling of acetonitriles and alcohols to synthesise α -alkylated arylacetonitriles or acetamides

Chen Li, Liang Bai, Min-Tong Ge, Ai-Bao Xia,* Ying Wang, Yuan-Rui Qiu and Dan-Qian Xu*

Catalytic Hydrogenation Research Centre, State Key Laboratory Breeding Base of Green Chemistry-Synthesis Technology,

Zhejiang University of Technology, Hangzhou, 310014, Zhejiang, China

Fax (+86) 0571 88320066;

E-mail: xiaaibao@zjut.edu.cn; chrc@zjut.edu.cn

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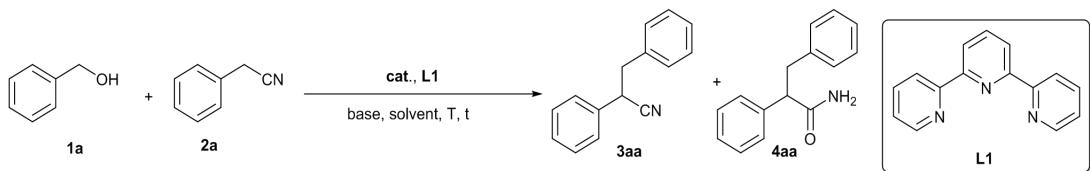
1. General information

NMR data were obtained on Bruker AVANCE III 500MHz for ¹H at 500 MHz and for ¹³C at 126 MHz with TMS as the internal standard; HRMS data were measured on an Agilent 6120 LC/TOF-MS with ESI source or Waters Premier GC/TOF-MS with EI source. Column chromatography and flash chromatography experiments were conducted using silica gel GF254 (200-300 mesh) eluting with ethyl ether and petroleum ether. TLC experiments were carried out on glass-backed silica plates. Unless otherwise noted, chemicals were used without purification as commercially available.

2. General procedures

Screening of the reaction conditions

Table S1 Optimization of Reaction Conditions^a

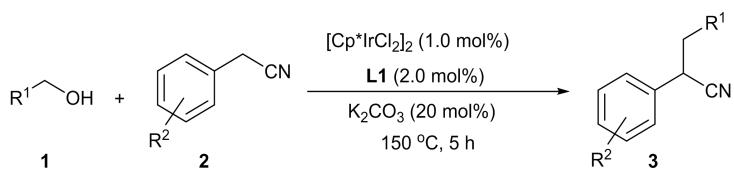


entry	cat.	base	base loading /mol%	solvent	temperature /°C	yield ^b /%	
						3aa	4aa
1	-	'BuOK	20	toluene	150	0	0
2	[CP*IrCl ₂] ₂	'BuOK	20	toluene	150	12	80
3	[Ir(cod)Cl ₂] ₂	'BuOK	20	toluene	150	41	28
4	IrCl ₃	'BuOK	20	toluene	150	48	18
5	RuCl ₂ (CO)(PPh ₃) ₃	'BuOK	20	toluene	150	35	8
6	[Cp*RuCl ₂] ₂	'BuOK	20	toluene	150	0	0
7	[CP*IrCl ₂] ₂	-	20	toluene	150	0	0
8	[CP*IrCl ₂] ₂	Cs ₂ CO ₃	20	toluene	150	80	13
9	[CP*IrCl ₂] ₂	'BuONa	20	toluene	150	24	61
10	[CP*IrCl ₂] ₂	KHCO ₃	20	toluene	150	83	7
11	[CP*IrCl ₂] ₂	KOH	20	toluene	150	40	49
12	[CP*IrCl ₂] ₂	K ₂ CO ₃	20	toluene	150	85	0
13	[CP*IrCl ₂] ₂	K ₂ CO ₃	10	toluene	150	65	0
14	[CP*IrCl ₂] ₂	K ₂ CO ₃	15	toluene	150	73	0
15	[CP*IrCl ₂] ₂	K ₂ CO ₃	30	toluene	150	82	7
16	[CP*IrCl ₂] ₂	K ₂ CO ₃	20	xylene	150	83	0
17	[CP*IrCl ₂] ₂	K ₂ CO ₃	20	DMSO	150	0	0
18	[CP*IrCl ₂] ₂	K₂CO₃	20	tert-amyl alcohol	150	95	0
19	[CP*IrCl ₂] ₂	K ₂ CO ₃	20	tert-amyl alcohol	100	45	0
20	[CP*IrCl ₂] ₂	K ₂ CO ₃	20	tert-amyl alcohol	120	62	0
21	[CP*IrCl ₂] ₂	K ₂ CO ₃	20	tert-amyl alcohol	140	82	0
22	[CP*IrCl ₂] ₂	'BuOK	20	toluene	150	12	80
23	[CP*IrCl ₂] ₂	'BuOK	10	toluene	150	85	5

24	[CP*IrCl ₂] ₂	'BuOK	15	toluene	150	73	10
25	[CP*IrCl ₂] ₂	'BuOK	30	toluene	150	10	73
26	[CP*IrCl ₂] ₂	'BuOK	20	xylene	150	15	75
27	[CP*IrCl ₂] ₂	'BuOK	20	DMSO	150	-	-
28	[CP*IrCl ₂] ₂	'BuOK	20	tert-amyl alcohol	150	0	90
29	[CP*IrCl ₂] ₂	'BuOK	20	tert-amyl alcohol	100	10	0
30	[CP*IrCl ₂] ₂	'BuOK	20	tert-amyl alcohol	120	55	7
31	[CP*IrCl ₂] ₂	'BuOK	20	tert-amyl alcohol	140	46	32

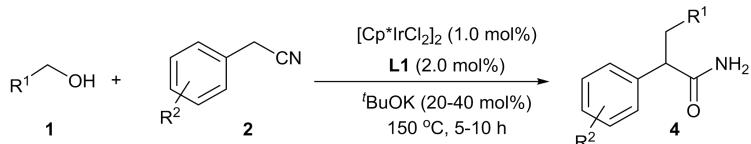
^aReaction conditions: benzonitrile **2a** (1.0 mmol), benzyl alcohol **1a** (1.1 mmol), cat. (1.0 mol%), **L1** (2.0 mmol%) and base (20 mol%) were heated in solvent (1.0 mL) under N₂. ^bIsolated yield.

General procedures for the synthesis of products 3



Alcohol **1** (1.1 mmol), arylacetonitrile **2** (1.0 mmol, 1.0 equiv.), [Cp*IrCl₂]₂ (1.0 mol%), **L1** (2.0 mol%), and K₂CO₃ (20 mol%) were heated at 150 °C in *tert*-amyl alcohol (1.0 mL) for 5 h under N₂. The mixture was cooled to room temperature when it was over. Then the mixture was extracted with EtOAc (10.0 mL × 3). The combined organic phases were washed with water and brine, dried (MgSO₄), filtered, and concentrated. The residue was purified by flash chromatography on silica gel with petroleum ether/ethyl acetate (20:1 to 5:1) to afford the products **3**.

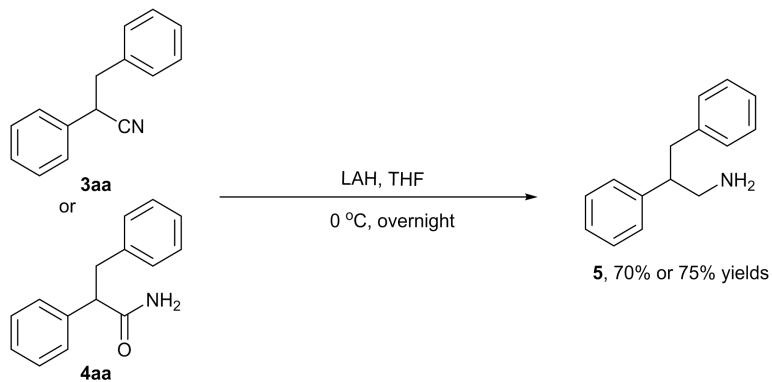
General procedures for the synthesis of products 4



Alcohol **1** (1.1 mmol), acetonitrile **2** (1.0 mmol, 1.0 equiv.), [Cp*IrCl₂]₂ (1.0 mol%), **L1** (2.0 mol%), and 'BuOK (20 mol%) were heated at 150 °C in *tert*-amyl alcohol (1.0 mL) for 5 h under N₂. The mixture was cooled to room temperature when it was over. Then the mixture was extracted with EtOAc (10.0 mL × 3). The combined organic phases were washed with water and brine, dried (MgSO₄), filtered, and concentrated. The residue was purified by flash chromatography on silica gel with petroleum ether/ethyl acetate (5:1 to 1:1) to afford the products **4**.

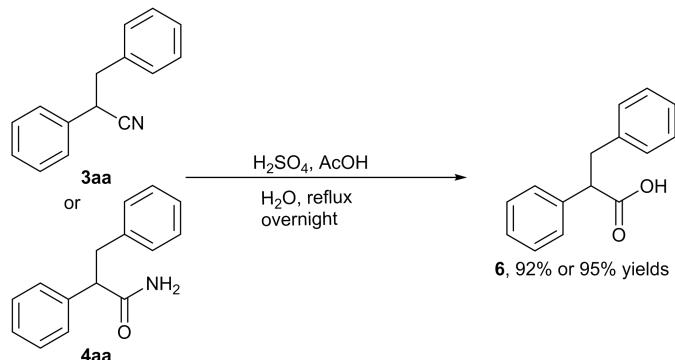
3. Functional group transformations and synthesis of drug fragment 7

Procedure for the synthesis of the product 5 from 3aa and 4aa



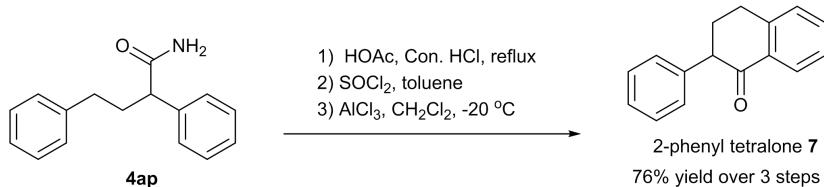
3aa (1.0 mmol, 207.0 mg) or **4aa** (1.0 mmol, 225.0 mg) in dry THF (10.0 mL) was dropwise added to a suspension of 2.5 mL LiAlH₄ (1.0 M) in dry THF at 0 °C under N₂. The mixture was stirred for 12 h. Then the mixture was very slowly quenched with 5.0 mL H₂O and diluted with DCM. The mixture was stirred another 0.5 h at room temperature. Then the mixture was extracted with EtOAc (10.0 mL × 3). The combined organic phases were washed with water and brine, dried (MgSO₄), filtered, and concentrated. The residue was purified by flash chromatography on silica gel with petroleum ether/ethyl acetate (1:1) to afford 2,3-diphenylpropan-1-amine **5** 147.7 mg in 70% yield or 158.3 mg in 75% yield, respectively.

Procedure for synthesis of the product **6** from **3aa** and **4aa**



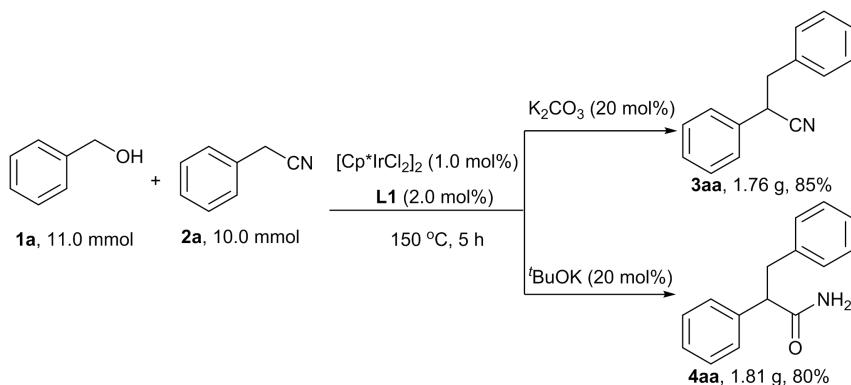
3aa (1.0 mmol, 207.0 mg) or **4aa** (1.0 mmol, 225.0 mg) in a mixture of concentrated H₂SO₄ (4.0 mL), H₂O (4.0 mL) and acetic acid (2.0 mL) was refluxed for 24 h. Then the mixture was quenched very slowly with 5.0 mL H₂O and diluted with 10.0 mL DCM. The mixture was stirred another 0.5 h at room temperature, then it was extracted with EtOAc (10.0 mL × 3). The combined organic phases were washed with water and brine, dried (MgSO₄), filtered, and concentrated. The residue was purified by flash chromatography on silica gel with petroleum ether/ethyl acetate (2:1) to afford 2,3-diphenylpropanoic acid **6** 207.9 mg in 92% yield or 214.7 mg in 95% yield, respectively.

Procedure for synthesis of the product **7**



Compound **4ap** (1.0 mmol, 239.0 mg) was mixed with 20.0 mL glacial acetic acid and 8.0 mL concentrated hydrochloric acid. The solution was refluxed for 24 h. After the reaction was completed, water was added to the reaction mixture and the resulting solution was extracted with DCM. The combined organic phases were washed with water and brine, dried (MgSO_4), filtered, and concentrated to obtain the carboxylic acid product. Then SOCl_2 (0.25 mL, 3.0 eq.) was slowly added into the cooled solution of the carboxylic acid product in 15.0 mL dry toluene at 0 °C. The mixture was stirred for another 15 min at 0 °C and refluxed for 4 h. The volatiles were removed under reduced pressure, and the resulting oil was dissolved in 10.0 mL dry DCM and was dropwise added into a stirred suspension of AlCl_3 (327.0 mg, 2.0 eq.) in 20.0 mL dry DCM at -20 °C within 40 min. The reaction was stirred at -20 °C for 2 h and quenched with 10.0 mL HCl (4.0 M). The mixture was extracted with EtOAc (20.0 mL × 3). The combined organic phases were washed with water and brine, dried (MgSO_4), filtered, and concentrated. The residue was purified by flash chromatography on silica gel with petroleum ether/ethyl acetate (10:1) to afford 2-phenyl tetralone **7** 168.7 mg in 76% yield over 3 steps.

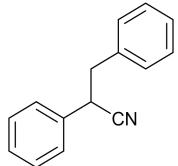
4. Multigram-scale synthesis of products **3aa** and **4aa**



Benzyl alcohol **1a** (11.0 mmol, 1.19 g), phenylacetonitrile **2a** (10.0 mmol, 1.17 g), catalyst $[\text{Cp}^*\text{IrCl}_2]_2$ (1.0 mol%), **L1** (2.0 mol%), K_2CO_3 (20 mol%) or ${}^t\text{BuOK}$ (20 mol%) were heated at 150 °C in *tert*-amyl alcohol (1.0 mL) for 5 h under N_2 . The mixture was cooled to room temperature when it was over. Then the mixture was extracted with EtOAc (20.0 mL × 3). The combined organic phases were washed with water and brine, dried (MgSO_4), filtered, and concentrated. The residue was purified by flash chromatography on silica gel with petroleum ether/ethyl acetate (20:3:1) to afford the product **3aa** 1.76 g in 85% yield or **4aa** 1.81 g in 80% yield.

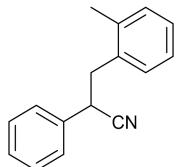
5. Characterization of products

2,3-diphenylpropanenitrile (**3aa**)



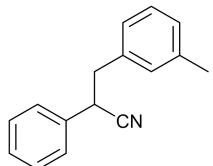
Yield: 95%; colorless solid; m.p. 46–48 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.42 – 7.26 (m, 8H), 7.20 – 7.14 (m, 2H), 4.03 (dd, *J* = 8.4, 6.4 Hz, 1H), 3.19 (qd, *J* = 13.6, 7.4 Hz, 2H); **13C NMR** (126 MHz, CDCl₃) δ 136.27, 135.22, 129.19 (2C), 128.99 (2C), 128.60 (2C), 128.18, 127.47 (2C), 127.36, 120.34, 42.16, 39.76; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₃NO ([M+Na]⁺) 230.0940, found 230.0944.

2-phenyl-3-(o-tolyl)propanenitrile (**3ab**)



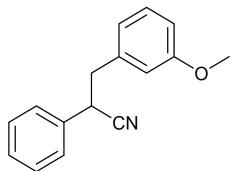
Yield: 85%; colorless oil; **1H NMR** (500 MHz, CDCl₃) δ 7.42 – 7.34 (m, 3H), 7.33 – 7.26 (m, 2H), 7.24 – 7.15 (m, 4H), 3.99 (dd, *J* = 8.7, 6.6 Hz, 1H), 3.21 (ddd, *J* = 20.4, 13.8, 7.7 Hz, 2H), 2.26 (s, 3H); **13C NMR** (126 MHz, CDCl₃) δ 136.25, 135.46, 134.67, 130.56, 130.03, 129.03 (2C), 128.20, 127.51, 127.37 (2C), 126.25, 120.45, 39.45, 38.73, 19.20; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₅N ([M+Na]⁺) 244.1097, found 244.1094.

2-phenyl-3-(m-tolyl)propanenitrile (**3ac**)



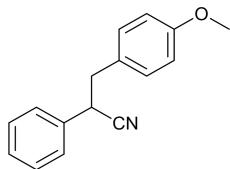
Yield: 92%; colorless oil; **1H NMR** (500 MHz, CDCl₃) δ 7.43 – 7.34 (m, 3H), 7.33 – 7.27 (m, 2H), 7.22 (t, *J* = 7.5 Hz, 1H), 7.12 (d, *J* = 7.6 Hz, 1H), 7.02 – 6.96 (m, 2H), 4.02 (dd, *J* = 8.6, 6.3 Hz, 1H), 3.15 (qd, *J* = 13.6, 7.5 Hz, 2H), 2.35 (s, 3H); **13C NMR** (126 MHz, CDCl₃) δ 138.21, 136.24, 135.38, 129.91, 128.96 (2C), 128.47, 128.14, 128.08, 127.44 (2C), 126.16, 120.38, 42.16, 39.81, 21.30; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₅N ([M+Na]⁺) 244.1097, found 244.1095.

3-(3-methoxyphenyl)-2-phenylpropanenitrile (**3ad**)



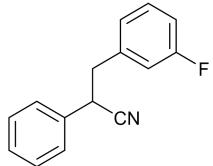
Yield: 92%; white solid; m.p. 37-39 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.42 – 7.33 (m, 3H), 7.33 – 7.27 (m, 2H), 7.24 (t, *J* = 7.9 Hz, 1H), 6.84 (dd, *J* = 8.3, 2.4 Hz, 1H), 6.77 (d, *J* = 7.5 Hz, 1H), 6.71 – 6.66 (m, 1H), 4.03 (dd, *J* = 8.3, 6.5 Hz, 1H), 3.77 (s, 3H), 3.16 (ddd, *J* = 20.0, 13.6, 7.4 Hz, 2H); **13C NMR** (126 MHz, CDCl₃) δ 159.62, 137.73, 135.19, 129.59, 128.98 (2C), 128.16, 127.45 (2C), 121.45, 120.35, 114.72, 112.92, 55.11, 42.15, 39.61; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₅NO ([M+Na]⁺) 260.1046, found 260.1048.

3-(4-methoxyphenyl)-2-phenylpropanenitrile (**3ae**)



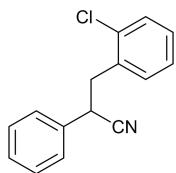
Yield: 96%; white solid; m.p. 84-85 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.41 – 7.32 (m, 3H), 7.29 – 7.25 (m, 2H), 7.10 – 7.03 (m, 2H), 6.88 – 6.82 (m, 2H), 3.98 (dd, *J* = 8.1, 6.5 Hz, 1H), 3.81 (s, 3H), 3.13 (qd, *J* = 13.7, 7.3 Hz, 2H); **13C NMR** (126 MHz, CDCl₃) δ 158.88, 135.28, 130.28 (2C), 128.97 (2C), 128.31, 128.13, 127.50 (2C), 120.45, 113.99 (2C), 55.22, 41.38, 40.04; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₅NO ([M+Na]⁺) 260.1046, found 260.1046.

3-(3-fluorophenyl)-2-phenylpropanenitrile (**3af**)



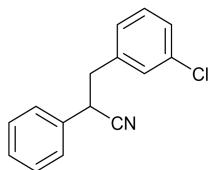
Yield: 90%; white solid; m.p. 46-48 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.41 – 7.35 (m, 3H), 7.31 – 7.25 (m, 3H), 7.03 – 6.92 (m, 2H), 6.88 – 6.81 (m, 1H), 4.03 (dd, *J* = 8.2, 6.5 Hz, 1H), 3.18 (qd, *J* = 13.7, 7.4 Hz, 2H); **13C NMR** (126 MHz, CDCl₃) δ 162.75 (d, *J*¹_{C-F} = 246.5 Hz, 1C), 138.59 (d, *J*³_{C-F} = 7.4 Hz, 1C), 134.81, 130.14 (d, *J*³_{C-F} = 8.3 Hz, 1C), 129.10 (2C), 128.37, 127.42 (2C), 124.93 (d, *J*⁴_{C-F} = 2.9 Hz, 1C), 120.05, 116.16 (d, *J*²_{C-F} = 21.4 Hz, 1C), 114.40 (d, *J*²_{C-F} = 21.0 Hz, 1C), 41.74 (d, *J*⁴_{C-F} = 1.6 Hz, 1C), 39.44; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₂FN ([M+Na]⁺) 248.0846, found 248.0844.

3-(2-chlorophenyl)-2-phenylpropanenitrile (**3ag**)



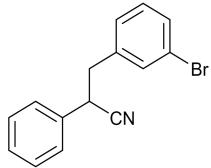
Yield: 94%; white solid; m.p. 66-67 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.45 – 7.34 (m, 6H), 7.29 – 7.22 (m, 3H), 4.20 (dd, *J* = 9.7, 6.1 Hz, 1H), 3.29 (ddd, *J* = 23.2, 13.6, 7.9 Hz, 2H); **13C NMR** (126 MHz, CDCl₃) δ 135.26, 134.17, 134.01, 131.81, 129.72, 129.11 (2C), 129.09, 128.30, 127.29 (2C), 127.15, 120.10, 40.44, 37.63; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₂ClN ([M+Na]⁺) 264.0550, found 264.0551.

3-(3-chlorophenyl)-2-phenylpropanenitrile (**3ah**)



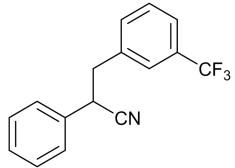
Yield: 97%; white solid; m.p. 67-68 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.43 – 7.34 (m, 3H), 7.31 – 7.24 (m, 4H), 7.15 – 7.11 (m, 1H), 7.08 – 7.02 (m, 1H), 4.02 (dd, *J* = 8.3, 6.5 Hz, 1H), 3.15 (qd, *J* = 13.7, 7.4 Hz, 2H); **13C NMR** (126 MHz, CDCl₃) δ 138.14, 134.76, 134.38, 129.88, 129.31, 129.11 (2C), 128.40, 127.65, 127.45, 127.43 (2C), 119.97, 41.69, 39.46; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₂ClN ([M+Na]⁺) 264.0550, found 264.0543.

3-(3-bromophenyl)-2-phenylpropanenitrile (**3ai**)



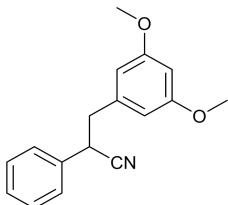
Yield: 85%; white solid; m.p. 83-84 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.45 – 7.34 (m, 4H), 7.31 – 7.25 (m, 3H), 7.22 – 7.15 (m, 1H), 7.12 – 7.07 (m, 1H), 4.01 (dd, *J* = 8.3, 6.5 Hz, 1H), 3.14 (qd, *J* = 13.7, 7.5 Hz, 2H); **13C NMR** (126 MHz, CDCl₃) δ 138.43, 134.74, 132.21, 130.57, 130.16, 129.11 (2C), 128.40, 127.92, 127.43 (2C), 122.55, 119.95, 41.64, 39.47; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₂BrN ([M+Na]⁺) 308.0045, found 308.0046.

2-phenyl-3-(3-(trifluoromethyl)phenyl)propanenitrile (**3aj**)



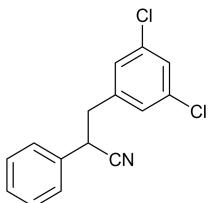
Yield: 85%; white solid; m.p. 65-66 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.55 (d, *J* = 7.8 Hz, 1H), 7.44 (t, *J* = 7.7 Hz, 1H), 7.40 – 7.34 (m, 4H), 7.31 (s, 1H), 7.28 – 7.24 (m, 2H), 4.09 – 4.01 (m, 1H), 3.30 – 3.19 (m, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 137.00, 134.52, 132.73, 130.92 (q, *J* = 32.4 Hz, 1C), 129.15 (2C), 129.12, 128.48, 127.45 (2C), 126.06 (q, *J* = 3.7 Hz, 1C), 124.32 (q, *J* = 3.8 Hz, 1C), 123.91 (q, *J* = 272.4 Hz, 1C), 119.85, 41.79, 39.40; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₂F₃N ([M+Na]⁺) 298.0814, found 298.0816.

3-(3,5-dimethoxyphenyl)-2-phenylpropanenitrile (**3ak**)



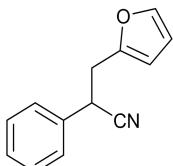
Yield: 83%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 7.42 – 7.27 (m, 5H), 6.39 (t, *J* = 2.2 Hz, 1H), 6.30 (d, *J* = 2.2 Hz, 2H), 4.01 (dd, *J* = 8.3, 6.5 Hz, 1H), 3.75 (s, 6H), 3.11 (ddd, *J* = 20.0, 13.5, 7.4 Hz, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 160.85 (2C), 138.49, 135.28, 129.04 (2C), 128.23, 127.52 (2C), 120.44, 107.20 (2C), 99.45, 55.30 (2C), 42.43, 39.57; **HRMS** (ES+) *m/z* calcd for C₁₇H₁₇NO₂ ([M+Na]⁺) 290.1151, found 290.1153.

3-(3,5-dichlorophenyl)-2-phenylpropanenitrile (**3al**)



Yield: 78%; white solid; m.p. 50-51 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.43 – 7.36 (m, 3H), 7.31 – 7.25 (m, 3H), 7.03 (d, *J* = 1.8 Hz, 2H), 4.02 (dd, *J* = 8.1, 6.6 Hz, 1H), 3.18 – 3.07 (m, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 139.32, 135.04 (2C), 134.31, 129.20 (2C), 128.58, 127.73 (3C), 127.36 (2C), 119.63, 41.32, 39.11; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₁Cl₂N ([M+Na]⁺) 298.0161, found 298.0154.

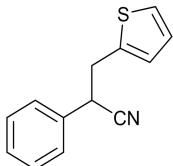
3-(furan-2-yl)-2-phenylpropanenitrile (**3am**)



Yield: 65%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 7.42 – 7.26 (m, 6H), 6.34 – 6.28 (m, 1H), 6.17 – 6.11 (m, 1H), 4.19 – 4.13 (m, 1H), 3.23 (ddd, *J* = 21.5, 15.0, 7.5 Hz, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 149.96, 142.14, 134.86, 129.08 (2C), 128.32,

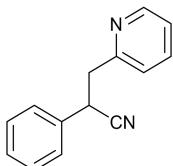
127.25 (2C), 120.07, 110.49, 108.20, 37.05, 34.62; **HRMS** (ES+) *m/z* calcd for C₁₃H₁₁NO ([M+Na]⁺) 220.0733, found 220.0732.

2-phenyl-3-(thiophen-2-yl)propanenitrile (**3an**)



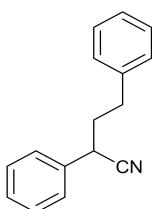
Yield: 76%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 7.42 – 7.30 (m, 5H), 7.20 (dd, *J* = 5.1, 1.1 Hz, 1H), 6.95 (dd, *J* = 5.1, 3.5 Hz, 1H), 6.90 – 6.85 (m, 1H), 4.06 (dd, *J* = 8.1, 6.4 Hz, 1H), 3.42 (ddd, *J* = 21.1, 14.7, 7.2 Hz, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 137.92, 134.74, 129.10 (2C), 128.41, 127.46 (2C), 127.10, 126.97, 124.91, 120.10, 40.06, 36.17; **HRMS** (ES+) *m/z* calcd for C₁₃H₁₁NS ([M+Na]⁺) 236.0504, found 236.0501.

2-phenyl-3-(pyridin-2-yl)propanenitrile (**3ao**)



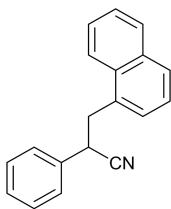
Yield: 83%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 8.61 (d, *J* = 4.8 Hz, 1H), 7.62 (td, *J* = 7.7, 1.7 Hz, 1H), 7.40 – 7.30 (m, 5H), 7.21 (dd, *J* = 7.5, 4.9 Hz, 1H), 7.13 (d, *J* = 7.7 Hz, 1H), 4.50 (dd, *J* = 9.2, 6.6 Hz, 1H), 3.33 (ddd, *J* = 20.4, 13.9, 7.9 Hz, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 156.16, 149.66, 136.72, 135.46, 129.07 (2C), 128.16, 127.39 (2C), 123.90, 122.35, 120.55, 44.15, 37.27; **HRMS** (ES+) *m/z* calcd for C₁₄H₁₂N₂ ([M+Na]⁺) 231.0893, found 231.0892.

2,4-diphenylbutanenitrile (**3ap**)



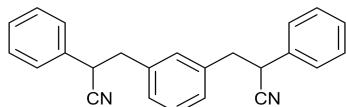
Yield: 81%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 7.45 – 7.40 (m, 2H), 7.38 – 7.32 (m, 5H), 7.31 – 7.20 (m, 3H), 3.77 (dd, *J* = 9.0, 6.1 Hz, 1H), 2.91 – 2.79 (m, 2H), 2.35 – 2.26 (m, 1H), 2.25 – 2.15 (m, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 139.77, 135.62, 129.10 (2C), 128.67 (2C), 128.41 (2C), 128.12, 127.26 (2C), 126.48, 120.62, 37.33, 36.53, 33.01; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₅N ([M+Na]⁺) 244.1097, found 244.1094.

3-(naphthalen-1-yl)-2-phenylpropanenitrile (**3aq**)



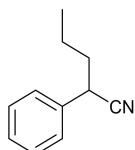
Yield: 90%; white solid; m.p. 84–85 °C; **1H NMR** (500 MHz, CDCl₃) δ 8.00 – 7.90 (m, 2H), 7.83 (d, *J* = 8.2 Hz, 1H), 7.62 – 7.51 (m, 2H), 7.45 – 7.31 (m, 7H), 4.20 (dd, *J* = 8.8, 6.6 Hz, 1H), 3.65 (qd, *J* = 14.0, 7.7 Hz, 2H); **13C NMR** (126 MHz, CDCl₃) δ 135.62, 133.97, 132.22, 131.31, 129.25, 129.12 (2C), 128.32, 128.29, 128.10, 127.35 (2C), 126.53, 125.78, 125.47, 122.60, 120.46, 39.56, 38.80; **HRMS** (ES+) *m/z* calcd for C₁₉H₁₅N ([M+Na]⁺) 280.1097, found 280.1097.

3,3'-(1,3-phenylene)bis(2-phenylpropanenitrile) (**3ar**)



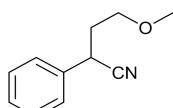
Yield: 85%; white solid; m.p. 145–146 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.42 – 7.33 (m, 10H), 7.25 – 7.20 (m, 1H), 7.18 (s, 1H), 7.15 – 7.11 (m, 2H), 4.51 – 4.43 (m, 2H), 3.19 – 3.07 (m, 4H). **13C NMR** (126 MHz, CDCl₃) δ 137.03, 136.99, 135.79 (2C), 130.01, 128.88 (4C), 128.24, 127.95 (2C), 127.82 (2C), 127.54 (4C), 120.92, 120.88, 40.39, 40.37, 37.84, 37.82. **HRMS** (ES+) *m/z* calcd for C₂₄H₂₀N₂ ([M+Na]⁺) 359.1519, found 359.1526.

2-phenylpentanenitrile (**3as**)



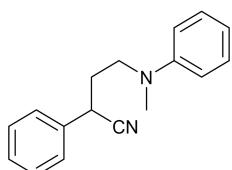
Yield: 68%; colorless oil; **1H NMR** (500 MHz, CDCl₃) δ 7.42 – 7.37 (m, 2H), 7.37 – 7.32 (m, 3H), 3.80 (dd, *J* = 8.6, 6.3 Hz, 1H), 1.99 – 1.80 (m, 2H), 1.61 – 1.43 (m, 2H), 0.98 (t, *J* = 7.4 Hz, 3H); **13C NMR** (126 MHz, CDCl₃) δ 136.03, 129.01 (2C), 127.95, 127.22 (2C), 120.89, 37.86, 37.14, 20.27, 13.39; **HRMS** (ES+) *m/z* calcd for C₁₁H₁₃N ([M+Na]⁺) 182.0940, found 182.0942.

4-methoxy-2-phenylbutanenitrile (**3at**)



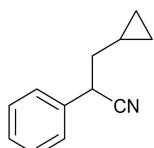
Yield: 85%; colorless oil; **1H NMR** (500 MHz, CDCl₃) δ 7.43 – 7.31 (m, 5H), 4.06 (dd, *J* = 8.8, 6.8 Hz, 1H), 3.56 (ddd, *J* = 9.8, 8.0, 4.4 Hz, 1H), 3.44 – 3.33 (m, 4H), 2.25 – 2.04 (m, 2H); **13C NMR** (126 MHz, CDCl₃) δ 135.55, 129.12 (2C), 128.12, 127.40 (2C), 120.76, 68.58, 58.81, 35.93, 33.80; **HRMS** (ES+) *m/z* calcd for C₁₁H₁₃NO ([M+Na]⁺) 198.0889, found 198.0885.

4-(methyl(phenyl)amino)-2-phenylbutanenitrile (3au**)**



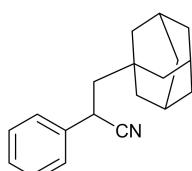
Yield: 82%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 7.45 – 7.40 (m, 2H), 7.39 – 7.35 (m, 3H), 7.30 – 7.24 (m, 2H), 6.78 (t, *J* = 7.3 Hz, 1H), 6.72 (d, *J* = 8.0 Hz, 2H), 3.89 (t, *J* = 7.4 Hz, 1H), 3.55 – 3.49 (m, 2H), 2.97 (s, 3H), 2.21 (dd, *J* = 14.5, 7.3 Hz, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 148.94, 135.33, 129.34 (2C), 129.22 (2C), 128.28, 127.25 (2C), 120.57, 117.13, 112.76 (2C), 50.05, 38.73, 34.81, 32.89; **HRMS** (ES+) *m/z* calcd for C₁₇H₁₈N₂ ([M+Na]⁺) 273.1362, found 273.1365.

3-cyclopropyl-2-phenylpropanenitrile (3av**)**



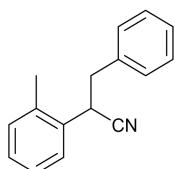
Yield: 80%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 7.41 – 7.31 (m, 5H), 3.88 (dd, *J* = 8.4, 6.3 Hz, 1H), 1.98 – 1.69 (m, 2H), 0.90 – 0.81 (m, 1H), 0.64 – 0.48 (m, 2H), 0.27 – 0.10 (m, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 135.96, 128.99 (2C), 127.98, 127.31 (2C), 121.03, 41.04, 37.84, 8.81, 4.75, 4.46; **HRMS** (ES+) *m/z* calcd for C₁₂H₁₃N ([M+Na]⁺) 194.0940, found 194.0936.

3-(adamantan-1-yl)-2-phenylpropanenitrile (3aw**)**



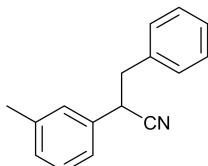
Yield: 64%; white solid; m.p. 49-50 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.41 – 7.29 (m, 5H), 3.82 (dd, *J* = 10.4, 3.3 Hz, 1H), 2.07 – 1.90 (m, 4H), 1.80 – 1.60 (m, 13H); **¹³C NMR** (126 MHz, CDCl₃) δ 138.08, 129.12 (2C), 127.76, 127.16 (2C), 122.36, 51.09, 42.21 (3C), 36.81 (3C), 33.03, 31.34, 28.48 (3C); **HRMS** (ES+) *m/z* calcd for C₁₉H₂₃N ([M+Na]⁺) 288.1723, found 288.1722.

3-phenyl-2-(o-tolyl)propanenitrile (3ba**)**



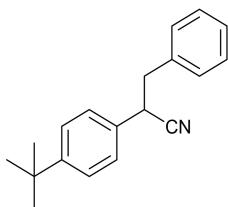
Yield: 75%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 7.48 – 7.44 (m, 1H), 7.37 – 7.25 (m, 5H), 7.23 – 7.18 (m, 3H), 4.18 (dd, *J* = 8.8, 5.9 Hz, 1H), 3.15 (ddd, *J* = 19.5, 13.6, 7.4 Hz, 2H), 2.29 (s, 3H); **¹³C NMR** (126 MHz, CDCl₃) δ 136.57, 135.06, 133.72, 130.96, 129.16 (2C), 128.71 (2C), 128.29, 127.70, 127.45, 126.91, 120.76, 41.01, 36.64, 19.06; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₅N ([M+Na]⁺) 244.1097, found 244.1092.

3-phenyl-2-(m-tolyl)propanenitrile (**3bb**)



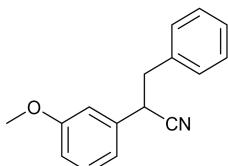
Yield: 92%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 7.36 – 7.25 (m, 4H), 7.21 – 7.14 (m, 3H), 7.13 (s, 1H), 7.09 (d, *J* = 7.7 Hz, 1H), 3.98 (dd, *J* = 8.6, 6.3 Hz, 1H), 3.17 (qd, *J* = 13.6, 7.5 Hz, 2H), 2.38 (s, 3H); **¹³C NMR** (126 MHz, CDCl₃) δ 138.88, 136.47, 135.21, 129.22 (2C), 128.94, 128.89, 128.62 (2C), 128.14, 127.37, 124.53, 120.48, 42.25, 39.81, 21.36; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₅N ([M+Na]⁺) 244.1097, found 244.1089.

2-(4-(tert-butyl)phenyl)-3-phenylpropanenitrile (**3bc**)



Yield: 84%; white solid; m.p. 66–67 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.43 – 7.39 (m, 2H), 7.37 – 7.28 (m, 3H), 7.28 – 7.19 (m, 4H), 4.00 (dd, *J* = 8.8, 6.3 Hz, 1H), 3.26 – 3.11 (m, 2H), 1.35 (s, 9H); **¹³C NMR** (126 MHz, CDCl₃) δ 151.30, 136.58, 132.30, 129.16 (2C), 128.63 (2C), 127.33, 127.09 (2C), 125.95 (2C), 120.50, 42.22, 39.47, 34.57, 31.27 (3C); **HRMS** (ES+) *m/z* calcd for C₁₉H₂₁N ([M+Na]⁺) 286.1566, found 286.1566.

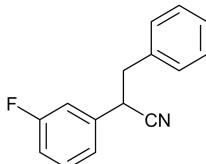
2-(3-methoxyphenyl)-3-phenylpropanenitrile (**3bd**)



Yield: 85%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 7.33 – 7.26 (m, 4H), 7.20 – 7.16 (m, 2H), 6.91 – 6.86 (m, 2H), 6.79 (t, *J* = 2.1 Hz, 1H), 3.99 (dd, *J* = 8.4, 6.4 Hz, 1H), 3.80 (s, 3H), 3.18 (qd, *J* = 13.6, 7.4 Hz, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 159.97, 136.64, 136.30, 130.06, 129.21 (2C), 128.62 (2C), 127.38, 120.27, 119.71,

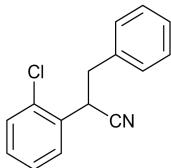
113.75, 113.17, 55.30, 42.08, 39.77; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₅NO ([M+Na]⁺) 260.1046, found 260.1044.

2-(3-fluorophenyl)-3-phenylpropanenitrile (**3be**)



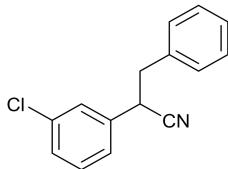
Yield: 62%; white solid; m.p. 64-65 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.37 – 7.27 (m, 4H), 7.18 – 7.13 (m, 2H), 7.10 – 7.02 (m, 2H), 7.02 – 6.97 (m, 1H), 4.02 (dd, *J* = 8.2, 6.5 Hz, 1H), 3.19 (ddd, *J* = 20.1, 13.6, 7.4 Hz, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 162.90 (d, *J*¹_{C-F} = 247.7 Hz, 1C), 137.48 (d, *J*³_{C-F} = 7.4 Hz, 1C), 135.82, 130.64 (d, *J*³_{C-F} = 8.4 Hz, 1C), 129.18 (2C), 128.72 (2C), 127.57, 123.24 (d, *J*⁴_{C-F} = 3.0 Hz, 1C), 119.81, 115.31 (d, *J*²_{C-F} = 21.1 Hz, 1C), 114.73 (d, *J*²_{C-F} = 22.7 Hz, 1C), 41.94, 39.45 (d, *J*⁴_{C-F} = 1.8 Hz); **HRMS** (ES+) *m/z* calcd for C₁₅H₁₂FN ([M+Na]⁺) 2248.0846, found 248.0838.

2-(2-chlorophenyl)-3-phenylpropanenitrile (**3bf**)



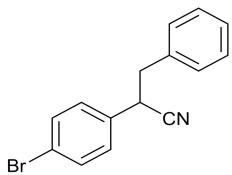
Yield: 60%; white solid; m.p. 65-66 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.52 – 7.43 (m, 2H), 7.37 – 7.30 (m, 5H), 7.28 – 7.24 (m, 2H), 4.55 (dd, *J* = 9.4, 4.8 Hz, 1H), 3.16 (ddd, *J* = 23.0, 13.6, 7.1 Hz, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 136.11, 133.04, 132.64, 130.03, 129.68, 129.25 (2C), 129.22, 128.69 (2C), 127.58, 127.53, 119.77, 40.14, 37.34; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₂ClN ([M+Na]⁺) 264.0550, found 264.0549.

2-(3-chlorophenyl)-3-phenylpropanenitrile (**3bg**)



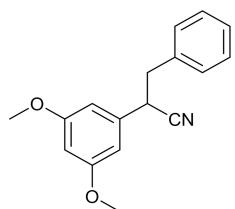
Yield: 92%; white solid; m.p. 64-65 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.36 – 7.26 (m, 6H), 7.18 – 7.13 (m, 3H), 4.00 (dd, *J* = 8.3, 6.5 Hz, 1H), 3.17 (ddd, *J* = 20.0, 13.6, 7.4 Hz, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 137.08, 135.79, 134.91, 130.29, 129.20 (2C), 128.73 (2C), 128.53, 127.72, 127.60, 125.74, 119.75, 41.97, 39.42; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₂ClN ([M+Na]⁺) 264.0550, found 264.0546.

2-(4-bromophenyl)-3-phenylpropanenitrile (**3bh**)



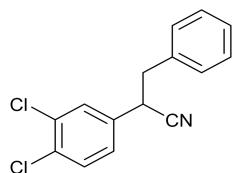
Yield: 90%; white solid; m.p. 67-68 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.52 – 7.47 (m, 2H), 7.35 – 7.27 (m, 3H), 7.16 – 7.09 (m, 4H), 3.99 (dd, *J* = 7.8, 6.8 Hz, 1H), 3.16 (ddd, *J* = 20.2, 13.6, 7.3 Hz, 2H); **13C NMR** (126 MHz, CDCl₃) δ 135.78, 134.16, 132.16 (2C), 129.24 (2C), 129.23 (2C), 128.72 (2C), 127.56, 122.32, 119.90, 41.95, 39.21; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₂BrN ([M+Na]⁺) 308.0045, found 308.0038.

2-(3,5-dimethoxyphenyl)-3-phenylpropanenitrile (**3bi**)



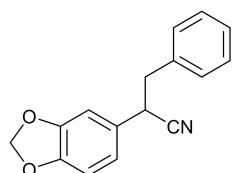
Yield: 85%; colorless oil; **1H NMR** (500 MHz, CDCl₃) δ 7.37 – 7.26 (m, 3H), 7.22 – 7.16 (m, 2H), 6.45 – 6.39 (m, 3H), 3.95 (dd, *J* = 8.3, 6.4 Hz, 1H), 3.77 (s, 6H), 3.17 (qd, *J* = 13.6, 7.4 Hz, 2H); **13C NMR** (126 MHz, CDCl₃) δ 161.16 (2C), 137.36, 136.35, 129.25 (2C), 128.65 (2C), 127.41, 120.26, 105.63 (2C), 100.08, 55.43 (2C), 41.99, 39.94; **HRMS** (ES+) *m/z* calcd for C₁₇H₁₇NO₂ ([M+Na]⁺) 290.1151, found 290.1153.

2-(3,4-dichlorophenyl)-3-phenylpropanenitrile (**3bj**)



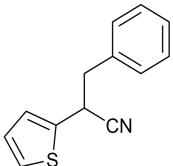
Yield: 82%; white solid; m.p. 88-89 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.44 (d, *J* = 8.3 Hz, 1H), 7.39 – 7.28 (m, 4H), 7.17 – 7.06 (m, 3H), 3.99 (dd, *J* = 7.8, 6.8 Hz, 1H), 3.16 (ddd, *J* = 20.2, 13.6, 7.3 Hz, 2H); **13C NMR** (126 MHz, CDCl₃) δ 135.43, 135.19, 133.20, 132.64, 130.95, 129.55, 129.22 (2C), 128.81 (2C), 127.73, 126.90, 119.45, 41.81, 38.88; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₁Cl₂N ([M+Na]⁺) 298.0161, found 298.0153.

2-(benzo[d][1,3]dioxol-5-yl)-3-phenylpropanenitrile (**3bk**)



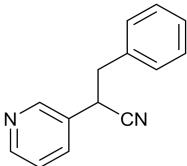
Yield: 83%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 7.35 – 7.26 (m, 3H), 7.20 – 7.14 (m, 2H), 6.80 – 6.76 (m, 2H), 6.71 (dd, *J* = 8.0, 1.7 Hz, 1H), 5.99 (s, 2H), 3.93 (dd, *J* = 8.2, 6.6 Hz, 1H), 3.15 (ddd, *J* = 20.1, 13.6, 7.4 Hz, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 148.15, 147.50, 136.22, 129.17 (2C), 128.83, 128.60 (2C), 127.36, 120.98, 120.40, 108.50, 107.82, 101.38, 42.20, 39.39; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₃NO₂ ([M+Na]⁺) 274.0838, found 274.0840.

3-phenyl-2-(thiophen-2-yl)propanenitrile (**3bl**)



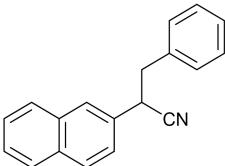
Yield: 55%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 7.35 – 7.27 (m, 4H), 7.25 – 7.21 (m, 2H), 7.03 – 6.96 (m, 2H), 4.34 – 4.28 (m, 1H), 3.33 – 3.22 (m, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 137.14, 135.91, 129.14 (2C), 128.71 (2C), 127.58, 127.06, 126.50, 125.65, 119.46, 42.19, 34.87; **HRMS** (ES+) *m/z* calcd for C₁₃H₁₁NS ([M+Na]⁺) 236.0504, found 236.0509.

3-phenyl-2-(pyridin-3-yl)propanenitrile (**3bm**)



Yield: 64%; white solid; m.p. 64–65 °C; **¹H NMR** (500 MHz, CDCl₃) δ 8.59 (dd, *J* = 4.8, 1.5 Hz, 1H), 8.47 (d, *J* = 2.3 Hz, 1H), 7.62 – 7.56 (m, 1H), 7.34 – 7.26 (m, 4H), 7.15 – 7.09 (m, 2H), 4.10 – 4.04 (m, 1H), 3.20 (ddd, *J* = 20.3, 13.6, 7.3 Hz, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 149.63, 148.81, 135.31, 135.03, 131.00, 129.22 (2C), 128.78 (2C), 127.70, 123.68, 119.39, 41.77, 37.21; **HRMS** (ES+) *m/z* calcd for C₁₄H₁₂N₂ ([M+Na]⁺) 231.0893, found 231.0894.

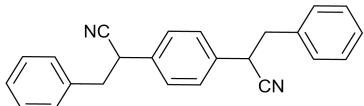
2-(naphthalen-2-yl)-3-phenylpropanenitrile (**3bn**)



Yield: 85%; white solid; m.p. 35–37 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.91 – 7.81 (m, 3H), 7.76 (s, 1H), 7.57 – 7.52 (m, 2H), 7.38 (dd, *J* = 8.5, 1.8 Hz, 1H), 7.35 – 7.26 (m, 3H), 7.23 – 7.17 (m, 2H), 4.20 (dd, *J* = 8.2, 6.5 Hz, 1H), 3.28 (qd, *J* = 13.7, 7.4 Hz, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 136.27, 133.24, 132.85, 132.50, 129.24 (2C), 128.98, 128.66 (2C), 127.90, 127.72, 127.42, 126.70, 126.64, 126.55, 124.95, 120.36,

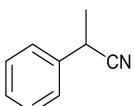
42.10, 39.97; **HRMS** (ES+) m/z calcd for $C_{19}H_{15}N$ ($[M+Na]^+$) 280.1097, found 280.1092.

2,2'-(1,4-phenylene)bis(3-phenylpropanenitrile) (**3bo**)



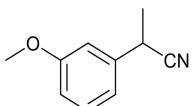
Yield: 88%; white solid; m.p. 147–149 °C; **1H NMR** (500 MHz, $CDCl_3$) δ 7.43 (s, 4H), 7.34 – 7.29 (m, 4H), 7.28 – 7.23 (m, 6H), 4.57 (t, $J = 7.7$ Hz, 2H), 3.17 (d, $J = 7.7$ Hz, 4H); **¹³C NMR** (126 MHz, $CDCl_3$) δ 136.90 (2C), 135.60 (2C), 129.19 (4C), 128.26 (4C), 128.11 (4C), 126.95 (2C), 120.81 (2C), 40.16 (2C), 37.43 (2C). **HRMS** (ES+) m/z calcd for $C_{24}H_{20}N_2$ ($[M+Na]^+$) 359.1519, found 359.1526.

2-phenylpropanenitrile (**3ca**)



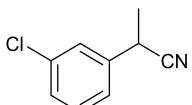
Yield: 46%; colorless oil; **1H NMR** (500 MHz, $CDCl_3$) δ 7.40 – 7.31 (m, 5H), 3.90 (q, $J = 7.3$ Hz, 1H), 1.64 (d, $J = 7.3$ Hz, 3H); **¹³C NMR** (126 MHz, $CDCl_3$) δ 137.06, 129.13 (2C), 128.03, 126.68 (2C), 121.56, 31.23, 21.44; **HRMS** (ES+) m/z calcd for C_9H_9N ($[M+Na]^+$) 154.0627, found 154.0624.

2-(3-methoxyphenyl)propanenitrile (**3cb**)



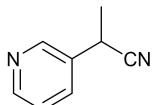
Yield: 57%; colorless oil; **1H NMR** (500 MHz, $CDCl_3$) δ 7.31 (t, $J = 7.9$ Hz, 1H), 6.98 – 6.86 (m, 3H), 3.88 (q, $J = 7.3$ Hz, 1H), 3.84 (s, 3H), 1.65 (d, $J = 7.3$ Hz, 3H); **¹³C NMR** (126 MHz, $CDCl_3$) δ 160.13, 138.55, 130.22, 121.52, 118.95, 113.36, 112.60, 55.32, 31.23, 21.38; **HRMS** (ES+) m/z calcd for $C_{10}H_{11}NO$ ($[M+Na]^+$) 184.0733, found 184.0737.

2-(3-chlorophenyl)propanenitrile (**3cc**)



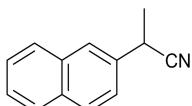
Yield: 55%; colorless oil; **1H NMR** (500 MHz, $CDCl_3$) δ 7.39 – 7.31 (m, 3H), 7.28 – 7.25 (m, 1H), 3.90 (q, $J = 7.3$ Hz, 1H), 1.65 (d, $J = 7.3$ Hz, 3H); **¹³C NMR** (126 MHz, $CDCl_3$) δ 138.92, 135.01, 130.47, 128.36, 127.01, 124.96, 120.93, 30.93, 21.27; **HRMS** (ES+) m/z calcd for C_9H_8ClN ($[M+Na]^+$) 188.0237, found 188.0235.

2-(pyridin-3-yl)propanenitrile (**3cd**)



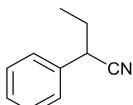
Yield: 45%; colorless oil; **¹H NMR** (500 MHz, CDCl₃) δ 8.77 – 8.56 (m, 2H), 7.76 (d, J = 7.9 Hz, 1H), 7.37 (dd, J = 7.8, 4.8 Hz, 1H), 3.97 (q, J = 7.3 Hz, 1H), 1.70 (d, J = 7.3 Hz, 3H); **¹³C NMR** (126 MHz, CDCl₃) δ 149.49, 148.14, 134.38, 132.94, 123.95, 120.48, 29.02, 21.22; **HRMS** (ES+) *m/z* calcd for C₈H₈N₂ ([M+Na]⁺) 155.0580, found 155.0578.

2-(naphthalen-2-yl)propanenitrile (**3ce**)



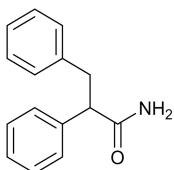
Yield: 63%; white solid; m.p. 60-61 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.92 – 7.84 (m, 4H), 7.58 – 7.51 (m, 2H), 7.45 (dd, J = 8.5, 1.8 Hz, 1H), 4.09 (q, J = 7.3 Hz, 1H), 1.75 (d, J = 7.3 Hz, 3H); **¹³C NMR** (126 MHz, CDCl₃) δ 134.35, 133.35, 132.80, 129.16, 127.87, 127.74, 126.75, 126.50, 125.60, 124.42, 121.59, 31.44, 21.43; **HRMS** (ES+) *m/z* calcd for C₁₃H₁₁N ([M+Na]⁺) 204.0784, found 204.0786.

2-phenylbutanenitrile (**3cf**)



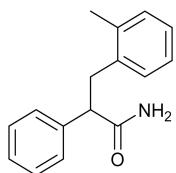
Yield: 54%; colorless oil °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.40 – 7.35 (m, 2H), 7.34 – 7.30 (m, 3H), 3.73 (dd, J = 9.5, 4.9 Hz, 1H), 1.98 – 1.90 (m, 2H), 1.07 (t, J = 7.4 Hz, 3H); **¹³C NMR** (126 MHz, CDCl₃) δ 135.72, 128.97 (2C), 127.97, 127.25 (2C), 120.71, 38.87, 29.16, 11.42; **HRMS** (ES+) *m/z* calcd for C₁₀H₁₁N ([M+Na]⁺) 168.0784, found 168.0788.

2,3-diphenylpropanamide (**4aa**)



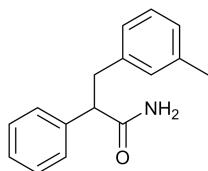
Yield: 90%; white solid; m.p. 127-128 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.35 – 7.26 (m, 5H), 7.25 – 7.20 (m, 2H), 7.19 – 7.15 (m, 1H), 7.13 – 7.09 (m, 2H), 5.69 (s, 1H), 5.41 (s, 1H), 3.66 (t, J = 7.5 Hz, 1H), 3.54 (dd, J = 13.7, 7.6 Hz, 1H), 3.01 (dd, J = 13.7, 7.4 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 175.24, 139.52, 139.40, 129.00 (2C), 128.82 (2C), 128.30 (2C), 128.06 (2C), 127.46, 126.25, 54.82, 39.37. **HRMS** (ES+) *m/z* calcd for C₁₅H₁₅NO ([M+Na]⁺) 248.1046, found 248.1050.

2-phenyl-3-(o-tolyl)propenamide (**4ab**)



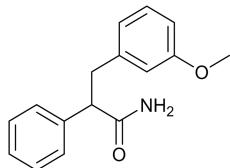
Yield: 83%; white solid; m.p. 149-150 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.33 – 7.26 (m, 5H), 7.14 – 7.01 (m, 4H), 5.65 (s, 1H), 5.39 (s, 1H), 3.64 (t, *J* = 7.3 Hz, 1H), 3.55 (dd, *J* = 13.9, 7.4 Hz, 1H), 3.01 (dd, *J* = 13.9, 7.1 Hz, 1H), 2.26 (s, 3H); **¹³C NMR** (126 MHz, CDCl₃) δ 175.27, 139.66, 137.68, 136.19, 130.21, 129.61, 128.78 (2C), 128.01 (2C), 127.42, 126.36, 125.82, 53.42, 36.58, 19.42; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₇NO ([M+Na]⁺) 262.1202, found 262.1207.

2-phenyl-3-(m-tolyl)propenamide (**4ac**)



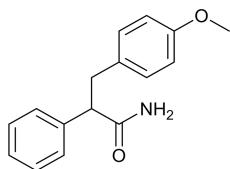
Yield: 77%; white solid; m.p. 107-108 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.34 – 7.26 (m, 5H), 7.14 – 7.09 (m, 1H), 7.02 – 6.89 (m, 3H), 5.67 (s, 1H), 5.41 (s, 1H), 3.65 (t, *J* = 7.5 Hz, 1H), 3.50 (dd, *J* = 13.7, 7.8 Hz, 1H), 2.97 (dd, *J* = 13.7, 7.1 Hz, 1H), 2.29 (s, 3H); **¹³C NMR** (126 MHz, CDCl₃) δ 175.27, 139.60, 139.48, 137.82, 129.80, 128.77 (2C), 128.17, 128.04 (2C), 127.39, 126.99, 125.95, 54.81, 39.35, 21.34; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₇NO ([M+Na]⁺) 262.1202, found 262.1208.

3-(3-methoxyphenyl)-2-phenylpropanamide (**4ad**)



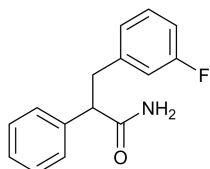
Yield: 86%; white solid; m.p. 97-98 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.33 – 7.25 (m, 5H), 7.14 (t, *J* = 7.9 Hz, 1H), 6.76 – 6.68 (m, 2H), 6.66 – 6.62 (m, 1H), 5.69 (s, 1H), 5.43 (s, 1H), 3.72 (s, 3H), 3.65 (t, *J* = 7.5 Hz, 1H), 3.51 (dd, *J* = 13.7, 7.5 Hz, 1H), 2.98 (dd, *J* = 13.7, 7.4 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 175.21, 159.48, 141.13, 139.46, 129.24, 128.80 (2C), 128.07 (2C), 127.44, 121.37, 114.60, 111.82, 55.08, 54.68, 39.40; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₇NO₂ ([M+Na]⁺) 278.1151, found 278.1156.

3-(4-methoxyphenyl)-2-phenylpropanamide (**4ae**)



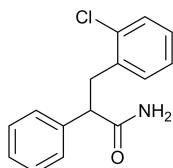
Yield: 94%; white solid; m.p. 157-158 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.34 – 7.25 (m, 5H), 7.05 – 7.00 (m, 2H), 6.79 – 6.73 (m, 2H), 5.51 (s, 1H), 5.35 (s, 1H), 3.76 (s, 3H), 3.61 (t, *J* = 7.5 Hz, 1H), 3.48 (dd, *J* = 13.8, 7.6 Hz, 1H), 2.96 (dd, *J* = 13.8, 7.3 Hz, 1H); **13C NMR** (126 MHz, CDCl₃) δ 175.20, 158.05, 139.47, 131.57, 129.94 (2C), 128.79 (2C), 128.06 (2C), 127.40, 113.71 (2C), 55.17, 55.11, 38.55; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₇NO₂ ([M+Na]⁺) 278.1151, found 278.1157.

3-(3-fluorophenyl)-2-phenylpropanamide (**4af**)



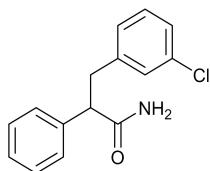
Yield: 85%; white solid; m.p. 133-134 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.34 – 7.25 (m, 5H), 7.20 – 7.14 (m, 1H), 6.91 – 6.78 (m, 3H), 5.65 (s, 1H), 5.40 (s, 1H), 3.64 (t, *J* = 7.5 Hz, 1H), 3.54 (dd, *J* = 13.7, 7.5 Hz, 1H), 2.99 (dd, *J* = 13.7, 7.5 Hz, 1H); **13C NMR** (126 MHz, CDCl₃) δ 174.87, 162.73 (d, *J*¹_{C-F} = 245.4 Hz, 1C), 142.05 (d, *J*³_{C-F} = 7.4 Hz, 1C), 139.04, 129.68 (d, *J*³_{C-F} = 8.3 Hz, 1C), 128.93 (2C), 128.01 (2C), 127.64, 124.71 (d, *J*⁴_{C-F} = 2.8 Hz, 1C), 115.83 (d, *J*²_{C-F} = 21.0 Hz, 1C), 113.17 (d, *J*²_{C-F} = 21.0 Hz, 1C), 54.51, 39.01; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₄FNO ([M+Na]⁺) 266.0952, found 266.0957.

3-(2-chlorophenyl)-2-phenylpropanamide (**4ag**)



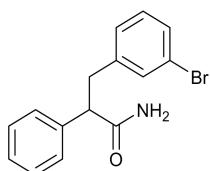
Yield: 83%; white solid; m.p. 174-175 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.33 – 7.31 (m, 4H), 7.29 – 7.27 (m, 2H), 7.15 – 7.06 (m, 3H), 5.52 (s, 1H), 5.36 (s, 1H), 3.84 (t, *J* = 7.4 Hz, 1H), 3.60 (dd, *J* = 13.6, 7.9 Hz, 1H), 3.14 (dd, *J* = 13.6, 7.0 Hz, 1H); **13C NMR** (126 MHz, CDCl₃) δ 174.81, 139.27, 136.92, 134.01, 131.85, 129.37, 128.82 (2C), 127.94 (2C), 127.87, 127.50, 126.64, 52.07, 37.52; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₄ClNO ([M+Na]⁺) 282.0656, found 282.0659.

3-(3-chlorophenyl)-2-phenylpropanamide (**4ah**)



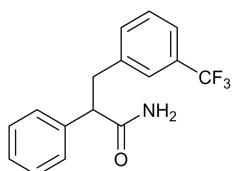
Yield: 95%; white solid; m.p. 121-122 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.35 – 7.25 (m, 5H), 7.16 – 7.09 (m, 3H), 7.00 – 6.95 (m, 1H), 5.74 (s, 1H), 5.44 (s, 1H), 3.63 (t, *J* = 7.5 Hz, 1H), 3.51 (dd, *J* = 13.7, 7.6 Hz, 1H), 2.96 (dd, *J* = 13.7, 7.4 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 174.85, 141.56, 138.99, 133.96, 129.50, 129.05, 128.91 (2C), 127.98 (2C), 127.62, 127.27, 126.46, 54.45, 38.95; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₄ClNO ([M+Na]⁺) 282.0656, found 282.0658.

3-(3-bromophenyl)-2-phenylpropanamide (4ai)



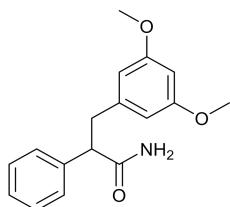
Yield: 65%; white solid; m.p. 106-107 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.33 – 7.25 (m, 7H), 7.09 – 7.05 (m, 1H), 7.04 – 6.99 (m, 1H), 5.80 (s, 1H), 5.44 (s, 1H), 3.62 (t, *J* = 7.5 Hz, 1H), 3.50 (dd, *J* = 13.7, 7.6 Hz, 1H), 2.95 (dd, *J* = 13.7, 7.4 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 174.92, 141.85, 138.94, 131.96, 129.81, 129.39, 128.92 (2C), 127.99 (2C), 127.74, 127.64, 122.26, 54.48, 38.92; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₄BrNO ([M+Na]⁺) 326.0151, found 326.0156.

2-phenyl-3-(3-(trifluoromethyl)phenyl)propanamide (4aj)



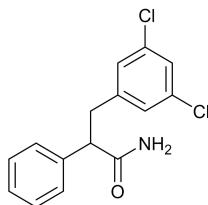
Yield: 89%; light yellow solid; m.p. 97-98 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.45 – 7.40 (m, 1H), 7.35 – 7.24 (m, 8H), 5.83 (s, 1H), 5.43 (s, 1H), 3.70 – 3.54 (m, 2H), 3.05 (dd, *J* = 13.4, 7.4 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 174.90, 140.38, 138.76, 132.52, 130.52 (q, *J* = 31.8 Hz, 1C), 128.97 (2C), 128.66, 128.03 (2C), 127.73, 125.70 (q, *J* = 3.7 Hz, 1C), 124.12 (q, *J* = 273.4 Hz, 1C), 123.14 (q, *J* = 3.8 Hz, 1C), 54.49, 39.08; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₄F₃NO ([M+Na]⁺) 316.0920, found 316.0925.

3-(3,5-dimethoxyphenyl)-2-phenylpropanamide (4ak)



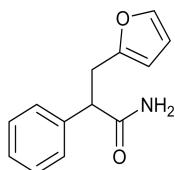
Yield: 87%; white solid; m.p. 115-116 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.35 – 7.26 (m, 5H), 6.30 – 6.23 (m, 3H), 5.84 (s, 1H), 5.46 (s, 1H), 3.70 (s, 6H), 3.64 (t, *J* = 7.5 Hz, 1H), 3.47 (dd, *J* = 13.7, 7.4 Hz, 1H), 2.95 (dd, *J* = 13.6, 7.5 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 175.45, 160.59 (2C), 141.84, 139.38, 128.82 (2C), 128.10 (2C), 127.47, 107.00 (2C), 98.45, 55.20 (2C), 54.61, 39.63; **HRMS** (ES+) *m/z* calcd for C₁₇H₁₉NO₃ ([M+Na]⁺) 308.1257, found 308.1263.

3-(3,5-dichlorophenyl)-2-phenylpropanamide (**4al**)



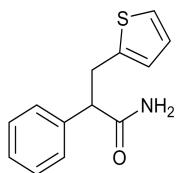
Yield: 64%; white solid; m.p. 123-124 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.36 – 7.25 (m, 5H), 7.19 – 7.15 (m, 1H), 6.99 (d, *J* = 1.7 Hz, 2H), 6.00 (s, 1H), 5.49 (s, 1H), 3.62 (t, *J* = 7.5 Hz, 1H), 3.49 (dd, *J* = 13.8, 7.6 Hz, 1H), 2.93 (dd, *J* = 13.8, 7.4 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 174.87, 142.84, 138.52, 134.63 (2C), 129.07 (2C), 127.95 (2C), 127.89, 127.56 (2C), 126.59, 54.16, 38.69; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₃Cl₂NO ([M+Na]⁺) 316.0266, found 316.0272.

3-(furan-2-yl)-2-phenylpropanamide (**4am**)



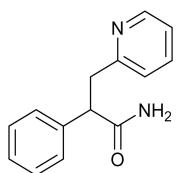
Yield: 80%; yellow solid; m.p. 100-101 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.35 – 7.24 (m, 6H), 6.21 (dd, *J* = 3.0, 1.9 Hz, 1H), 5.92 (d, *J* = 2.7 Hz, 1H), 5.84 (s, 1H), 5.52 (s, 1H), 3.83 (t, *J* = 7.6 Hz, 1H), 3.52 (dd, *J* = 15.2, 7.6 Hz, 1H), 3.06 (dd, *J* = 15.2, 7.5 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 175.02, 153.12, 141.13, 139.04, 128.84 (2C), 127.86 (2C), 127.56, 110.23, 106.56, 51.45, 31.70; **HRMS** (ES+) *m/z* calcd for C₁₃H₁₃NO₂ ([M+Na]⁺) 238.0838, found 238.0842.

2-phenyl-3-(thiophen-2-yl)propenamide (**4an**)



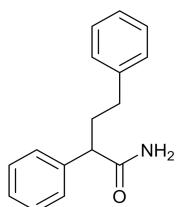
Yield: 87%; light yellow solid; m.p. 117-119 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.36 – 7.27 (m, 5H), 7.10 – 7.06 (m, 1H), 6.89 – 6.82 (m, 1H), 6.74 – 6.71 (m, 1H), 5.78 (s, 1H), 5.48 (s, 1H), 3.80 – 3.66 (m, 2H), 3.24 (dd, *J* = 14.1, 6.7 Hz, 1H); **13C NMR** (126 MHz, CDCl₃) δ 174.85, 141.73, 138.97, 128.89 (2C), 128.04 (2C), 127.65, 126.66, 125.71, 123.70, 54.87, 33.41; **HRMS** (ES+) *m/z* calcd for C₁₃H₁₃NOS ([M+Na]⁺) 254.0610, found 254.0616.

2-phenyl-3-(pyridin-2-yl)propanamide (**4ao**)



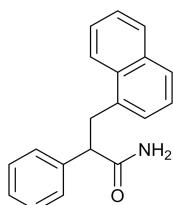
Yield: 75%; light grey solid; m.p. 162-163 °C; **1H NMR** (500 MHz, CDCl₃) δ 8.51 (dd, *J* = 4.8, 0.8 Hz, 1H), 7.56 – 7.48 (m, 1H), 7.38 – 7.22 (m, 5H), 7.13 – 7.03 (m, 2H), 5.83 (s, 1H), 5.45 (s, 1H), 4.18 (dd, *J* = 8.4, 6.6 Hz, 1H), 3.63 (dd, *J* = 14.1, 8.5 Hz, 1H), 3.17 (dd, *J* = 14.1, 6.6 Hz, 1H); **13C NMR** (126 MHz, CDCl₃) δ 175.13, 159.14, 149.04, 139.55, 136.37, 128.77 (2C), 127.98 (2C), 127.35, 124.16, 121.43, 51.88, 41.56; **HRMS** (ES+) *m/z* calcd for C₁₄H₁₄N₂O ([M+Na]⁺) 249.0998, found 249.0998.

2,4-diphenylbutanamide (**4ap**)



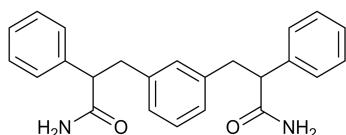
Yield: 74%; white solid; m.p. 95-96 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.39 – 7.35 (m, 2H), 7.34 – 7.27 (m, 5H), 7.23 – 7.15 (m, 3H), 5.88 (s, 1H), 5.43 (s, 1H), 3.38 (dd, *J* = 7.9, 7.1 Hz, 1H), 2.60 (t, *J* = 7.6 Hz, 2H), 2.52 (ddt, *J* = 14.0, 8.3, 6.9 Hz, 1H), 2.13 (ddd, *J* = 15.2, 13.5, 7.5 Hz, 1H); **13C NMR** (126 MHz, CDCl₃) δ 175.89, 141.37, 139.59, 128.95 (2C), 128.49 (2C), 128.36 (2C), 128.03 (2C), 127.45, 125.94, 51.69, 34.22, 33.44; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₇NO ([M+Na]⁺) 262.1202, found 262.1211.

3-(naphthalen-1-yl)-2-phenylpropanamide (**4aq**)



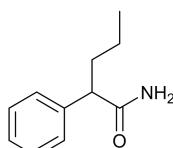
Yield: 85%; white solid; m.p. 151-152 °C; **¹H NMR** (500 MHz, CDCl₃) δ 8.07 (d, *J* = 8.4 Hz, 1H), 7.90 – 7.84 (m, 1H), 7.71 (d, *J* = 8.2 Hz, 1H), 7.57 – 7.47 (m, 2H), 7.36 – 7.26 (m, 6H), 7.18 (d, *J* = 6.8 Hz, 1H), 5.64 (s, 1H), 5.31 (s, 1H), 4.06 (dd, *J* = 14.0, 7.5 Hz, 1H), 3.83 (t, *J* = 7.2 Hz, 1H), 3.43 (dd, *J* = 14.0, 6.8 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 175.24, 139.70, 135.30, 133.91, 131.65, 128.96, 128.84 (2C), 127.97 (2C), 127.47, 127.43, 127.09, 126.04, 125.43 (2C), 123.47, 53.58, 36.51; **HRMS** (ES+) *m/z* calcd for C₁₉H₁₇NO ([M+Na]⁺) 298.1202, found 298.1210.

3,3'-(1,3-phenylene)bis(2-phenylpropanamide) (**4ar**)



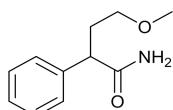
Yield: 83%; white solid; m.p. 143-144 °C; **¹H NMR** (500 MHz, DMSO) δ 7.44 – 7.40 (m, 2H), 7.39 – 7.34 (m, 4H), 7.33 – 7.27 (m, 4H), 7.25 – 7.20 (m, 2H), 7.10 – 7.03 (m, 2H), 6.97 – 6.93 (m, 2H), 6.80 (s, 2H), 3.74 (dd, *J* = 9.0, 6.2 Hz, 2H), 3.26 (dd, *J* = 13.6, 9.2 Hz, 2H), 2.78 (dd, *J* = 13.6, 6.0 Hz, 2H). **¹³C NMR** (126 MHz, DMSO) δ 174.46 (2C), 141.29 (2C), 140.10 (2C), 129.88, 128.63 (4C), 128.20 (5C), 127.05 (2C), 126.90 (2C), 53.15 (2C), 39.18 (2C). **HRMS** (ES+) *m/z* calcd for C₂₄H₂₄N₂O₂ ([M+Na]⁺) 395.1730, found 395.1735.

2-phenylpentanamide (**4as**)



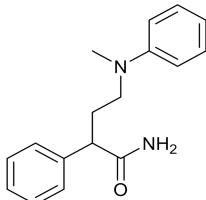
Yield: 80%; white solid; m.p. 85-86 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.36 – 7.26 (m, 5H), 5.92 (s, 1H), 5.47 (s, 1H), 3.46 – 3.37 (m, 1H), 2.12 (ddt, *J* = 13.1, 10.0, 6.4 Hz, 1H), 1.87 – 1.69 (m, 1H), 1.39 – 1.19 (m, 2H), 0.91 (t, *J* = 7.3 Hz, 3H); **¹³C NMR** (126 MHz, CDCl₃) δ 176.37, 139.99, 128.83 (2C), 127.94 (2C), 127.27, 52.56, 35.00, 20.78, 13.83; **HRMS** (ES+) *m/z* calcd for C₁₁H₁₅NO ([M+Na]⁺) 200.1046, found 200.1052.

4-methoxy-2-phenylbutanamide (**4at**)



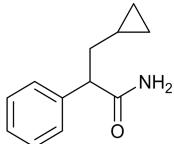
Yield: 81%; white solid; m.p. 86-87 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.35 – 7.30 (m, 4H), 7.29 – 7.25 (m, 1H), 5.89 (s, 1H), 5.66 (s, 1H), 3.65 (t, *J* = 7.6 Hz, 1H), 3.39 (ddd, *J* = 9.7, 6.3, 5.4 Hz, 1H), 3.32 – 3.22 (m, 4H), 2.48 – 2.34 (m, 1H), 2.03 – 1.93 (m, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 175.86, 139.59, 128.81 (2C), 128.00 (2C), 127.32, 69.95, 58.47, 48.65, 32.90; **HRMS** (ES+) *m/z* calcd for C₁₁H₁₅NO₂ ([M+Na]⁺) 216.0995, found 216.1001.

4-(methyl(phenyl)amino)-2-phenylbutanamide (**4au**)



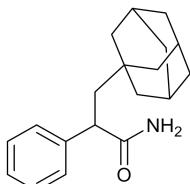
Yield: 81%; brown solid; m.p. 141-142 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.42 – 7.28 (m, 5H), 7.24 – 7.17 (m, 2H), 6.74 – 6.63 (m, 3H), 5.91 (s, 1H), 5.42 (s, 1H), 3.45 (t, *J* = 7.5 Hz, 1H), 3.38 – 3.25 (m, 2H), 2.90 (s, 3H), 2.56 – 2.43 (m, 1H), 2.09 – 1.97 (m, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 175.83, 149.35, 139.47, 129.16 (2C), 129.07 (2C), 127.94 (2C), 127.64, 116.37, 112.48 (2C), 50.74, 50.00, 38.11, 29.95; **HRMS** (ES+) *m/z* calcd for C₁₇H₂₀N₂O ([M+Na]⁺) 291.1468, found 291.1470.

3-cyclopropyl-2-phenylpropanamide (**4av**)



Yield: 82%; white solid; m.p. 105-106 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.37 – 7.31 (m, 4H), 7.29 – 7.25 (m, 1H), 6.04 (s, 1H), 5.53 (s, 1H), 3.52 (t, *J* = 7.5 Hz, 1H), 1.95 (dt, *J* = 14.4, 7.3 Hz, 1H), 1.76 (ddd, *J* = 14.1, 7.8, 6.7 Hz, 1H), 0.70 – 0.55 (m, 1H), 0.46 – 0.32 (m, 2H), 0.16 – 0.00 (m, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 176.42, 140.07, 128.80 (2C), 128.02 (2C), 127.28, 53.15, 38.13, 9.33, 4.70, 4.40; **HRMS** (ES+) *m/z* calcd for C₁₂H₁₅NO ([M+Na]⁺) 212.1046, found 212.1050.

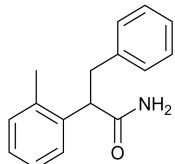
3-(adamantan-1-yl)-2-phenylpropanamide (**4aw**)



Yield: 64%; white solid; m.p. 144-145 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.33 – 7.29 (m, 4H), 7.27 – 7.23 (m, 1H), 5.70 (s, 1H), 5.50 (s, 1H), 3.55 (dd, *J* = 7.2, 5.3 Hz, 1H), 2.23 (dd, *J* = 14.2, 7.2 Hz, 1H), 1.95 – 1.90 (m, 3H), 1.71 – 1.64 (m, 3H), 1.63 – 1.57 (m, 3H), 1.54 – 1.43 (m, 7H); **¹³C NMR** (126 MHz, CDCl₃) δ 176.69, 142.15, 128.81

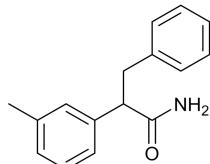
(2C), 127.77 (2C), 127.00, 47.38, 47.21, 42.47 (3C), 36.99 (3C), 32.95, 28.62 (3C); **HRMS** (ES+) m/z calcd for C₁₉H₂₅NO ([M+Na]⁺) 306.1828, found 306.1832.

3-phenyl-2-(o-tolyl)propenamide (**4ba**)



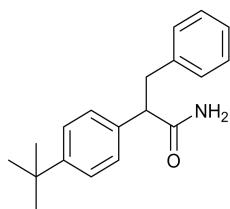
Yield: 68%; white solid; m.p. 130-131 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.25 – 7.05 (m, 9H), 5.74 (s, 1H), 5.29 (s, 1H), 3.93 (t, J = 7.4 Hz, 1H), 3.57 (dd, J = 13.7, 6.9 Hz, 1H), 2.97 (dd, J = 13.7, 7.9 Hz, 1H), 2.18 (s, 3H); **13C NMR** (126 MHz, CDCl₃) δ 175.52, 139.67, 137.66, 136.15, 130.67, 128.94 (2C), 128.24 (2C), 127.55, 127.29, 126.66, 126.19, 50.43, 38.84, 19.63; **HRMS** (ES+) m/z calcd for C₁₆H₁₇NO ([M+Na]⁺) 262.1202, found 262.1210.

3-phenyl-2-(m-tolyl)propenamide (**4bb**)



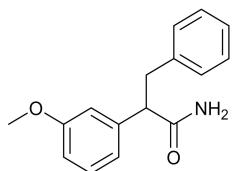
Yield: 75%; white solid; m.p. 93-94 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.27 – 7.06 (m, 9H), 5.75 (s, 1H), 5.44 (s, 1H), 3.63 (t, J = 7.4 Hz, 1H), 3.53 (dd, J = 13.6, 7.9 Hz, 1H), 2.99 (dd, J = 13.6, 7.0 Hz, 1H), 2.34 (s, 3H); **13C NMR** (126 MHz, CDCl₃) δ 175.43, 139.68, 139.41, 138.50, 129.00 (2C), 128.67 (2C), 128.29 (2C), 128.21, 126.22, 125.09, 54.73, 39.29, 21.41; **HRMS** (ES+) m/z calcd for C₁₆H₁₇NO ([M+Na]⁺) 262.1202, found 262.1209.

2-(4-(tert-butyl)phenyl)-3-phenylpropanamide (**4bc**)



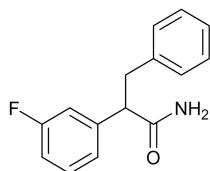
Yield: 88%; white solid; m.p. 121-122 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.37 – 7.33 (m, 2H), 7.29 – 7.21 (m, 5H), 7.21 – 7.12 (m, 2H), 5.67 (s, 1H), 5.40 (s, 1H), 3.71 – 3.61 (m, 1H), 3.54 (dd, J = 13.6, 8.2 Hz, 1H), 2.99 (dd, J = 13.6, 6.6 Hz, 1H), 1.32 (s, 9H); **13C NMR** (126 MHz, CDCl₃) δ 175.52, 150.37, 139.76, 136.39, 128.99 (2C), 128.29 (2C), 127.59 (2C), 126.21, 125.72 (2C), 54.38, 39.32, 34.47, 31.33 (3C); **HRMS** (ES+) m/z calcd for C₁₉H₂₃NO ([M+Na]⁺) 304.1672, found 304.1677.

2-(3-methoxyphenyl)-3-phenylpropanamide (**4bd**)



Yield: 92%; white solid; m.p. 84-85 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.16 – 7.00 (m, 6H), 6.80 – 6.69 (m, 3H), 5.61 (s, 1H), 5.35 (s, 1H), 3.68 (s, 3H), 3.53 (t, *J* = 7.5 Hz, 1H), 3.42 (dd, *J* = 13.7, 7.6 Hz, 1H), 2.90 (dd, *J* = 13.7, 7.3 Hz, 1H); **13C NMR** (126 MHz, CDCl₃) δ 175.08, 159.84, 140.94, 139.50, 129.77, 128.96 (2C), 128.27 (2C), 126.22, 120.39, 113.69, 112.85, 55.20, 54.73, 39.19; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₇NO₂ ([M+Na]⁺) 278.1151, found 278.1158.

2-(3-fluorophenyl)-3-phenylpropanamide (**4be**)



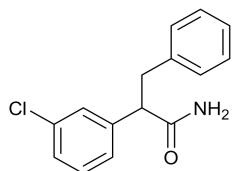
Yield: 80%; white solid; m.p. 141-142 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.32 – 7.21 (m, 4H), 7.14 – 7.03 (m, 4H), 7.00 – 6.92 (m, 1H), 5.66 (s, 1H), 5.39 (s, 1H), 3.64 (t, *J* = 7.5 Hz, 1H), 3.50 (dd, *J* = 13.7, 7.8 Hz, 1H), 2.99 (dd, *J* = 13.7, 7.3 Hz, 1H); **13C NMR** (126 MHz, CDCl₃) δ 174.54, 162.93 (d, *J*¹_{C-F} = 246.6 Hz, 1C), 141.73 (d, *J*³_{C-F} = 7.3 Hz, 1C), 139.01, 130.24 (d, *J*³_{C-F} = 8.3 Hz, 1C), 128.93 (2C), 128.40 (2C), 126.44, 123.75 (d, *J*⁴_{C-F} = 2.9 Hz, 1C), 115.01 (d, *J*²_{C-F} = 21.9 Hz, 1C), 114.44 (d, *J*²_{C-F} = 21.1 Hz, 1C), 54.53, 39.45; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₄FNO ([M+Na]⁺) 266.0952, found 266.0959.

2-(2-chlorophenyl)-3-phenylpropanamide (**4bf**)



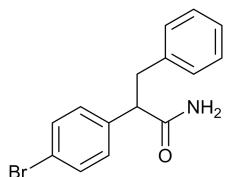
Yield: 66%; white solid; m.p. 124-125 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.60 – 7.55 (m, 1H), 7.39 – 7.35 (m, 1H), 7.31 – 7.18 (m, 7H), 5.71 (s, 1H), 5.47 (s, 1H), 4.29 (dd, *J* = 8.6, 6.2 Hz, 1H), 3.52 (dd, *J* = 13.7, 8.7 Hz, 1H), 3.01 (dd, *J* = 13.7, 6.1 Hz, 1H); **13C NMR** (126 MHz, CDCl₃) δ 174.08, 139.21, 137.00, 133.58, 129.62, 129.00 (2C), 128.98, 128.59, 128.39 (2C), 127.50, 126.40, 49.94, 38.29; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₄ClNO ([M+Na]⁺) 282.0656, found 282.0663.

2-(3-chlorophenyl)-3-phenylpropanamide (**4bg**)



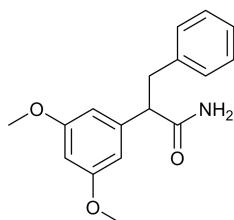
Yield: 84%; light yellow solid; m.p. 122-123 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.32 (s, 1H), 7.28 – 7.17 (m, 6H), 7.14 – 7.09 (m, 2H), 5.76 (s, 1H), 5.46 (s, 1H), 3.62 (t, *J* = 7.5 Hz, 1H), 3.48 (dd, *J* = 13.7, 8.0 Hz, 1H), 2.98 (dd, *J* = 13.7, 7.0 Hz, 1H); **13C NMR** (126 MHz, CDCl₃) δ 174.53, 141.34, 138.98, 134.55, 130.01, 128.95 (2C), 128.43 (2C), 128.16, 127.67, 126.48, 126.20, 54.44, 39.46; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₄ClNO ([M+Na]⁺) 282.0656, found 282.0664.

2-(4-bromophenyl)-3-phenylpropanamide (**4bh**)



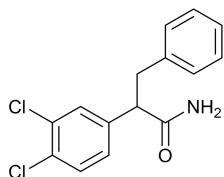
Yield: 68%; white solid; m.p. 169-170 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.47 – 7.42 (m, 2H), 7.26 – 7.16 (m, 5H), 7.13 – 7.08 (m, 2H), 5.48 (s, 1H), 5.33 (s, 1H), 3.60 (t, *J* = 7.5 Hz, 1H), 3.49 (dd, *J* = 13.6, 7.6 Hz, 1H), 2.97 (dd, *J* = 13.6, 7.4 Hz, 1H); **13C NMR** (126 MHz, CDCl₃) δ 174.46, 138.99, 138.29, 131.89 (2C), 129.74 (2C), 128.95 (2C), 128.41 (2C), 126.44, 121.46, 54.28, 39.51; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₄BrNO ([M+Na]⁺) 326.0151, found 326.0155.

2-(3,5-dimethoxyphenyl)-3-phenylpropanamide (**4bi**)



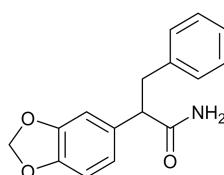
Yield: 77%; light yellow solid; m.p. 134-135 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.26 – 7.11 (m, 5H), 6.45 (d, *J* = 2.2 Hz, 2H), 6.36 (t, *J* = 2.2 Hz, 1H), 5.64 (s, 1H), 5.47 (s, 1H), 3.76 (s, 6H), 3.59 (t, *J* = 7.4 Hz, 1H), 3.50 (dd, *J* = 13.7, 7.5 Hz, 1H), 2.99 (dd, *J* = 13.7, 7.3 Hz, 1H); **13C NMR** (126 MHz, CDCl₃) δ 174.92, 161.01 (2C), 141.70, 139.51, 128.97 (2C), 128.30 (2C), 126.24, 106.17 (2C), 99.32, 55.34 (2C), 54.91, 39.04; **HRMS** (ES+) *m/z* calcd for C₁₇H₁₉NO₃ ([M+Na]⁺) 308.1257, found 308.1264.

2-(3,4-dichlorophenyl)-3-phenylpropanamide (**4bj**)



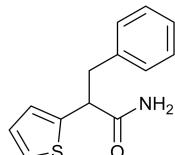
Yield: 77%; white solid; m.p. 128-129 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.43 – 7.36 (m, 2H), 7.30 – 7.19 (m, 3H), 7.17 – 7.09 (m, 3H), 5.68 (s, 1H), 5.44 (s, 1H), 3.59 (t, *J* = 7.5 Hz, 1H), 3.45 (dd, *J* = 13.7, 8.0 Hz, 1H), 2.96 (dd, *J* = 13.7, 7.1 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 174.06, 139.48, 138.61, 132.73, 131.59, 130.63, 129.96, 128.91 (2C), 128.52 (2C), 127.36, 126.62, 53.90, 39.61; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₃Cl₂NO ([M+Na]⁺) 316.0266, found 316.0271.

2-(benzo[d][1,3]dioxol-5-yl)-3-phenylpropanamide (**4bk**)



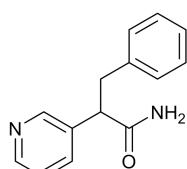
Yield: 77%; white solid; m.p. 139-140 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.26 – 7.21 (m, 2H), 7.20 – 7.15 (m, 1H), 7.14 – 7.09 (m, 2H), 6.86 – 6.83 (m, 1H), 6.75 – 6.67 (m, 2H), 5.95 (dd, *J* = 5.8, 1.4 Hz, 2H), 5.64 (s, 1H), 5.42 (s, 1H), 3.57 (t, *J* = 7.5 Hz, 1H), 3.48 (dd, *J* = 13.6, 7.4 Hz, 1H), 2.95 (dd, *J* = 13.6, 7.6 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 175.29, 147.99, 146.90, 139.43, 133.13, 128.96 (2C), 128.30 (2C), 126.25, 121.43, 108.35, 108.20, 101.09, 54.36, 39.39; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₅NO₃ ([M+Na]⁺) 292.0944, found 292.0946.

3-phenyl-2-(thiophen-2-yl)propenamide (**4bl**)



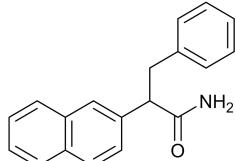
Yield: 67%; white solid; m.p. 122-123 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.29 – 7.14 (m, 6H), 6.94 (dt, *J* = 10.4, 3.2 Hz, 2H), 5.82 (s, 1H), 5.59 (s, 1H), 3.98 (t, *J* = 7.6 Hz, 1H), 3.51 (dd, *J* = 13.7, 7.5 Hz, 1H), 3.09 (dd, *J* = 13.7, 7.7 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 174.45, 141.65, 138.89, 128.95 (2C), 128.38 (2C), 126.87, 126.51, 125.76, 124.89, 49.97, 40.54; **HRMS** (ES+) *m/z* calcd for C₁₃H₁₃NOS ([M+Na]⁺) 254.0610, found 254.0613.

3-phenyl-2-(pyridin-3-yl)propenamide (**4bm**)



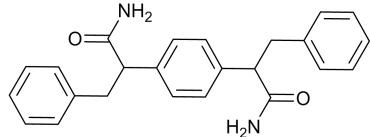
Yield: 62%; white solid; m.p. 187-188 °C; **¹H NMR** (500 MHz, DMSO) δ 8.49 (d, *J* = 1.9 Hz, 1H), 8.42 (dd, *J* = 4.7, 1.6 Hz, 1H), 7.79 (dt, *J* = 7.9, 1.9 Hz, 1H), 7.54 (s, 1H), 7.33 (dd, *J* = 7.8, 4.8 Hz, 1H), 7.26 – 7.12 (m, 5H), 6.92 (s, 1H), 3.82 (dd, *J* = 8.4, 7.1 Hz, 1H), 3.31 (dd, *J* = 13.6, 8.7 Hz, 1H), 2.90 (dd, *J* = 13.6, 6.8 Hz, 1H); **¹³C NMR** (126 MHz, DMSO) δ 173.91, 149.65, 148.40, 139.91, 136.51, 135.46, 129.31 (2C), 128.58 (2C), 126.54, 123.85, 50.65, 38.76; **HRMS** (ES+) *m/z* calcd for C₁₄H₁₄N₂O ([M+Na]⁺) 249.0998, found 249.0999.

2-(naphthalen-2-yl)-3-phenylpropanamide (**4bn**)



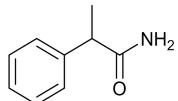
Yield: 77%; white solid; m.p. 143-144 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.86 – 7.76 (m, 3H), 7.72 (s, 1H), 7.53 – 7.44 (m, 3H), 7.25 – 7.12 (m, 5H), 5.68 (s, 1H), 5.45 (s, 1H), 3.83 (t, *J* = 7.5 Hz, 1H), 3.64 (dd, *J* = 13.8, 7.4 Hz, 1H), 3.12 (dd, *J* = 13.8, 7.5 Hz, 1H); **¹³C NMR** (126 MHz, CDCl₃) δ 175.15, 139.50, 136.85, 133.44, 132.70, 129.02 (2C), 128.65, 128.34 (2C), 127.80, 127.67, 127.05, 126.31, 126.28, 126.01, 125.87, 54.91, 39.29; **HRMS** (ES+) *m/z* calcd for C₁₉H₁₇NO ([M+Na]⁺) 298.1202, found 298.1207.

2,2'-(1,4-phenylene)bis(3-phenylpropanamide) (**4bo**)



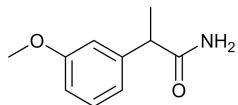
Yield: 82%; white solid; m.p. 156-158 °C; **¹H NMR** (500 MHz, DMSO) δ 7.41 (s, 2H), 7.30 (s, 4H), 7.26 – 7.14 (m, 10H), 6.77 (s, 2H), 3.73 (dd, *J* = 9.1, 5.9 Hz, 2H), 3.29 (dd, *J* = 13.2, 9.7 Hz, 2H), 2.83 (dd, *J* = 13.4, 5.4 Hz, 2H); **¹³C NMR** (126 MHz, DMSO) δ 174.42 (2C), 140.53 (2C), 139.60 (2C), 129.27 (4C), 128.52 (4C), 127.95 (4C), 126.39 (2C), 52.87 (2C), 39.08 (2C); **HRMS** (ES+) *m/z* calcd for C₂₄H₂₄N₂O₂ ([M+Na]⁺) 395.1730, found 395.1732.

2-phenylpropanamide (**4ca**)



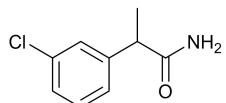
Yield: 71%; white solid; m.p. 88-89 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.38 – 7.26 (m, 5H), 6.05 (s, 1H), 5.45 (s, 1H), 3.60 (q, *J* = 7.2 Hz, 1H), 1.52 (d, *J* = 7.2 Hz, 3H); **¹³C NMR** (126 MHz, CDCl₃) δ 176.95, 141.25, 128.91 (2C), 127.56 (2C), 127.31, 46.57, 18.28; **HRMS** (ES+) *m/z* calcd for C₉H₁₁NO ([M+Na]⁺) 172.0733, found 172.0737.

2-(3-methoxyphenyl)propenamide (4cb**)**



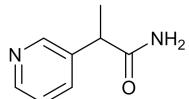
Yield: 57%; white solid; m.p. 83-84 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.33 – 7.22 (m, 1H), 6.95 – 6.77 (m, 3H), 5.84 (s, 1H), 5.43 (s, 1H), 3.81 (s, 3H), 3.58 (q, *J* = 7.2 Hz, 1H), 1.52 (d, *J* = 7.2 Hz, 3H); **13C NMR** (126 MHz, CDCl₃) δ 176.61, 159.99, 142.83, 129.97, 119.90, 113.37, 112.63, 55.22, 46.63, 18.19; **HRMS** (ES+) *m/z* calcd for C₁₀H₁₃NO₂ ([M+Na]⁺) 202.0838, found 202.0841.

2-(3-chlorophenyl)propenamide (4cc**)**



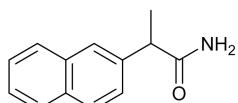
Yield: 62%; white solid; m.p. 82-83 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.35 – 7.18 (m, 4H), 5.94 (s, 1H), 5.47 (s, 1H), 3.57 (q, *J* = 7.2 Hz, 1H), 1.51 (d, *J* = 7.2 Hz, 3H); **13C NMR** (126 MHz, CDCl₃) δ 175.98, 143.19, 134.69, 130.18, 127.78, 127.54, 125.77, 46.27, 18.36; **HRMS** (ES+) *m/z* calcd for C₉H₁₀ClNO ([M+Na]⁺) 206.343, found 206.348.

2-(pyridin-3-yl)propenamide (4cd**)**



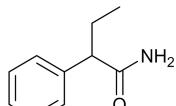
Yield: 68%; white solid; m.p. 98-99 °C; **1H NMR** (500 MHz, CDCl₃) δ 8.49 (s, 2H), 7.71 (d, *J* = 7.8 Hz, 1H), 7.37 – 7.18 (m, 1H), 6.12 (s, 2H), 3.61 (q, *J* = 7.1 Hz, 1H), 1.50 (d, *J* = 7.1 Hz, 3H); **13C NMR** (126 MHz, CDCl₃) δ 175.89, 149.00, 148.50, 137.02, 134.97, 123.81, 43.84, 18.59; **HRMS** (ES+) *m/z* calcd for C₈H₁₀N₂O ([M+Na]⁺) 173.0685, found 173.0688.

2-(naphthalen-2-yl)propenamide (4ce**)**



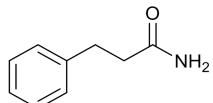
Yield: 45%; white solid; m.p. 125-126 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.92 – 7.73 (m, 4H), 7.58 – 7.40 (m, 3H), 5.85 (s, 1H), 5.44 (s, 1H), 3.78 (q, *J* = 7.2 Hz, 1H), 1.62 (d, *J* = 7.2 Hz, 3H); **13C NMR** (126 MHz, CDCl₃) δ 176.66, 138.74, 133.53, 132.64, 128.79, 127.73, 127.69, 126.41, 126.28, 126.02, 125.67, 46.75, 18.32; **HRMS** (ES+) *m/z* calcd for C₁₃H₁₃NO ([M+Na]⁺) 222.0889, found 222.0891.

2-phenylbutanamide (4af**)**



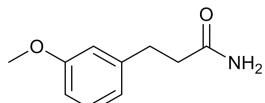
Yield: 72%; white solid; m.p. 74-75 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.37 – 7.26 (m, 5H), 5.98 (s, 1H), 5.51 (s, 1H), 3.29 (t, *J* = 7.6 Hz, 1H), 2.25 – 2.11 (m, 1H), 1.86 – 1.74 (m, 1H), 0.90 (t, *J* = 7.4 Hz, 3H); **¹³C NMR** (126 MHz, CDCl₃) δ 176.31, 139.80, 128.81 (2C), 127.96 (2C), 127.28, 54.53, 26.09, 12.25; **HRMS** (ES+) *m/z* calcd for C₁₀H₁₃NO ([M+Na]⁺) 186.0889, found 186.0893.

3-phenylpropanamide (**4da**)



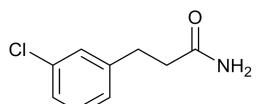
Yield: 45%; white solid; m.p. 100-101 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.34 – 7.16 (m, 5H), 6.04 (s, 1H), 5.63 (s, 1H), 3.00 – 2.91 (m, 2H), 2.58 – 2.47 (m, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 175.01, 140.69, 128.57 (2C), 128.31 (2C), 126.30, 37.50, 31.40; **HRMS** (ES+) *m/z* calcd for C₉H₁₁NO ([M+Na]⁺) 172.0733, found 172.0734.

3-(3-methoxyphenyl)propenamide (**4db**)



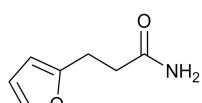
Yield: 52%; white solid; m.p. 105-107 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.25 – 7.17 (m, 1H), 6.87 – 6.71 (m, 3H), 5.84 (s, 1H), 5.53 (s, 1H), 3.80 (s, 3H), 2.99 – 2.89 (m, 2H), 2.58 – 2.47 (m, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 174.70, 159.75, 142.32, 129.57, 120.64, 114.11, 111.59, 55.16, 37.40, 31.43; **HRMS** (ES+) *m/z* calcd for C₁₀H₁₃NO₂ ([M+Na]⁺) 202.0838, found 202.0841.

3-(3-chlorophenyl)propenamide (**4dc**)



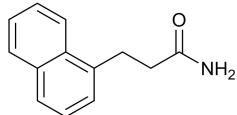
Yield: 51%; white solid; m.p. 103-104 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.26 – 7.16 (m, 3H), 7.13 – 7.06 (m, 1H), 5.83 (s, 1H), 5.52 (s, 1H), 3.02 – 2.90 (m, 2H), 2.57 – 2.48 (m, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 174.16, 142.74, 134.25, 129.81, 128.45, 126.59, 126.50, 37.05, 30.89; **HRMS** (ES+) *m/z* calcd for C₉H₁₀ClNO ([M+Na]⁺) 206.0343, found 206.0343.

3-(furan-2-yl)propenamide (**4dd**)



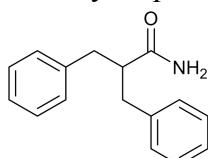
Yield: 42%; white solid; m.p. 96-97 °C; **¹H NMR** (500 MHz, CDCl₃) δ 7.32 (d, *J* = 1.1 Hz, 1H), 6.30 (dd, *J* = 3.1, 1.9 Hz, 1H), 6.06 (dd, *J* = 3.1, 0.7 Hz, 1H), 5.49 (s, 2H), 3.13 – 2.93 (m, 2H), 2.65 – 2.53 (m, 2H); **¹³C NMR** (126 MHz, CDCl₃) δ 173.98, 154.17, 141.26, 110.29, 105.55, 34.11, 23.75; **HRMS** (ES+) *m/z* calcd for C₇H₉NO₂ ([M+Na]⁺) 162.0525, found 162.0526.

3-(naphthalen-1-yl)propenamide (4de**)**



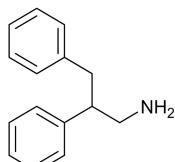
Yield: 54%; white solid; m.p. 95-96 °C; **1H NMR** (500 MHz, CDCl₃) δ 8.10 – 8.03 (m, 1H), 7.92 – 7.85 (m, 1H), 7.80 – 7.71 (m, 1H), 7.58 – 7.47 (m, 2H), 7.45 – 7.34 (m, 2H), 5.70 (s, 1H), 5.43 (s, 1H), 3.49 – 3.33 (m, 2H), 2.71 – 2.58 (m, 2H); **13C NMR** (126 MHz, CDCl₃) δ 174.65, 136.69, 133.92, 131.54, 128.92, 127.14, 126.13, 126.12, 125.64, 125.62, 123.42, 36.70, 28.46; **HRMS** (ES+) *m/z* calcd for C₁₃H₁₃NO ([M+Na]⁺) 222.0889, found 222.0892.

2-benzyl-3-phenylpropanamide (4df**)**



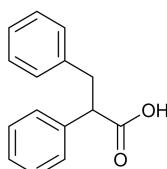
Yield: 58%; white solid; m.p. 121-122 °C; **1H NMR** (500 MHz, CDCl₃) δ 7.32 – 7.27 (m, 5H), 7.25 – 7.20 (m, 3H), 7.18 – 7.14 (m, 2H), 5.77 (s, 1H), 4.40 (d, *J* = 5.7 Hz, 2H), 3.01 (t, *J* = 7.6 Hz, 2H), 2.53 (t, *J* = 7.6 Hz, 2H); **13C NMR** (126 MHz, CDCl₃) δ 172.03, 140.75, 138.11, 128.66 (2C), 128.56 (2C), 128.41 (2C), 127.74 (2C), 127.46, 126.27, 43.59, 38.48, 31.73; **HRMS** (ES+) *m/z* calcd for C₁₆H₁₇NO ([M+Na]⁺) 262.1202, found 262.1209.

2,3-diphenylpropan-1-amine (5**)**



white solid; **1H NMR** (500 MHz, CDCl₃) δ 7.33 – 7.28 (m, 2H), 7.25 – 7.20 (m, 3H), 7.20 – 7.14 (m, 3H), 7.10 – 7.05 (m, 2H), 3.03 – 2.88 (m, 5H), 1.94 (s, 2H); **13C NMR** (126 MHz, CDCl₃) δ 142.67, 140.03, 129.04 (2C), 128.56 (2C), 128.18 (2C), 128.02 (2C), 126.65, 125.95, 51.05, 46.73, 40.77; **HRMS** (ES+) *m/z* calcd for C₁₅H₁₇N ([M+H]⁺) 212.1434, found 212.1433.

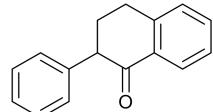
2,3-diphenylpropanoic acid (6**)**



white solid; **1H NMR** (500 MHz, CDCl₃) δ 7.31 – 7.28 (m, 4H), 7.28 – 7.14 (m, 4H), 7.11 – 7.07 (m, 2H), 3.85 (dd, *J* = 8.1, 7.3 Hz, 1H), 3.40 (dd, *J* = 13.8, 8.4 Hz, 1H),

3.02 (dd, $J = 13.8, 7.0$ Hz, 1H); **^{13}C NMR** (126 MHz, CDCl_3) δ 179.30, 138.68, 137.93, 128.93 (2C), 128.72 (2C), 128.39 (2C), 128.12 (2C), 127.64, 126.47, 53.46, 39.27; **HRMS** (ES+) m/z calcd for $\text{C}_{15}\text{H}_{14}\text{O}_2$ ($[\text{M}+\text{Na}]^+$) 249.0886, found 249.0887.

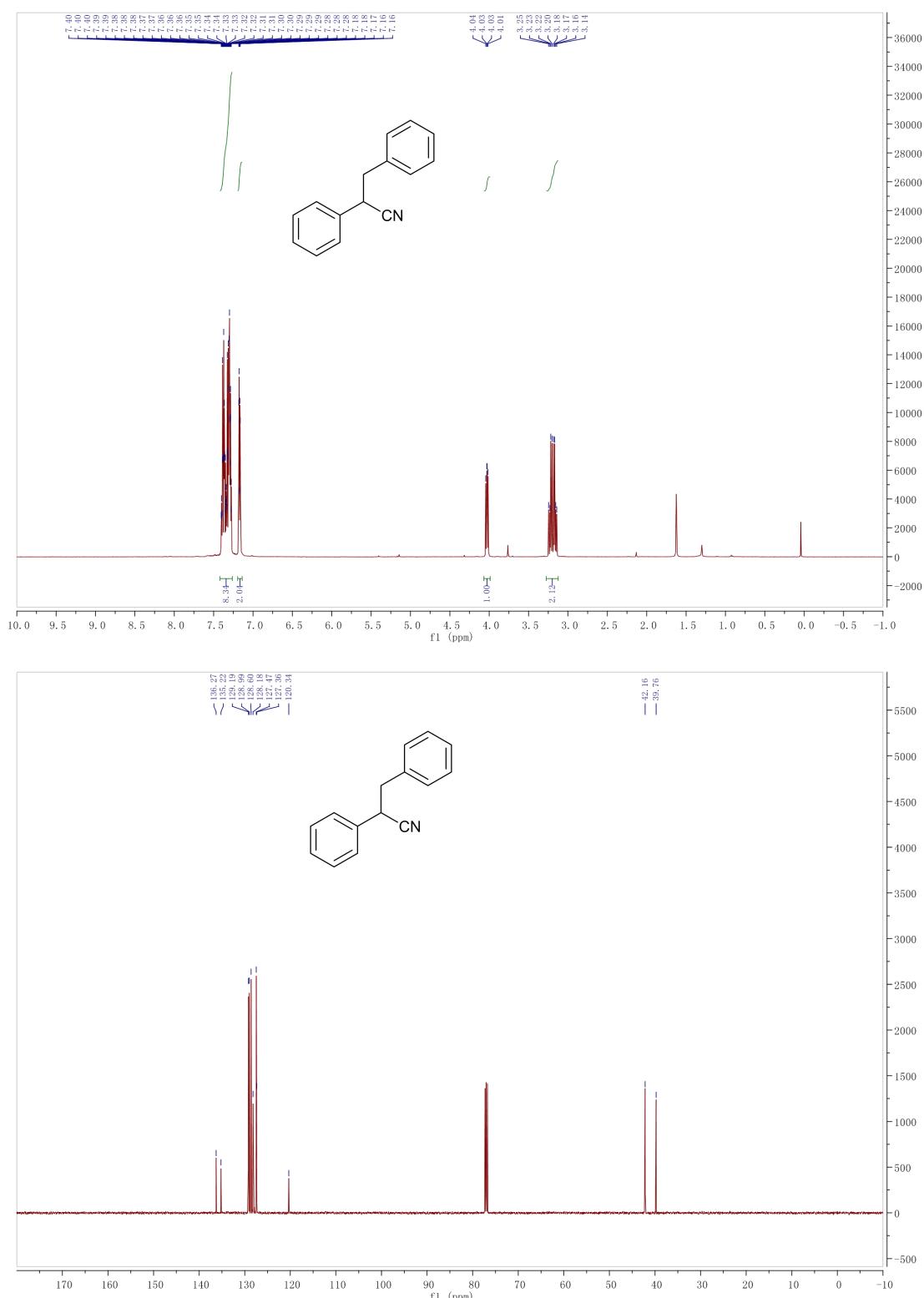
2-phenyl-3,4-dihydronaphthalen-1(2H)-one (7)



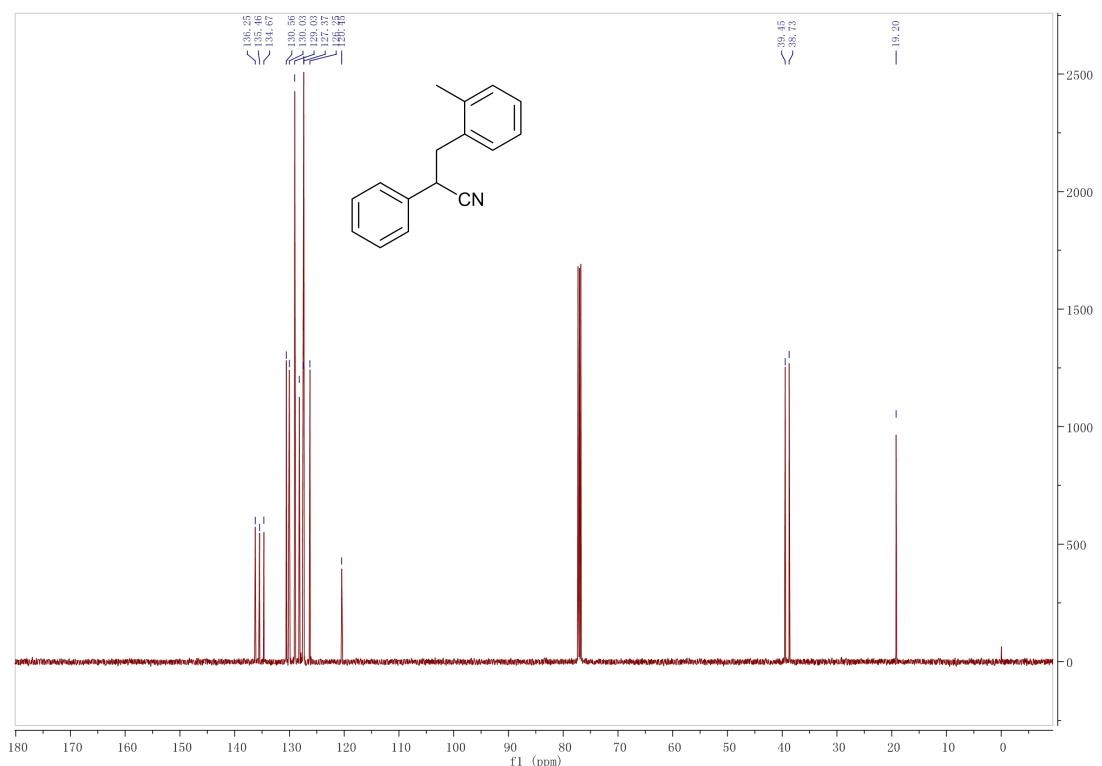
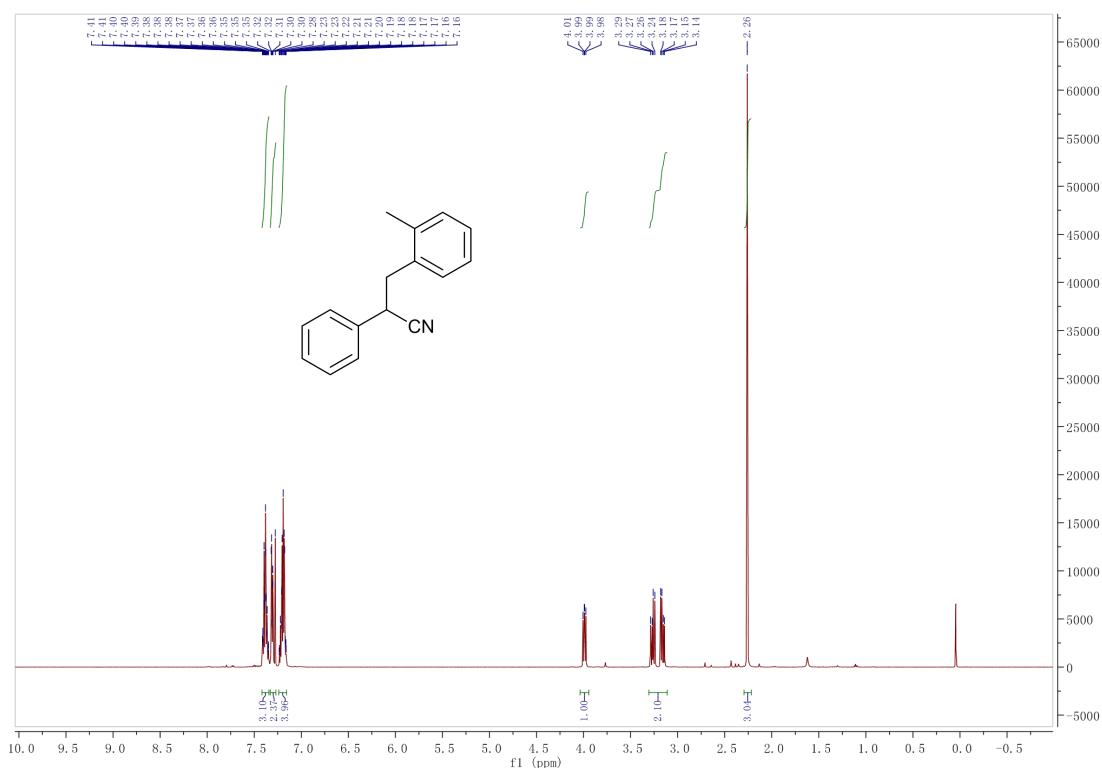
white solid; **^1H NMR** (500 MHz, CDCl_3) δ 8.13 (dd, $J = 7.8, 0.8$ Hz, 1H), 7.57 – 7.50 (m, 1H), 7.41 – 7.34 (m, 3H), 7.33 – 7.27 (m, 2H), 7.25 – 7.18 (m, 2H), 3.83 (dd, $J = 9.2, 6.7$ Hz, 1H), 3.22 – 3.01 (m, 2H), 2.54 – 2.41 (m, 2H); **^{13}C NMR** (126 MHz, CDCl_3) δ 198.21, 144.06, 139.73, 133.43, 132.86, 128.77, 128.52 (2C), 128.42 (2C), 127.79, 126.92, 126.77, 54.39, 31.17, 28.76; **HRMS** (ES+) m/z calcd for $\text{C}_{16}\text{H}_{14}\text{O}$ ($[\text{M}+\text{Na}]^+$) 245.0937, found 245.0941.

6. ^1H , ^{13}C NMR spectra

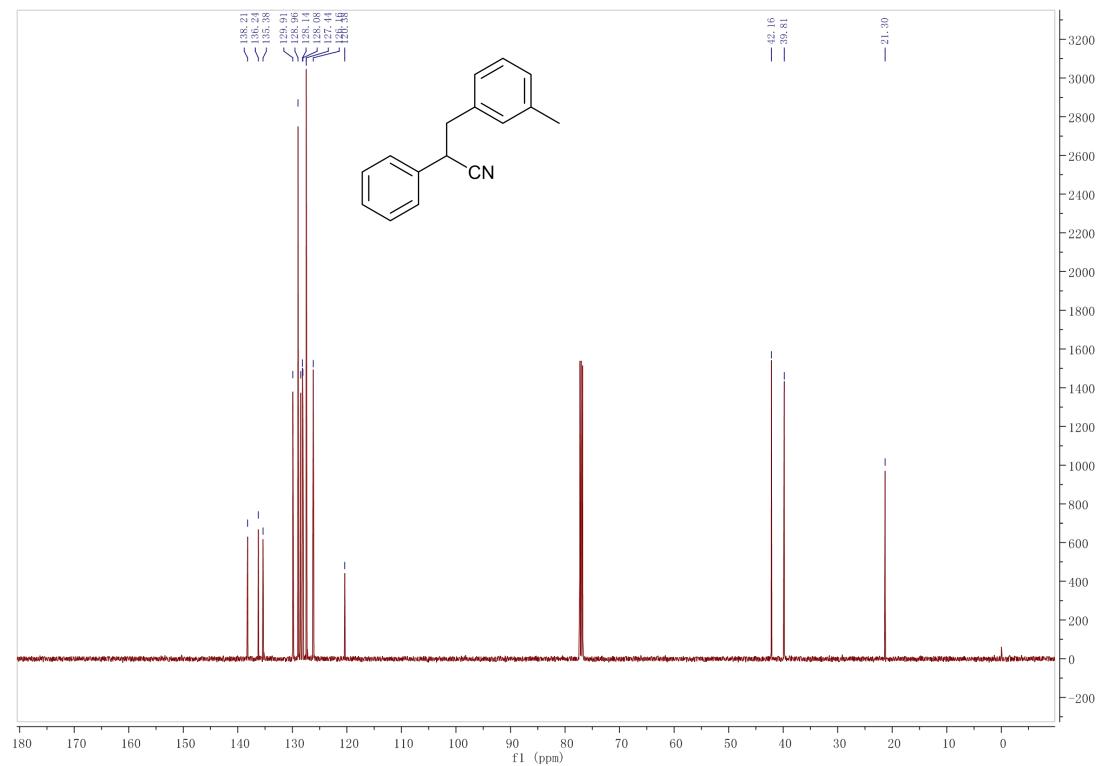
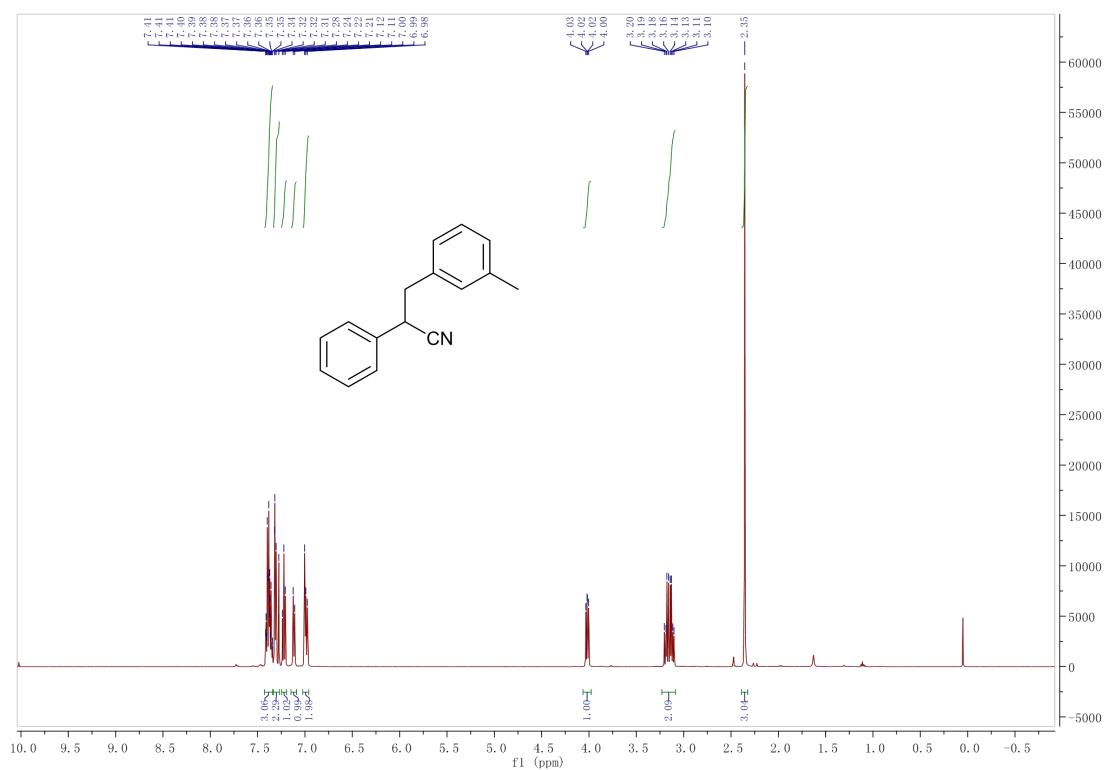
2,3-diphenylpropanenitrile (**3aa**)



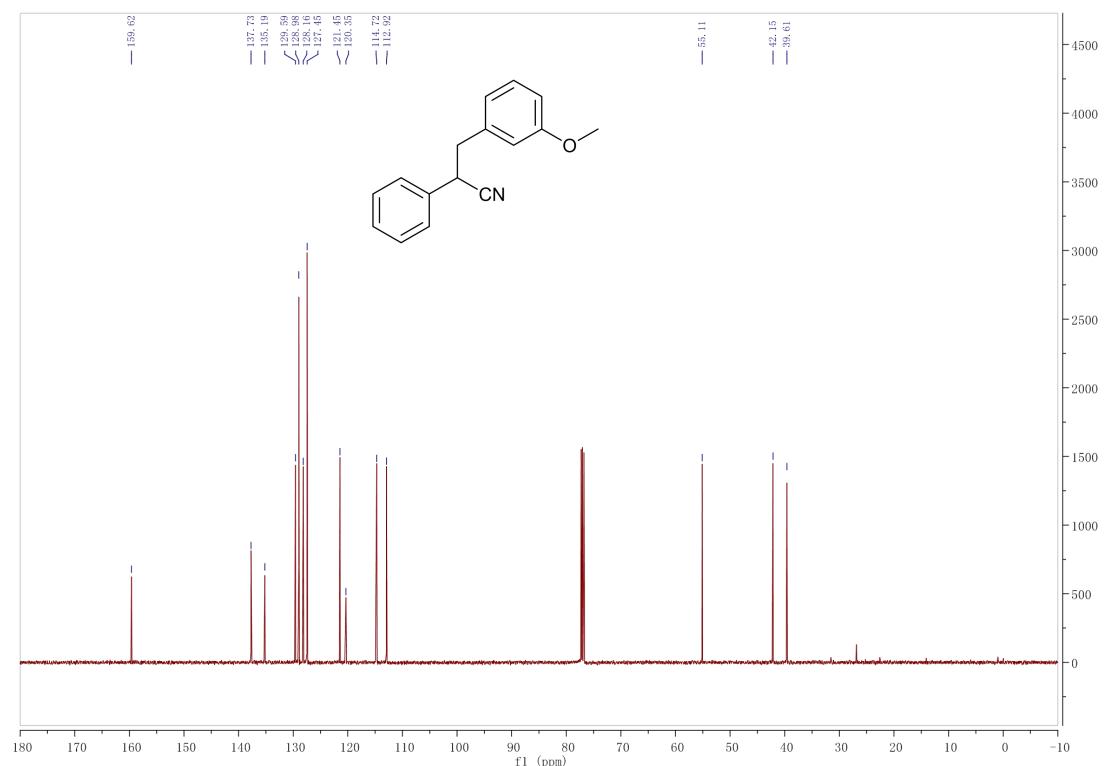
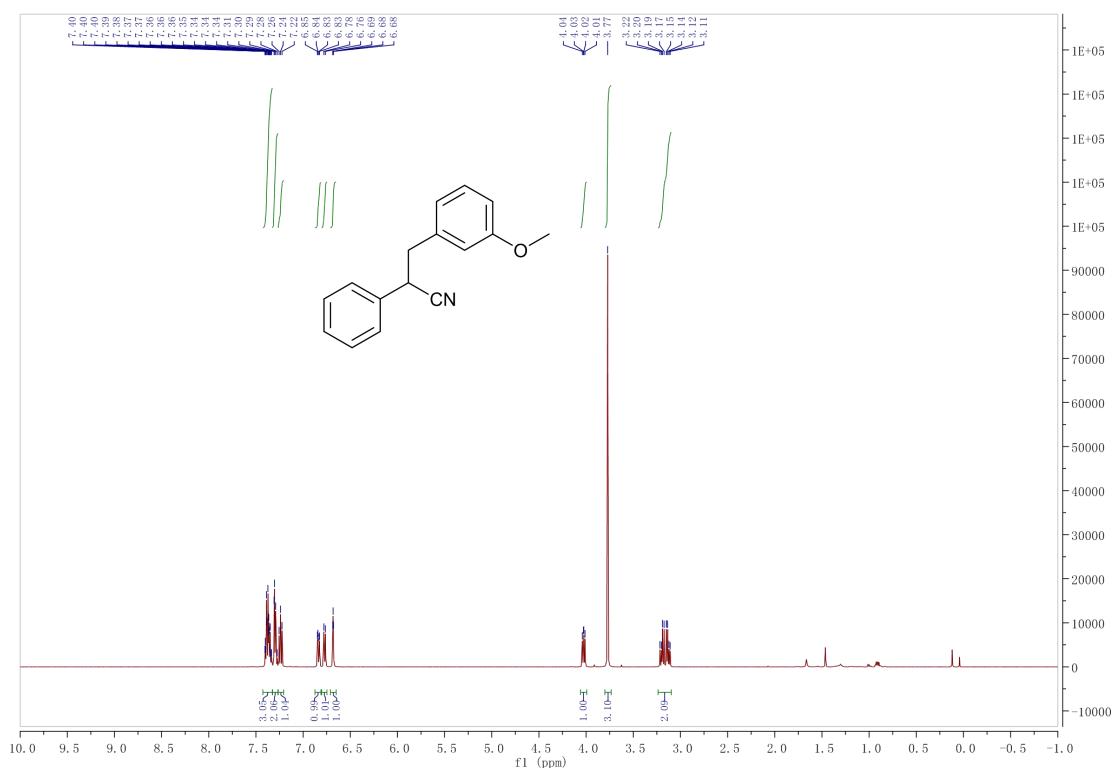
2-phenyl-3-(o-tolyl)propanenitrile (**3ab**)



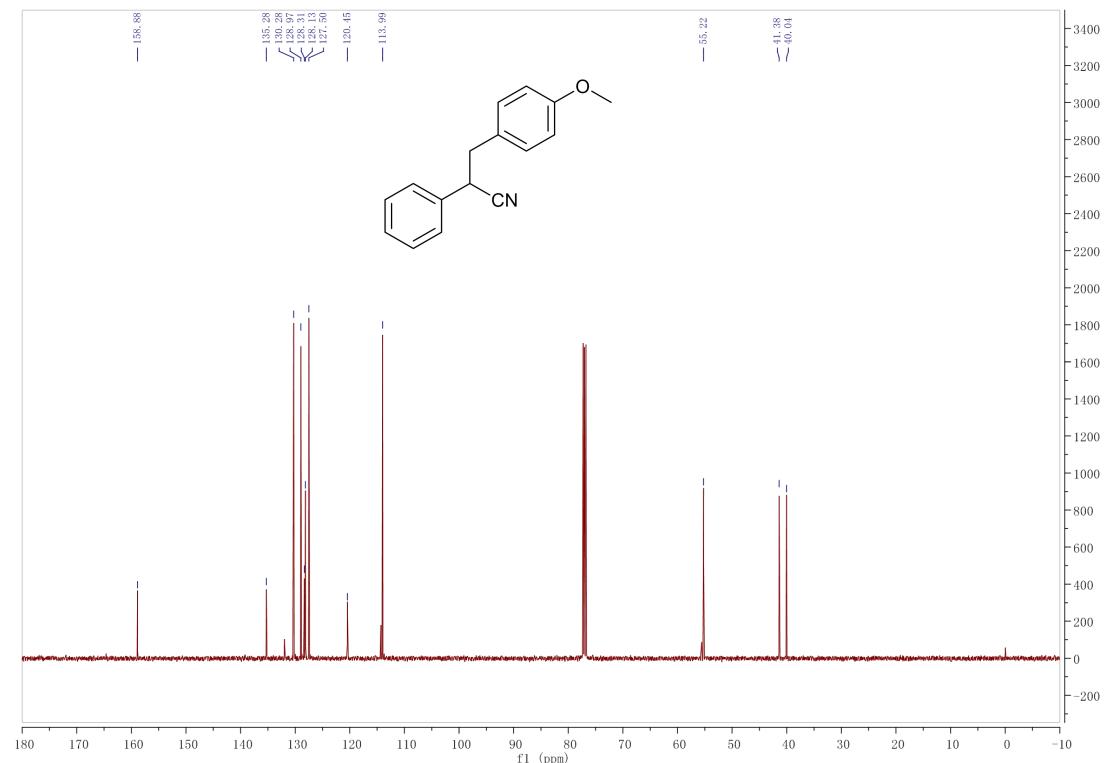
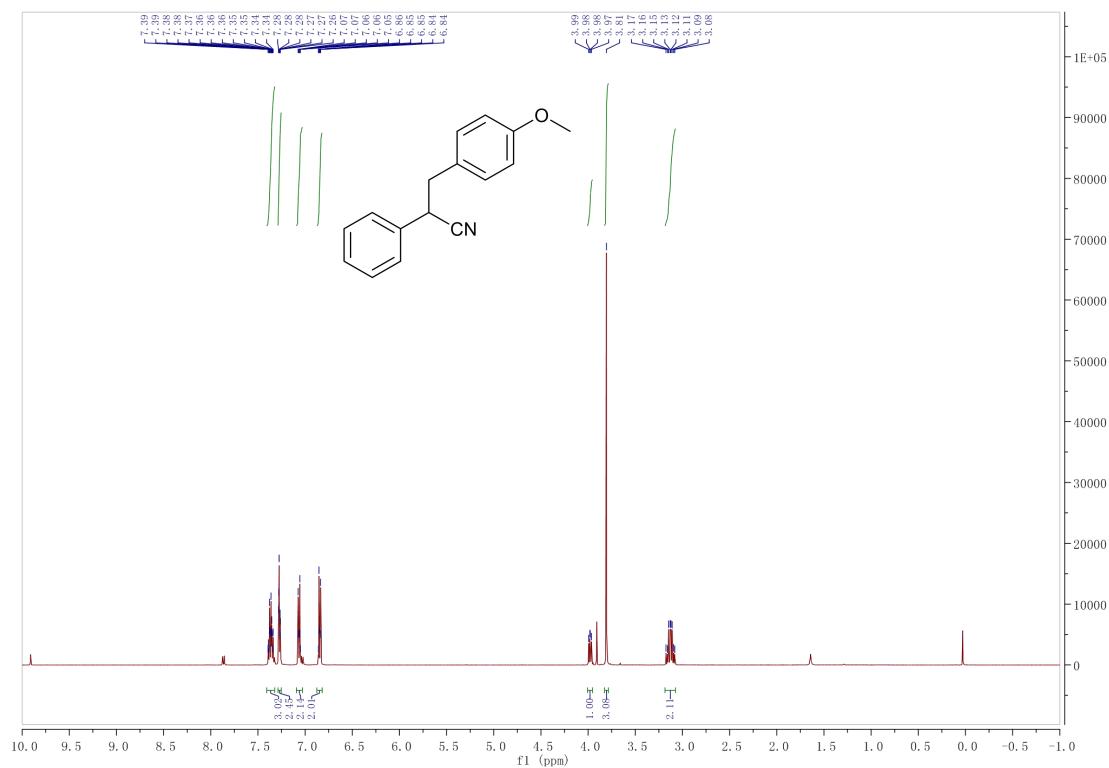
2-phenyl-3-(m-tolyl)propanenitrile (3ac**)**



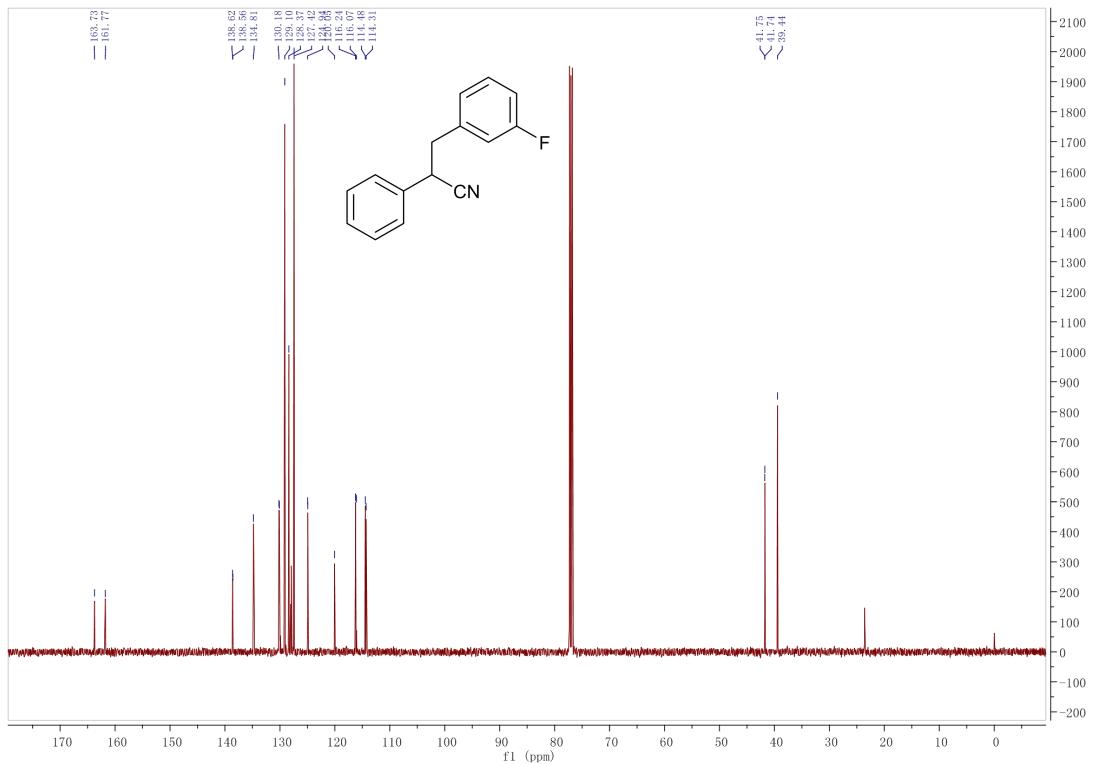
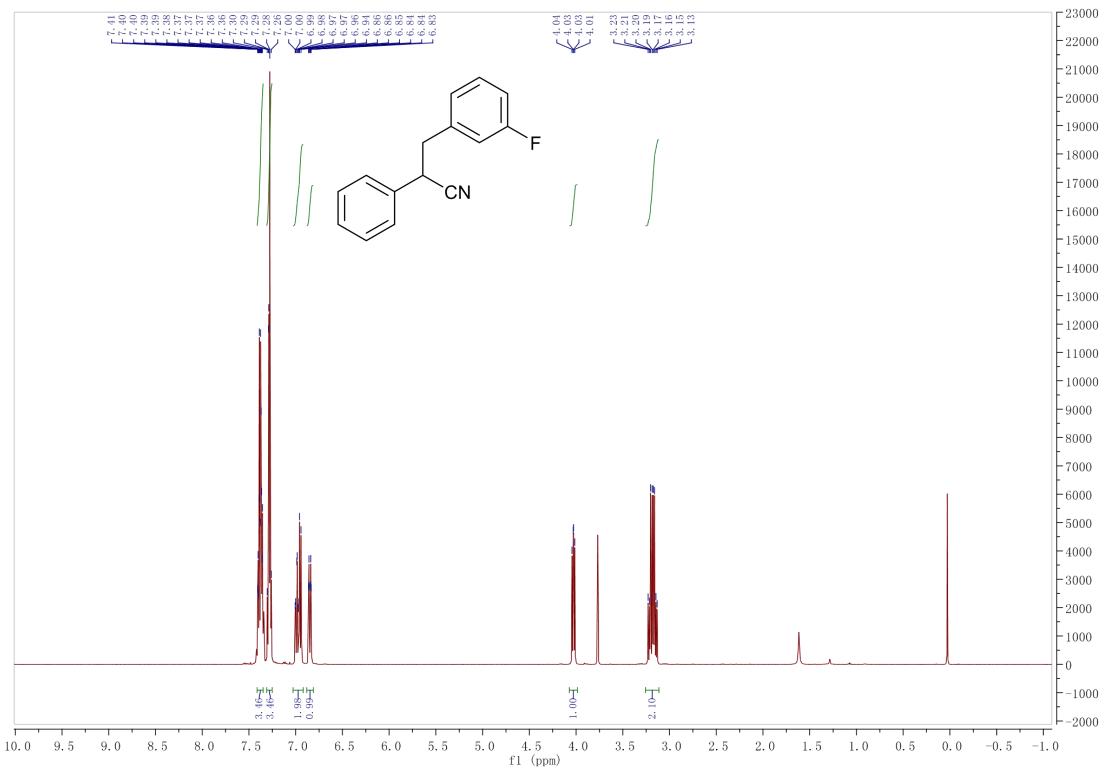
3-(3-methoxyphenyl)-2-phenylpropanenitrile (3ad**)**



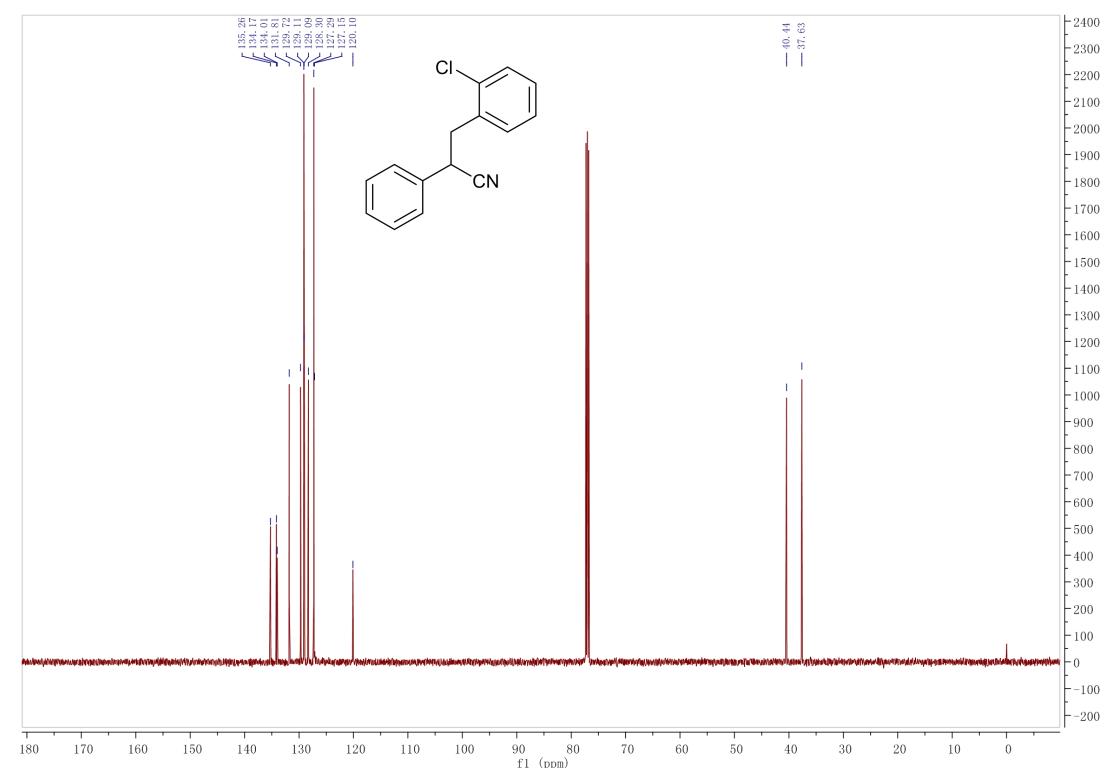
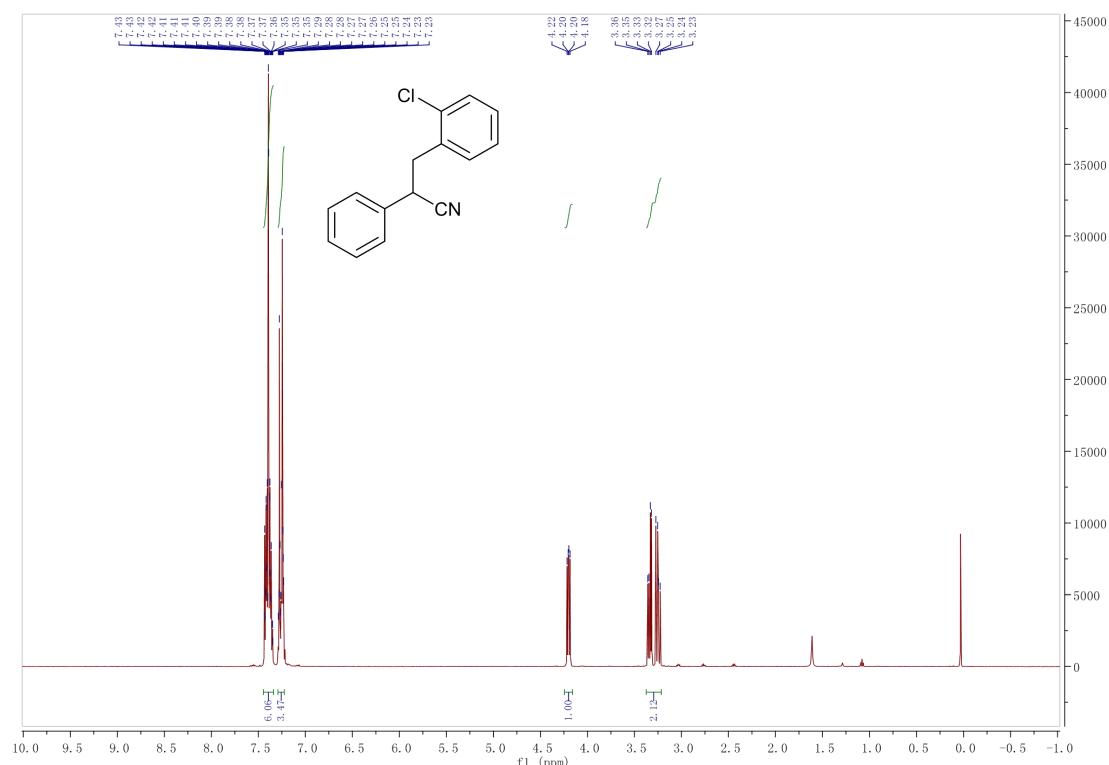
3-(4-methoxyphenyl)-2-phenylpropanenitrile (3ae**)**



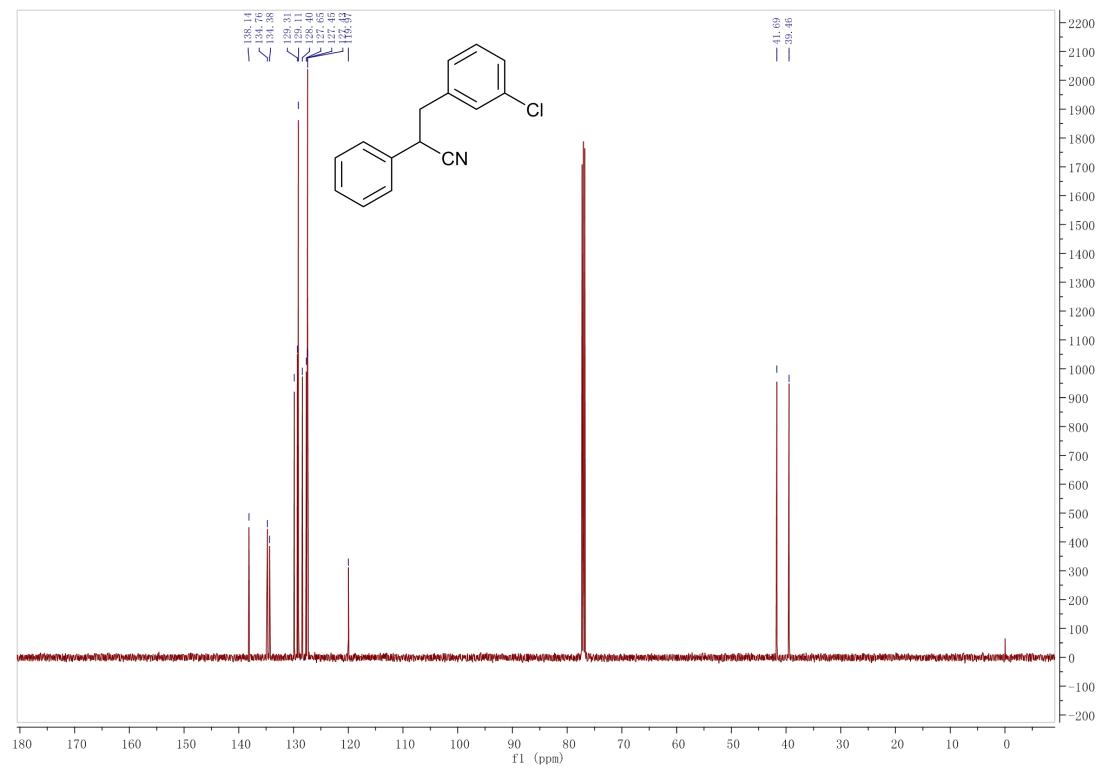
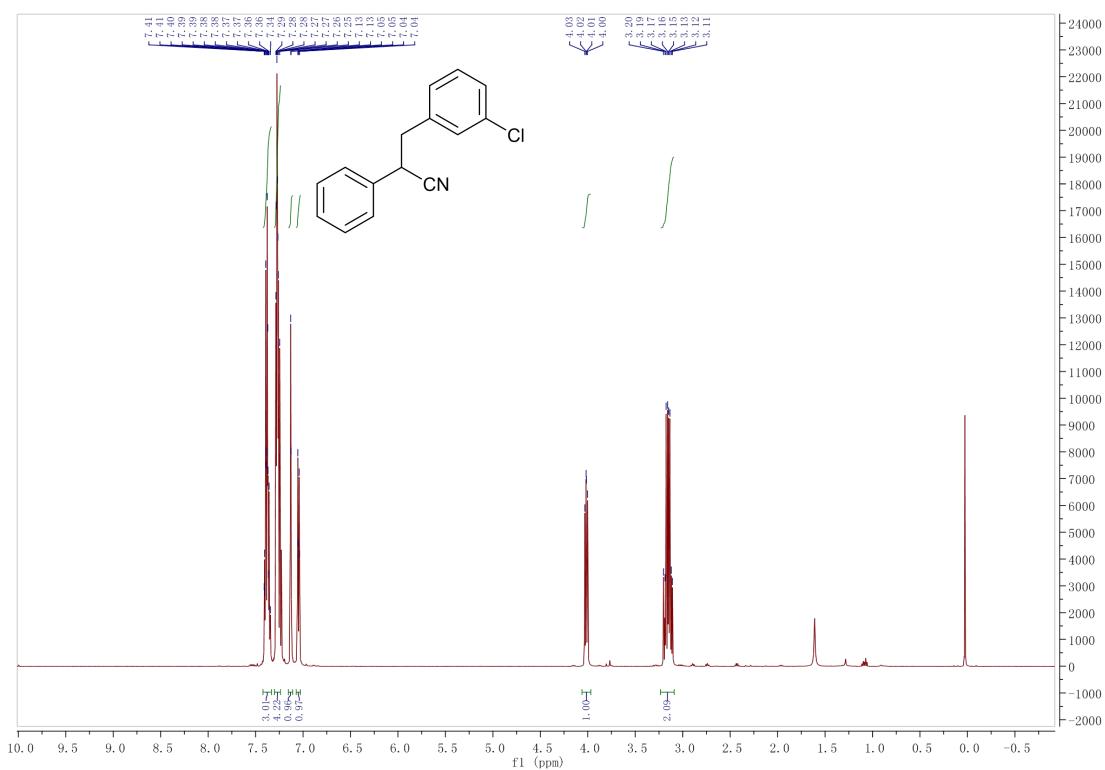
3-(3-fluorophenyl)-2-phenylpropanenitrile (3af**)**



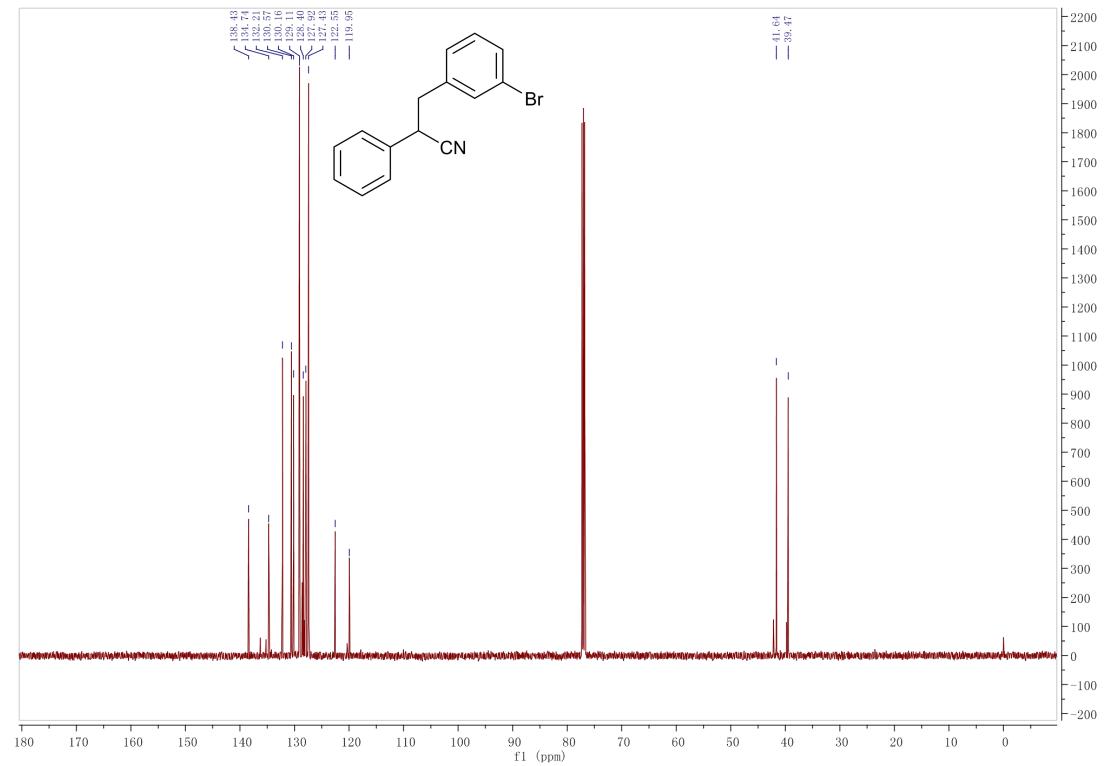
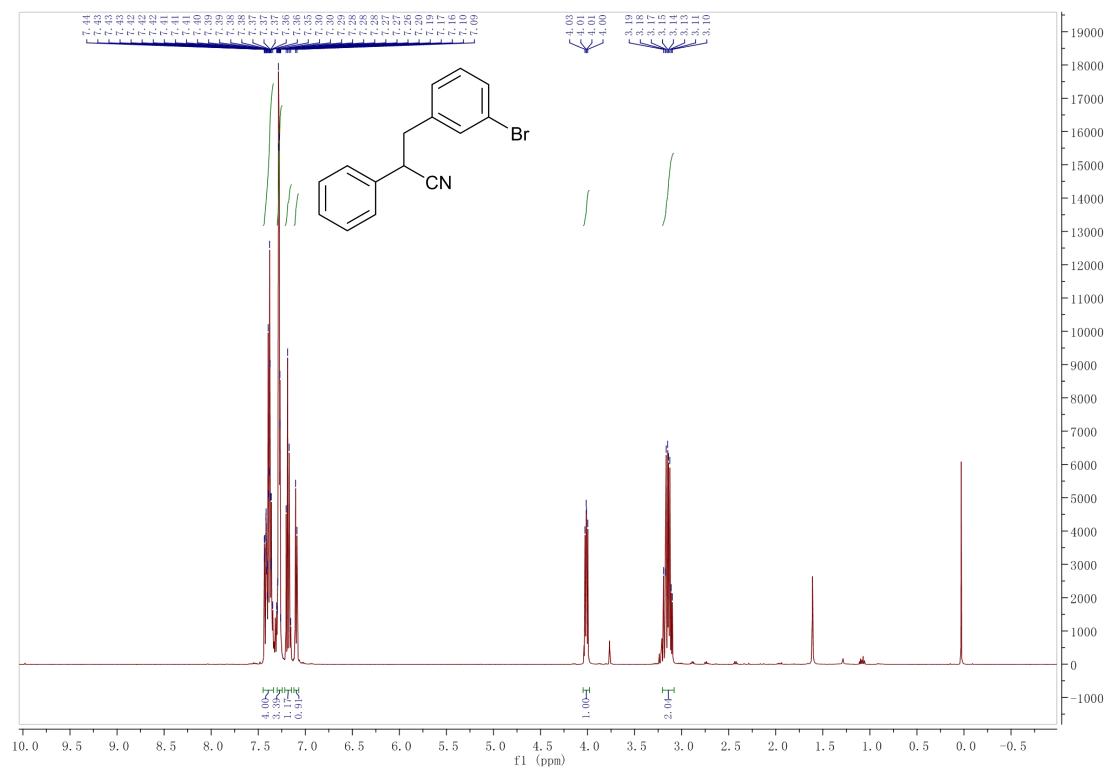
3-(2-chlorophenyl)-2-phenylpropanenitrile (3ag**)**



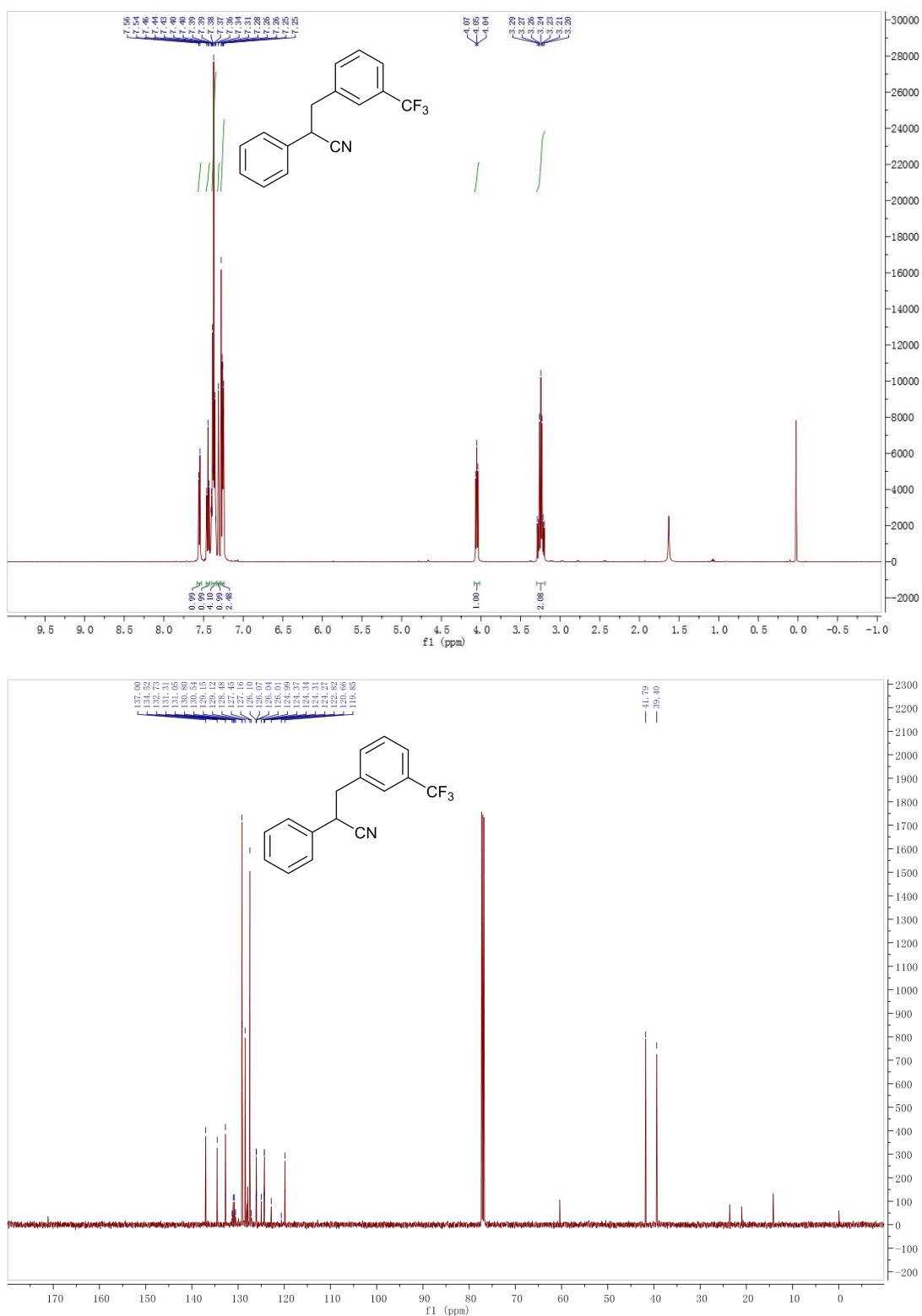
3-(3-chlorophenyl)-2-phenylpropanenitrile (3ah**)**



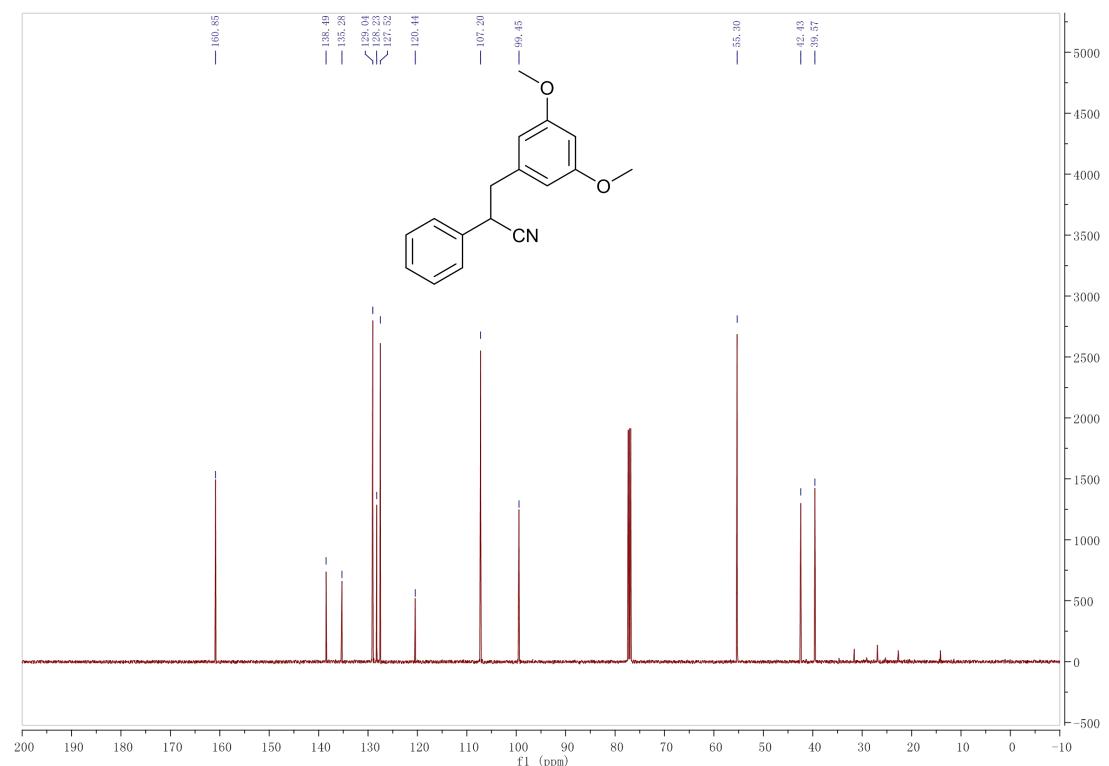
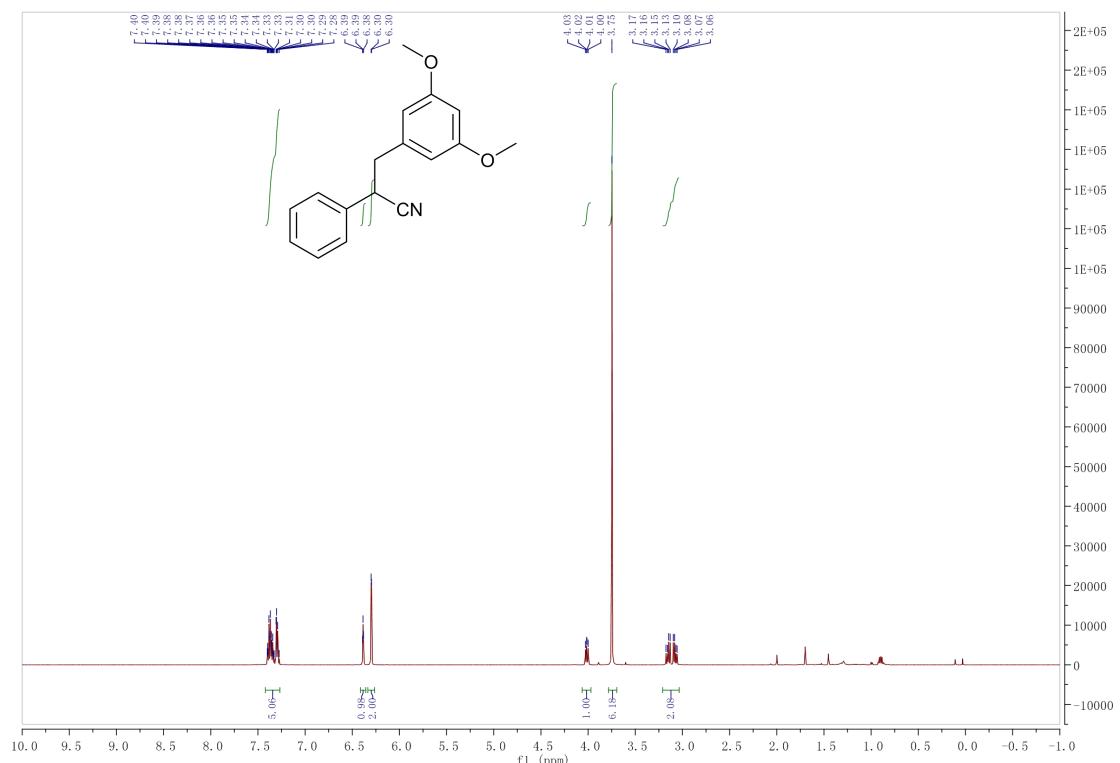
3-(3-bromophenyl)-2-phenylpropanenitrile (3ai**)**



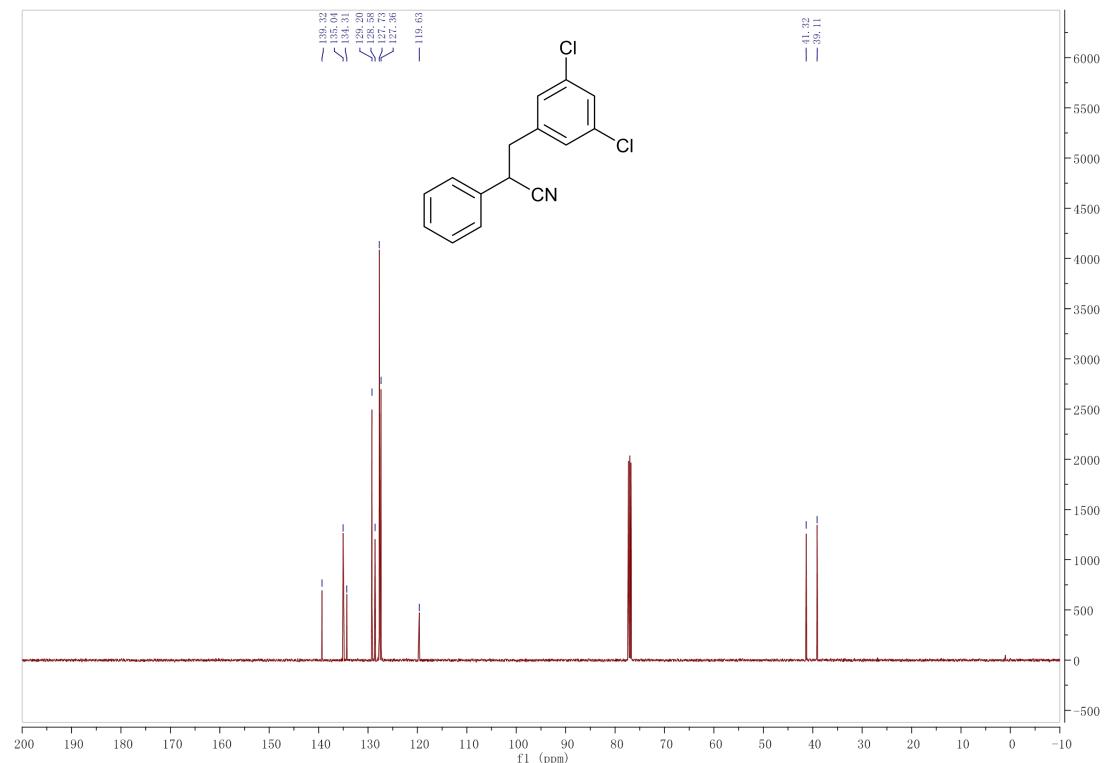
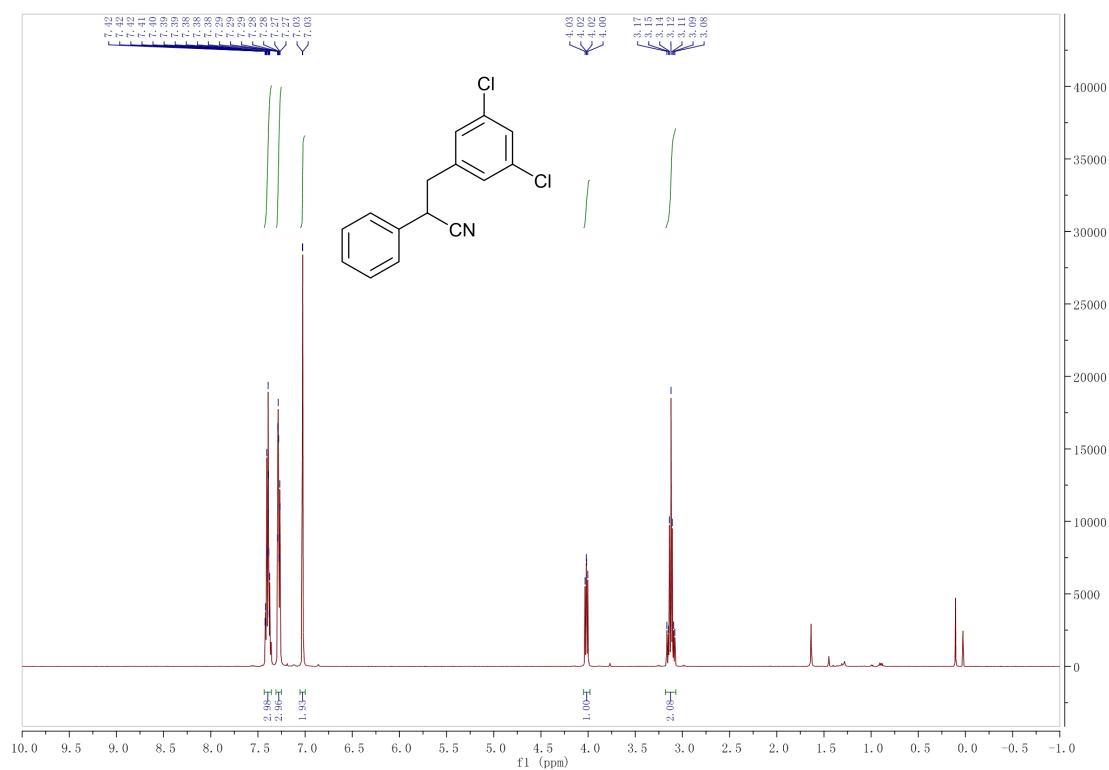
2-phenyl-3-(3-(trifluoromethyl)phenyl)propanenitrile (3aj**)**



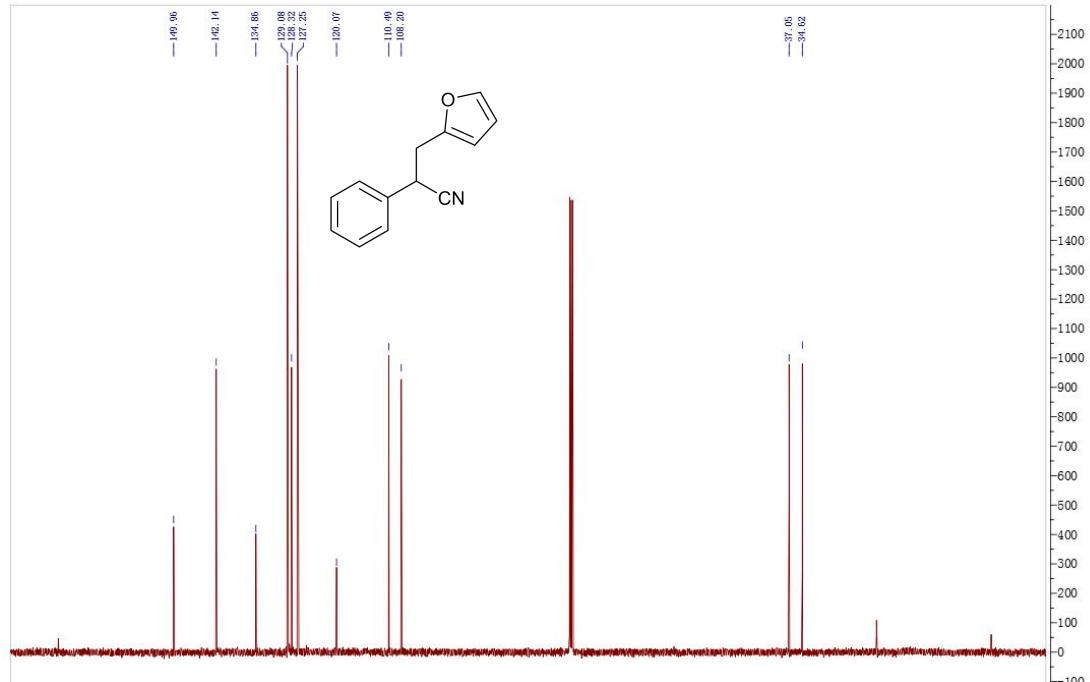
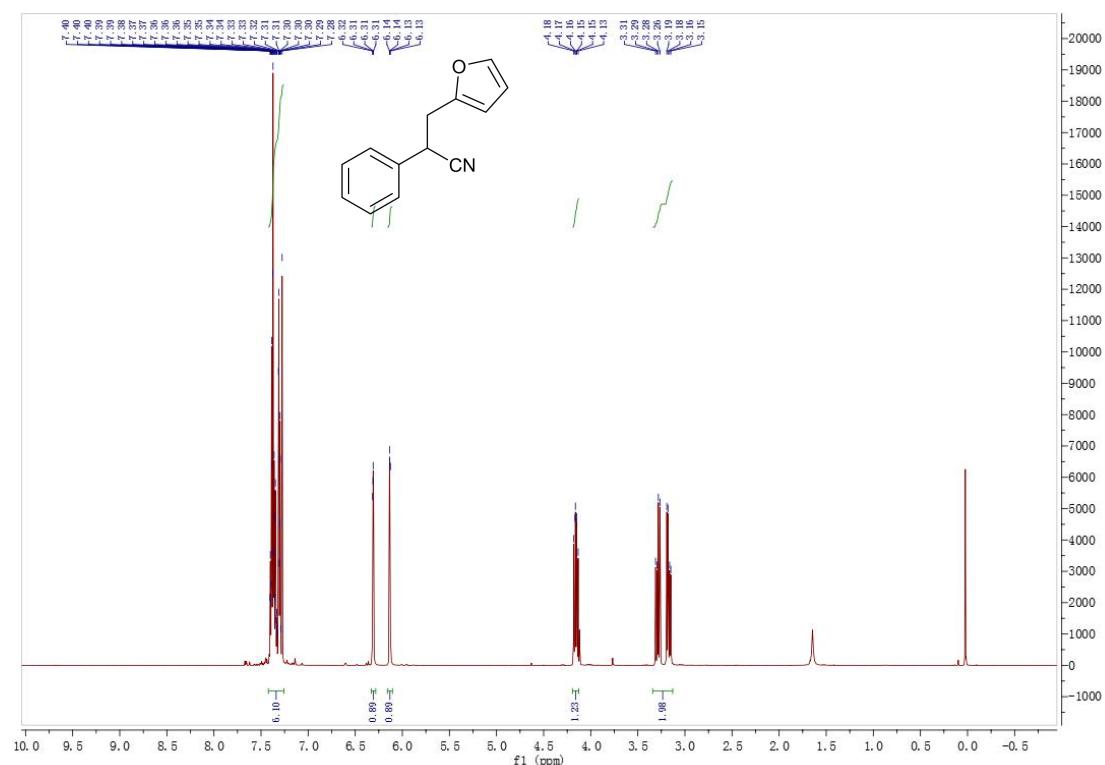
3-(3,5-dimethoxyphenyl)-2-phenylpropanenitrile (3ak**)**



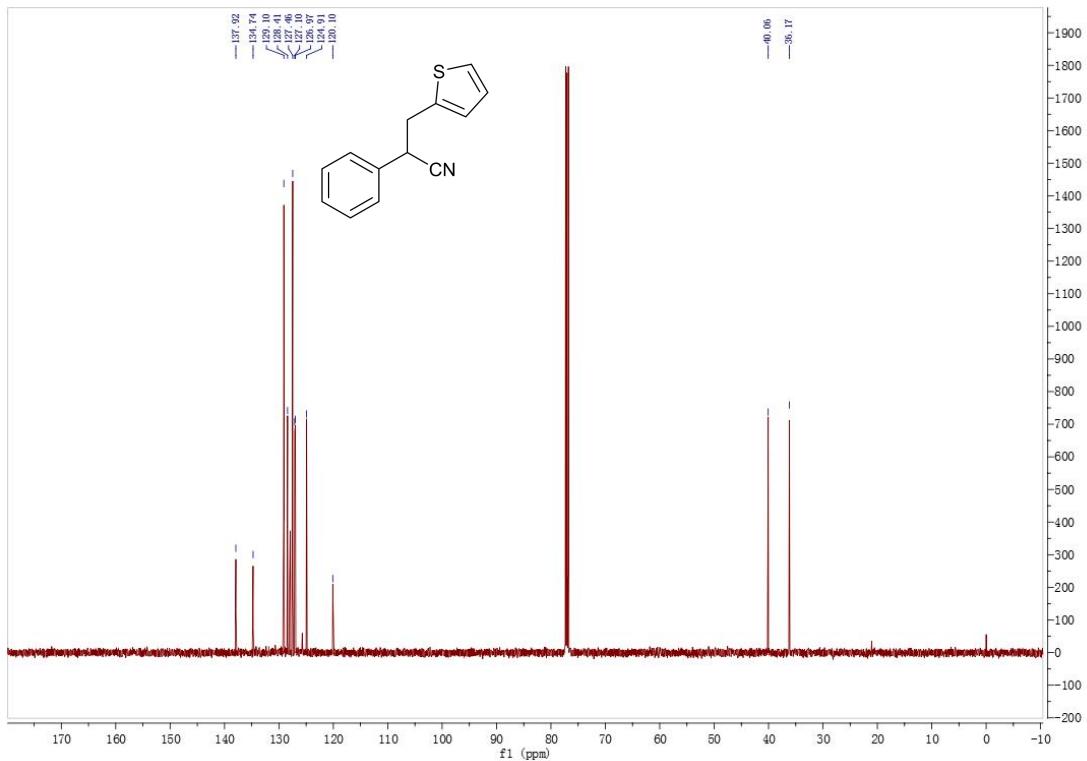
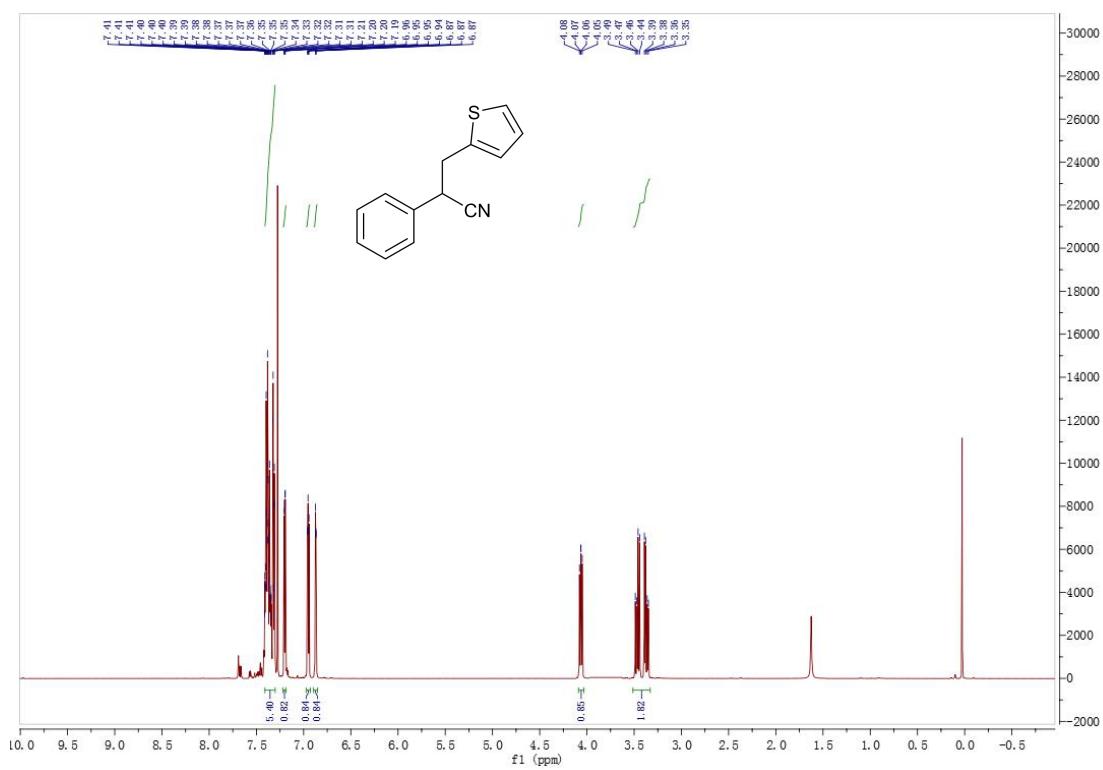
3-(3,5-dichlorophenyl)-2-phenylpropanenitrile (3al**)**



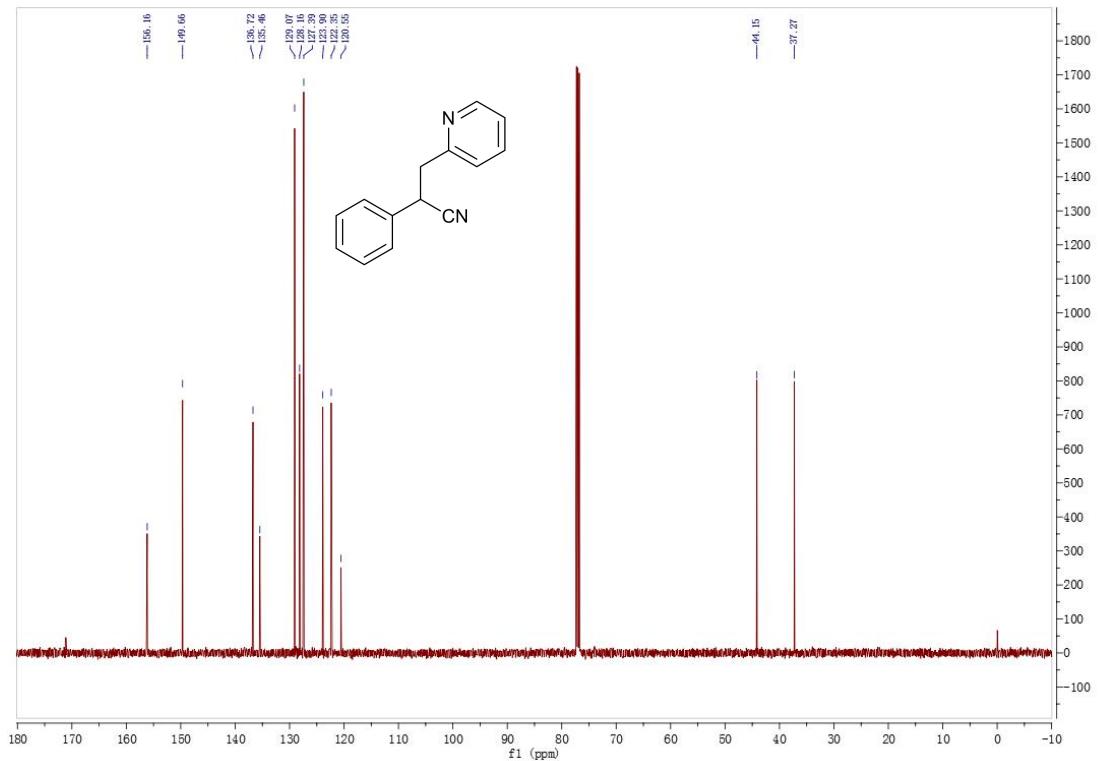
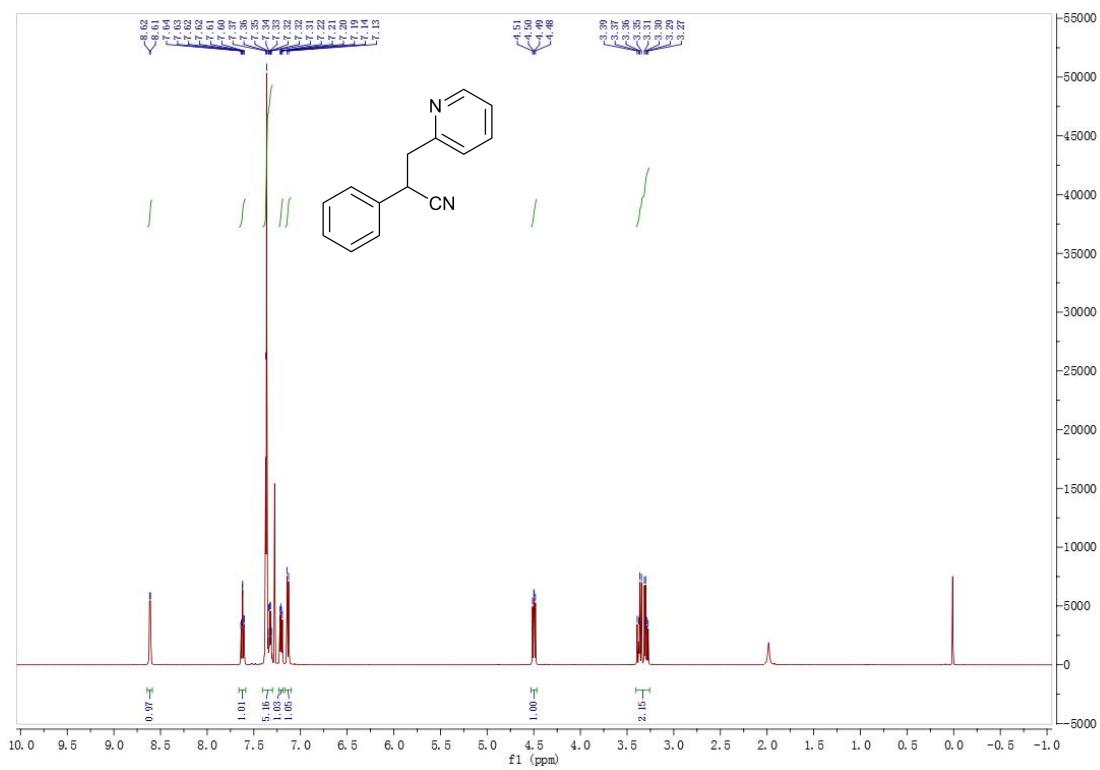
3-(furan-2-yl)-2-phenylpropanenitrile (3am**)**



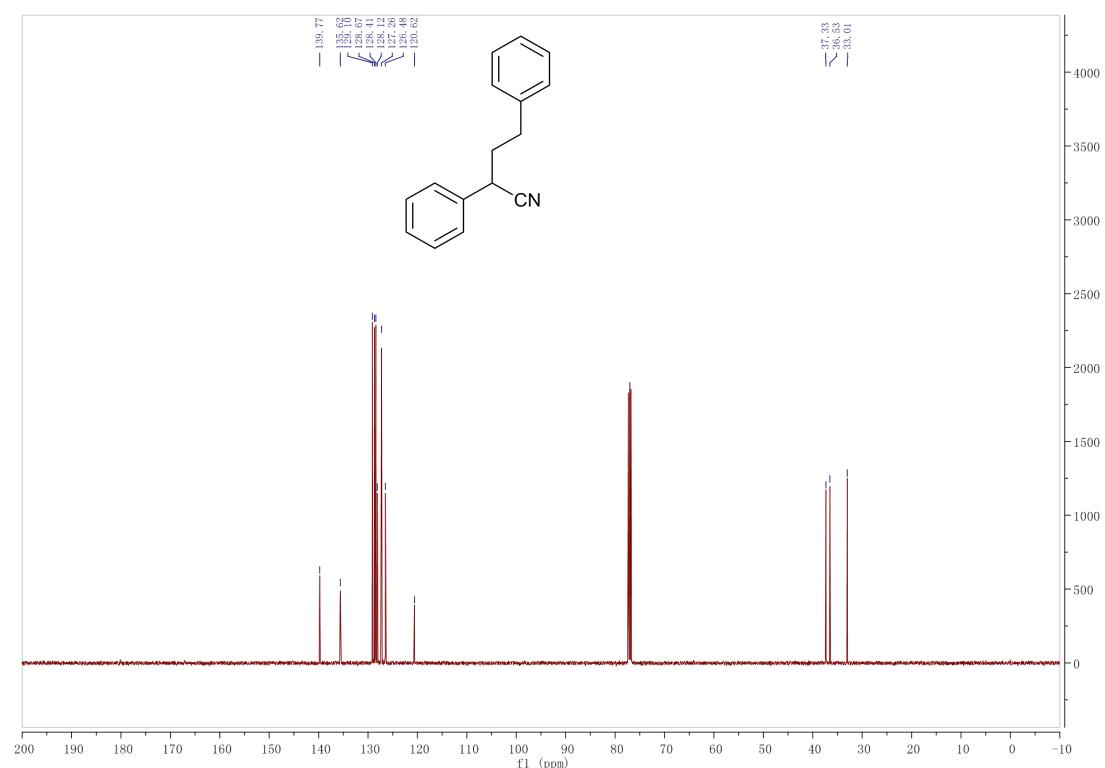
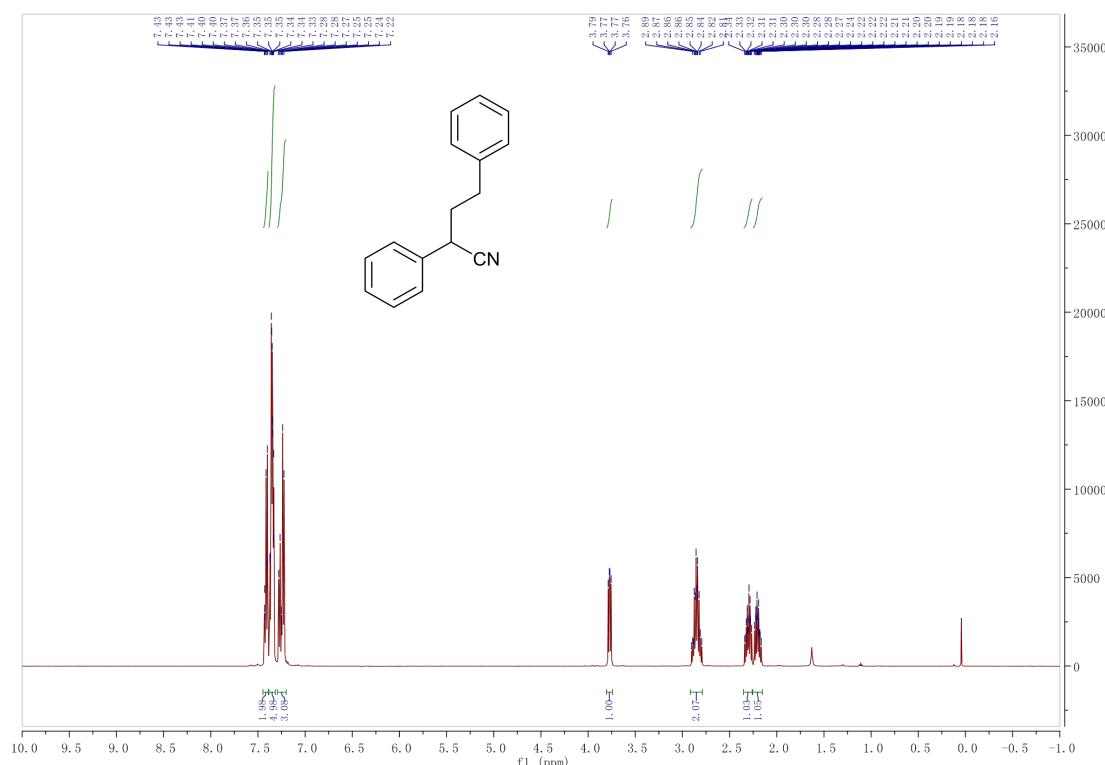
2-phenyl-3-(thiophen-2-yl)propanenitrile (**3an**)



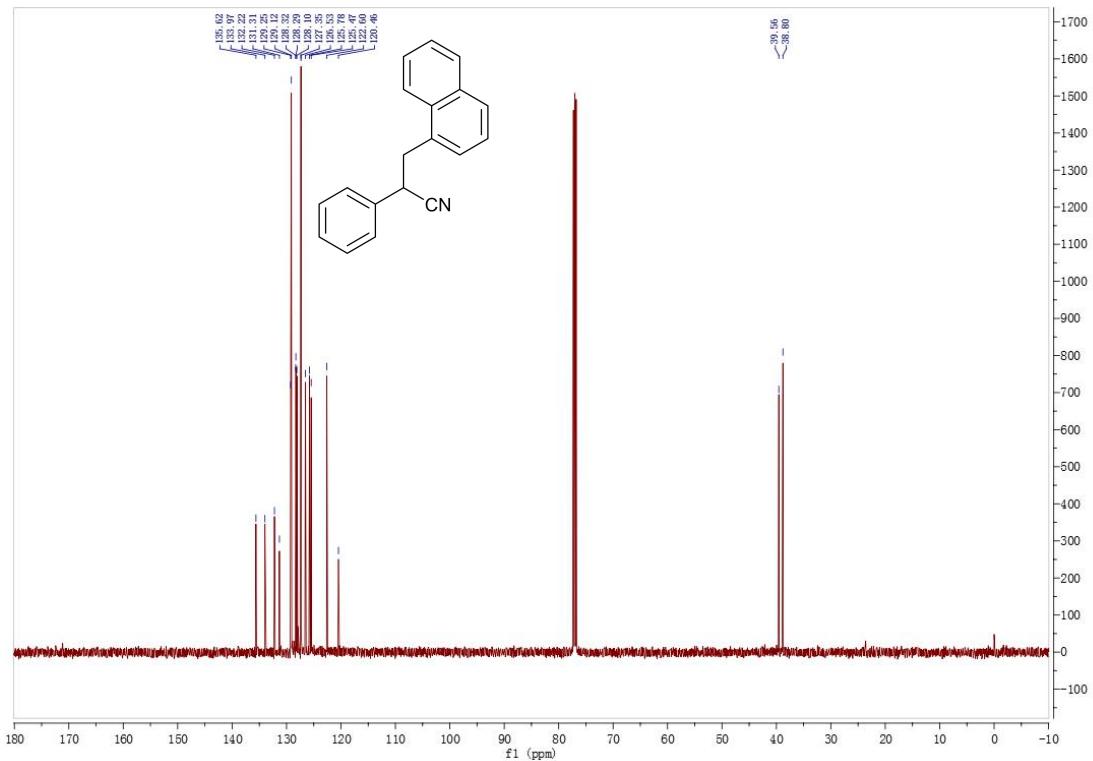
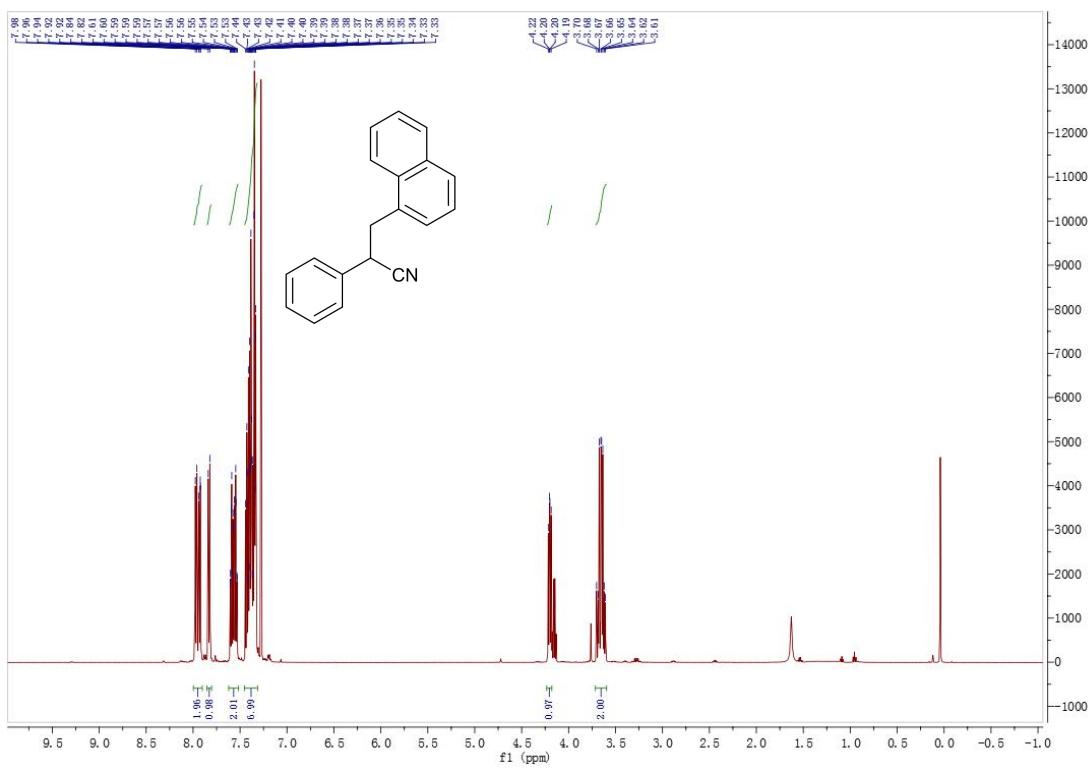
2-phenyl-3-(pyridin-2-yl)propanenitrile (3ao**)**



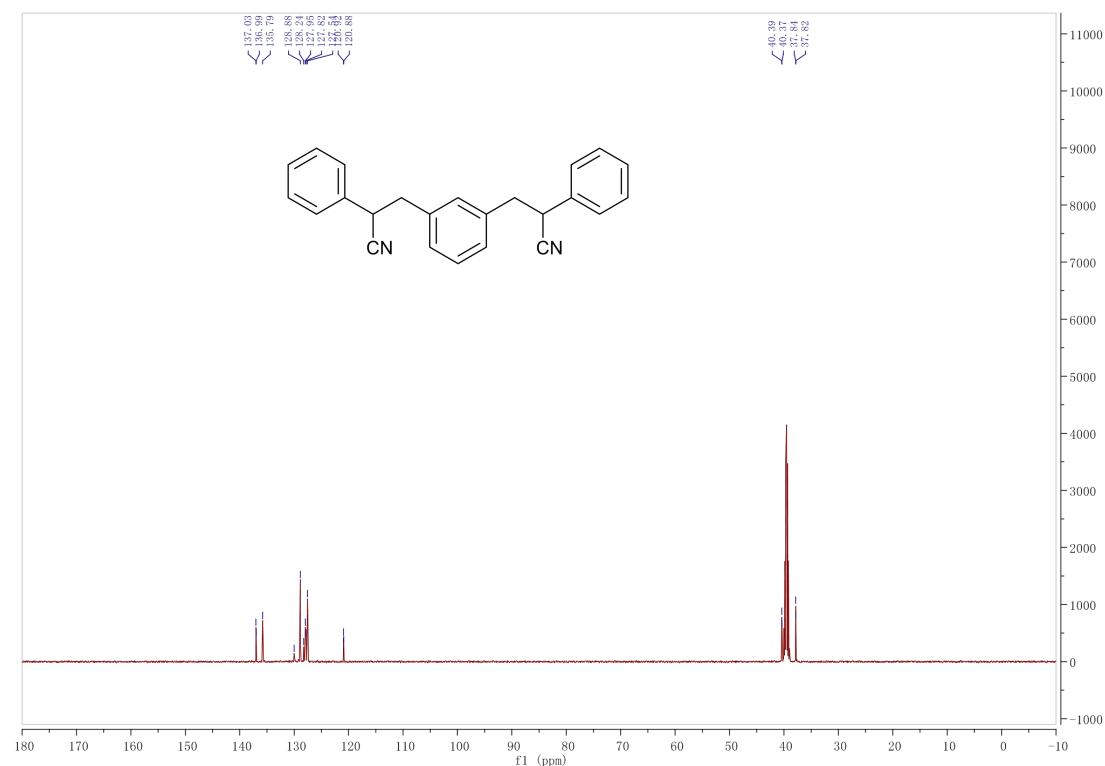
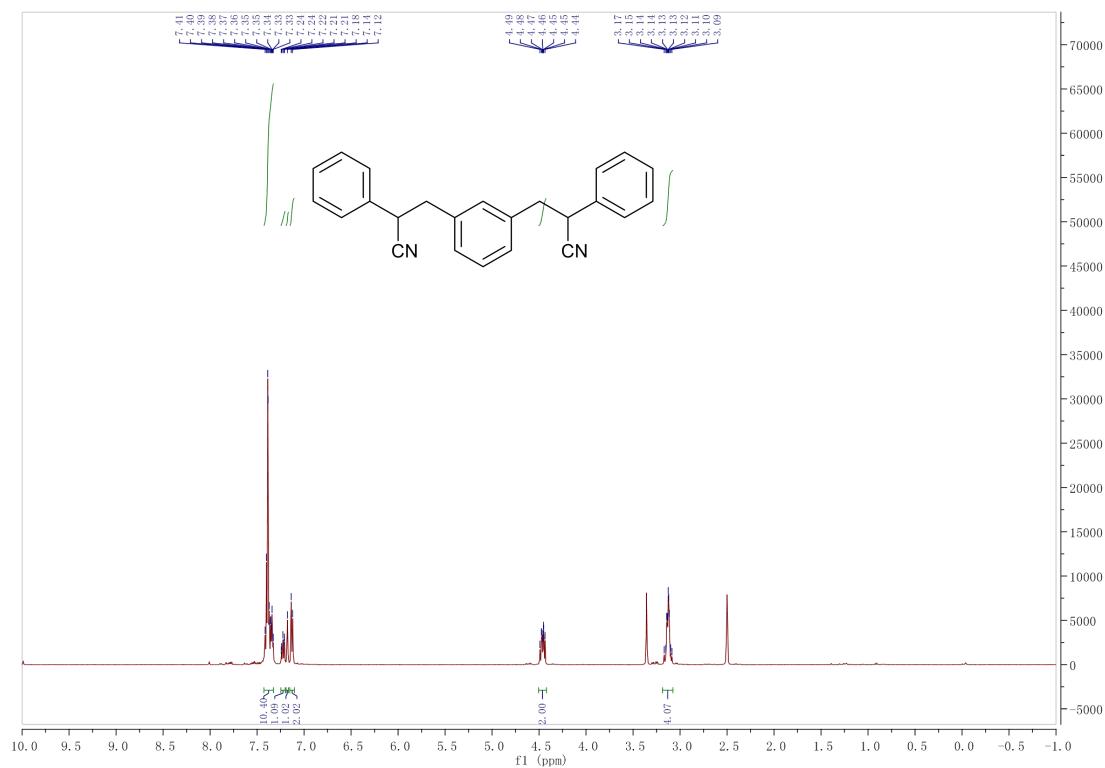
2,4-diphenylbutanenitrile (3ap**)**



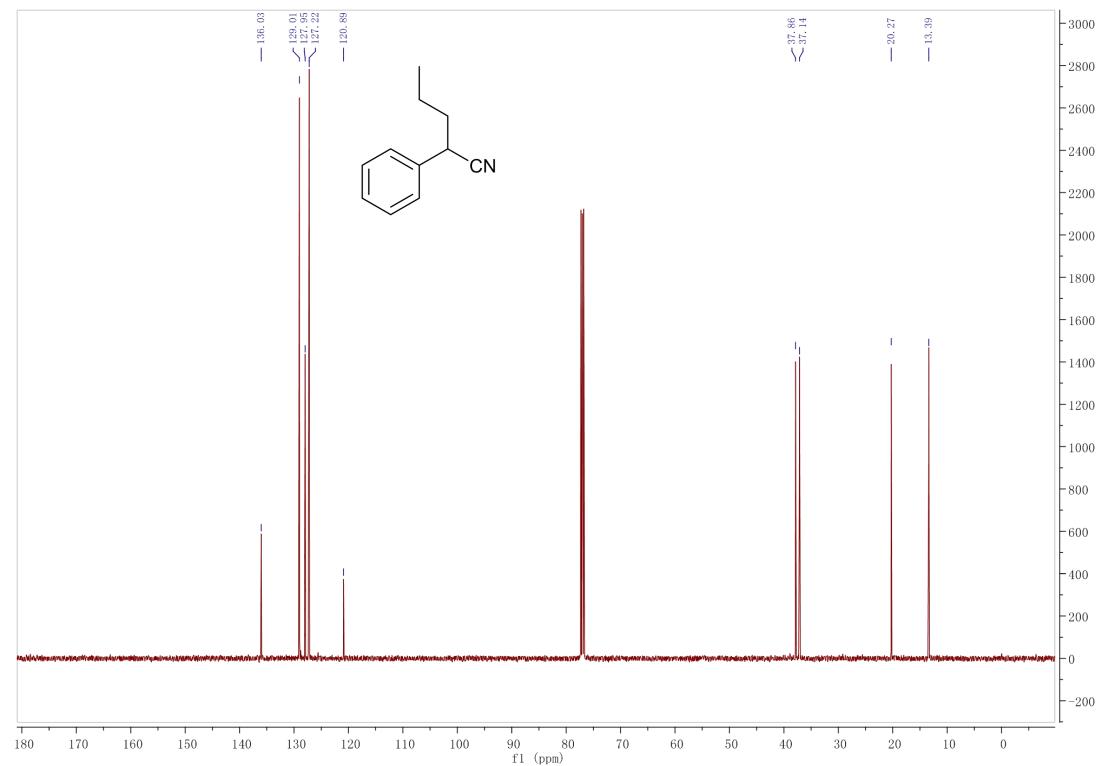
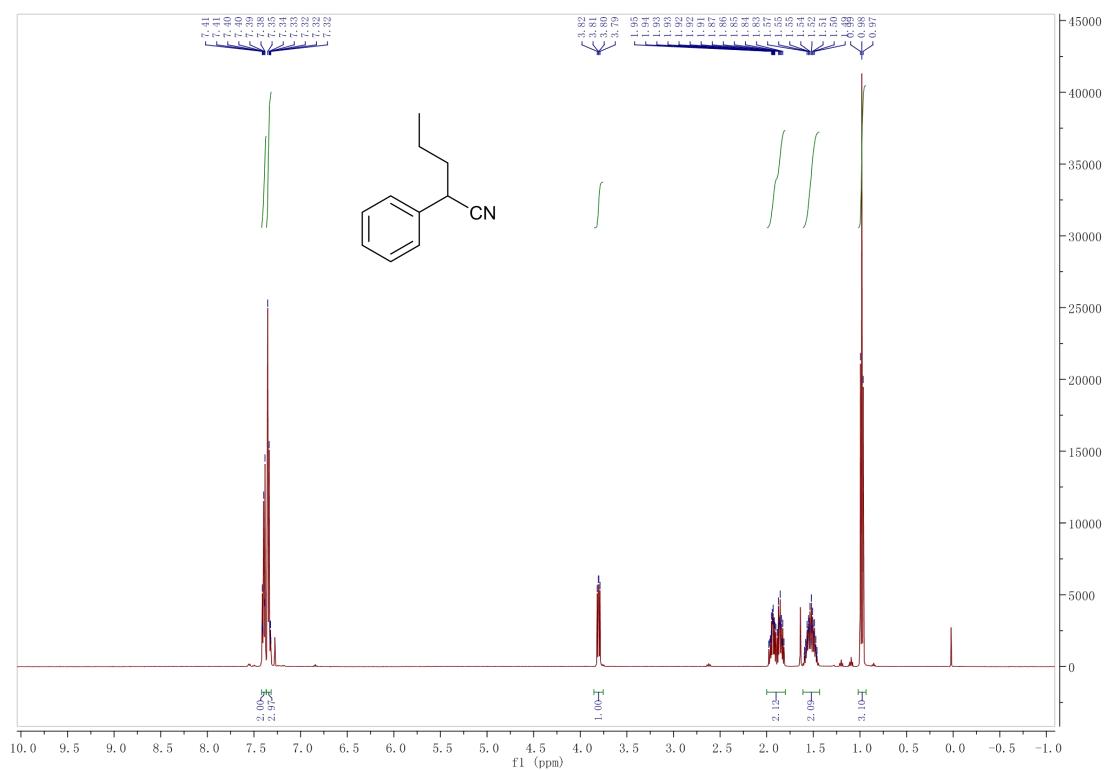
3-(naphthalen-1-yl)-2-phenylpropanenitrile (3aq**)**



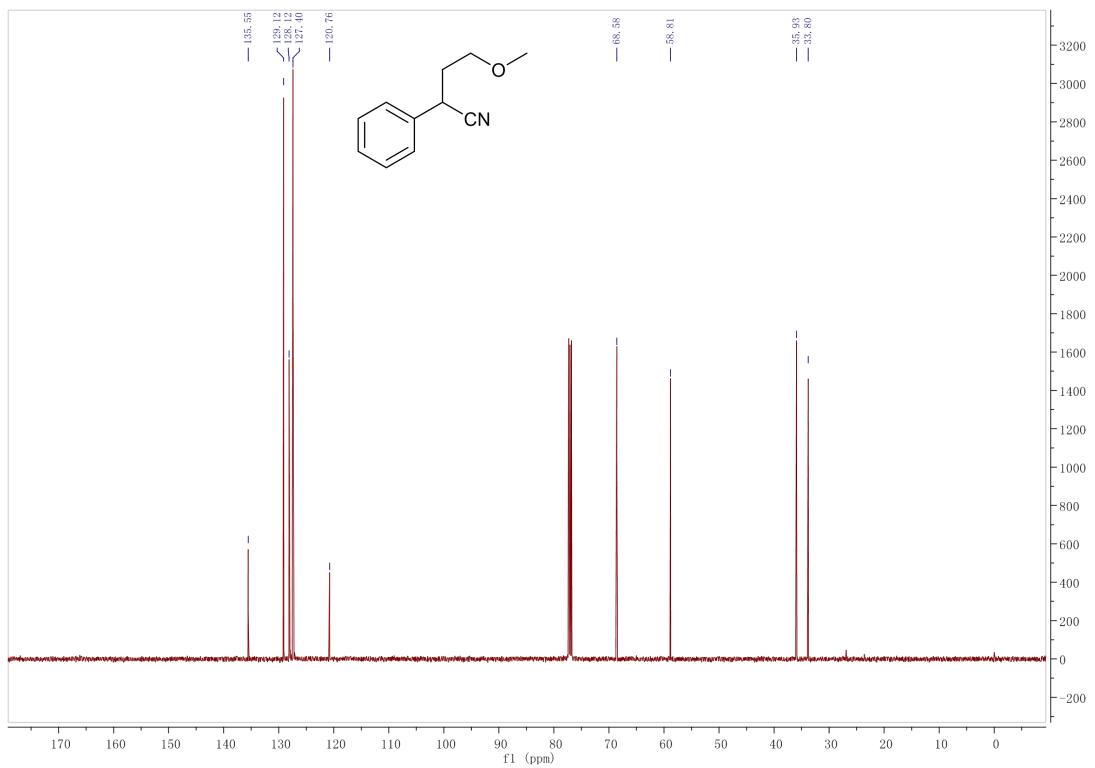
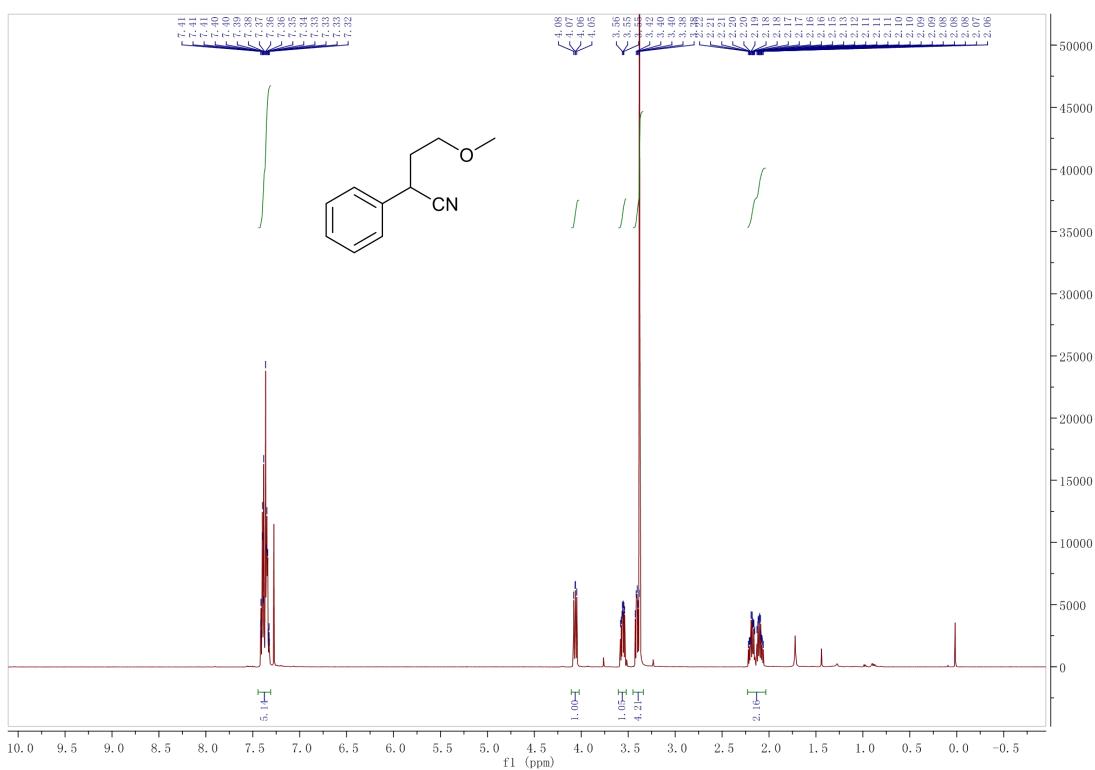
3,3'-(1,3-phenylene)bis(2-phenylpropanenitrile) (3ar**)**



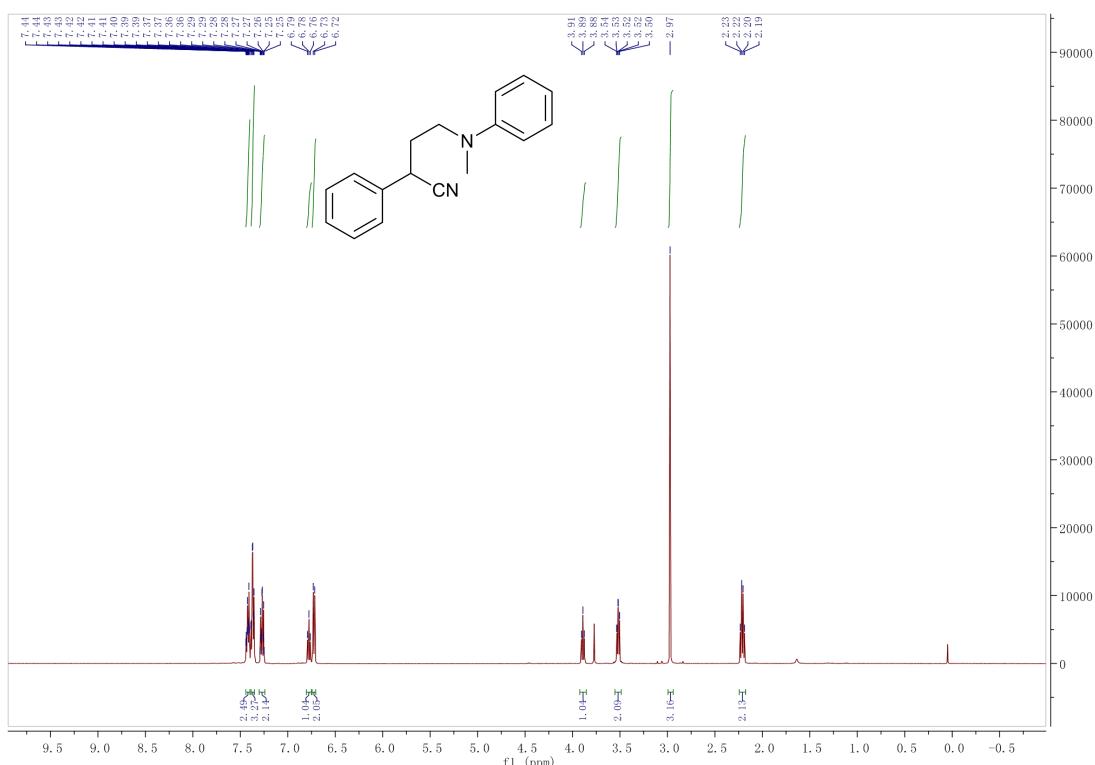
2-phenylpentanenitrile (3as**)**



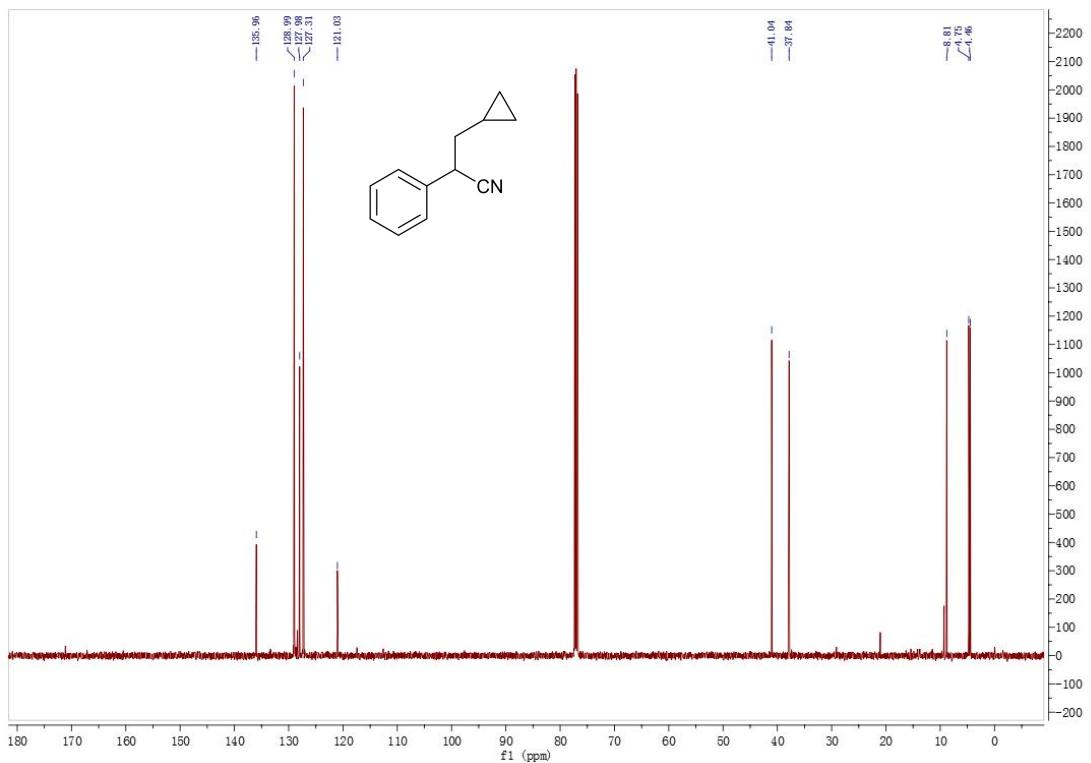
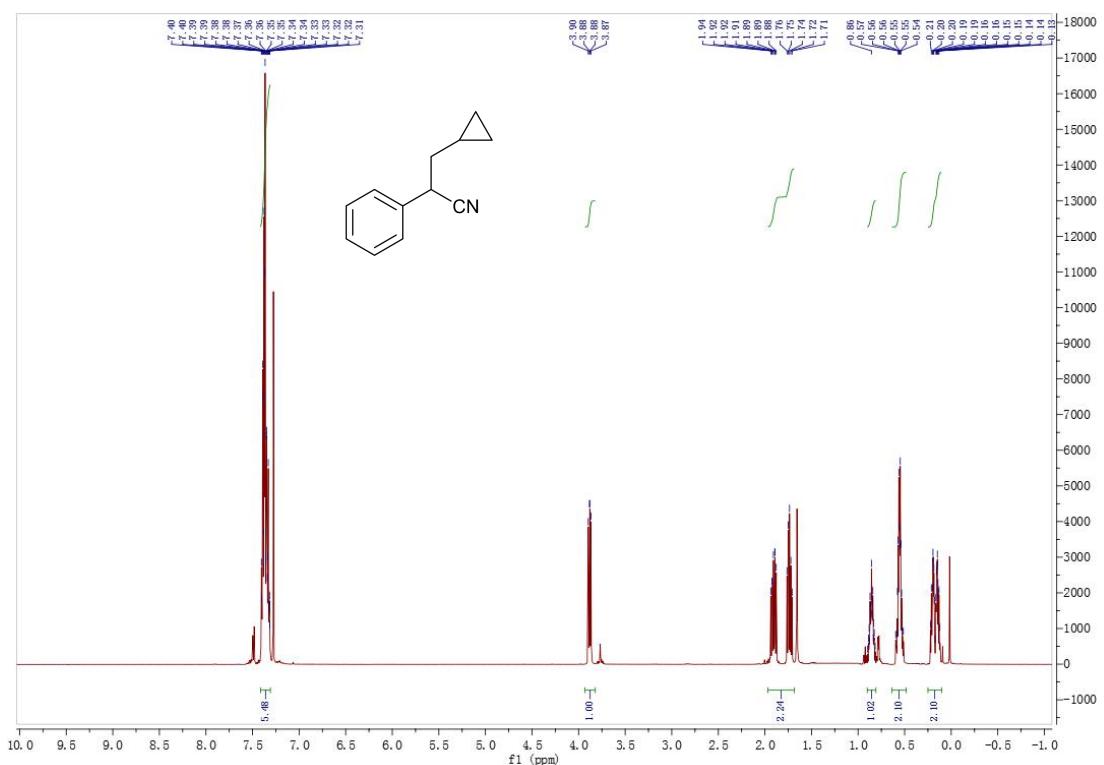
4-methoxy-2-phenylbutanenitrile (3at**)**



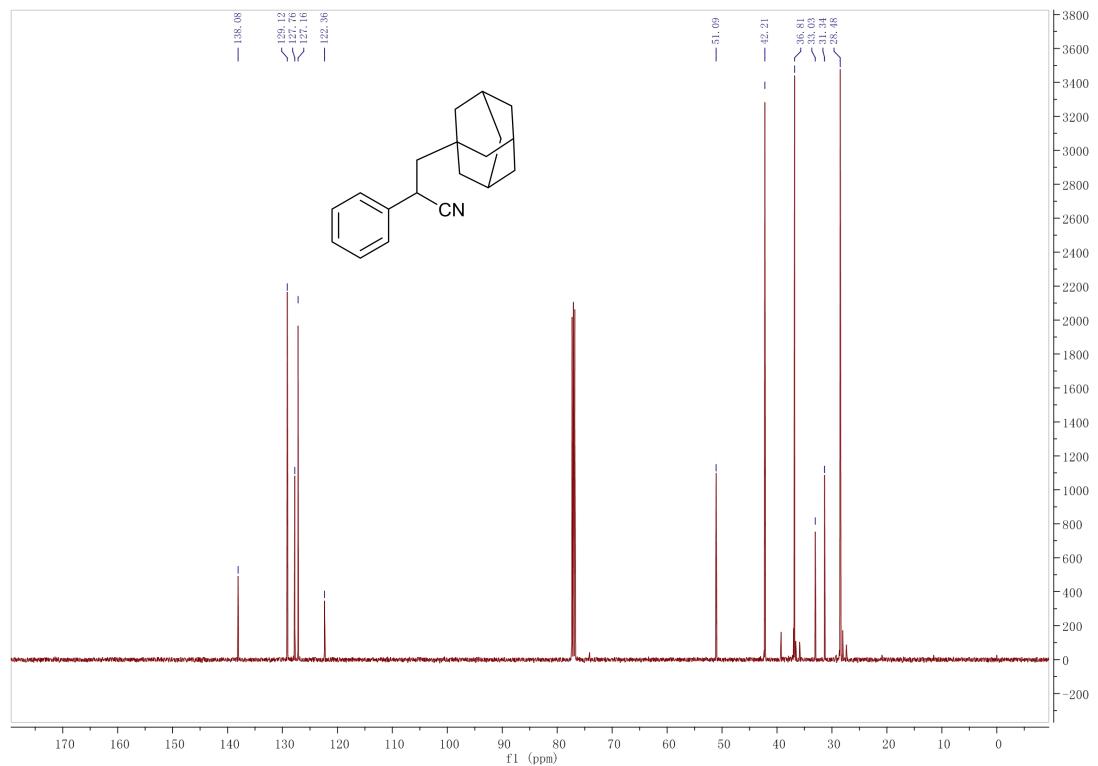
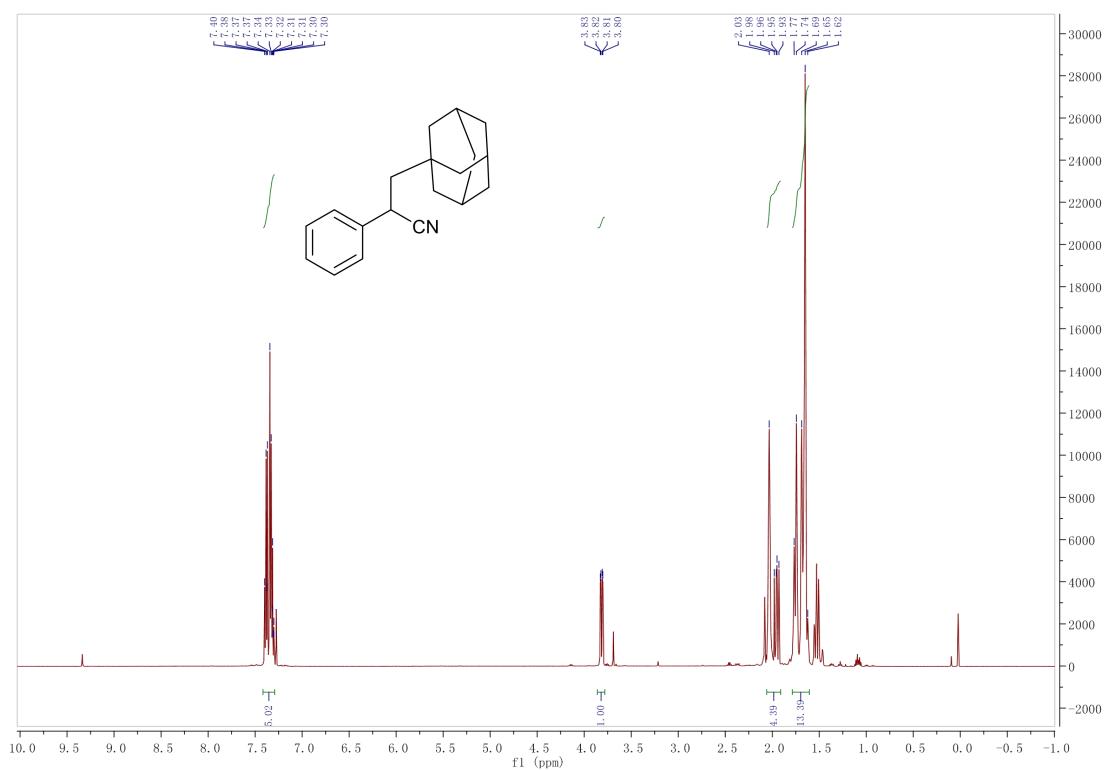
4-(methyl(phenyl)amino)-2-phenylbutanenitrile (3au**)**



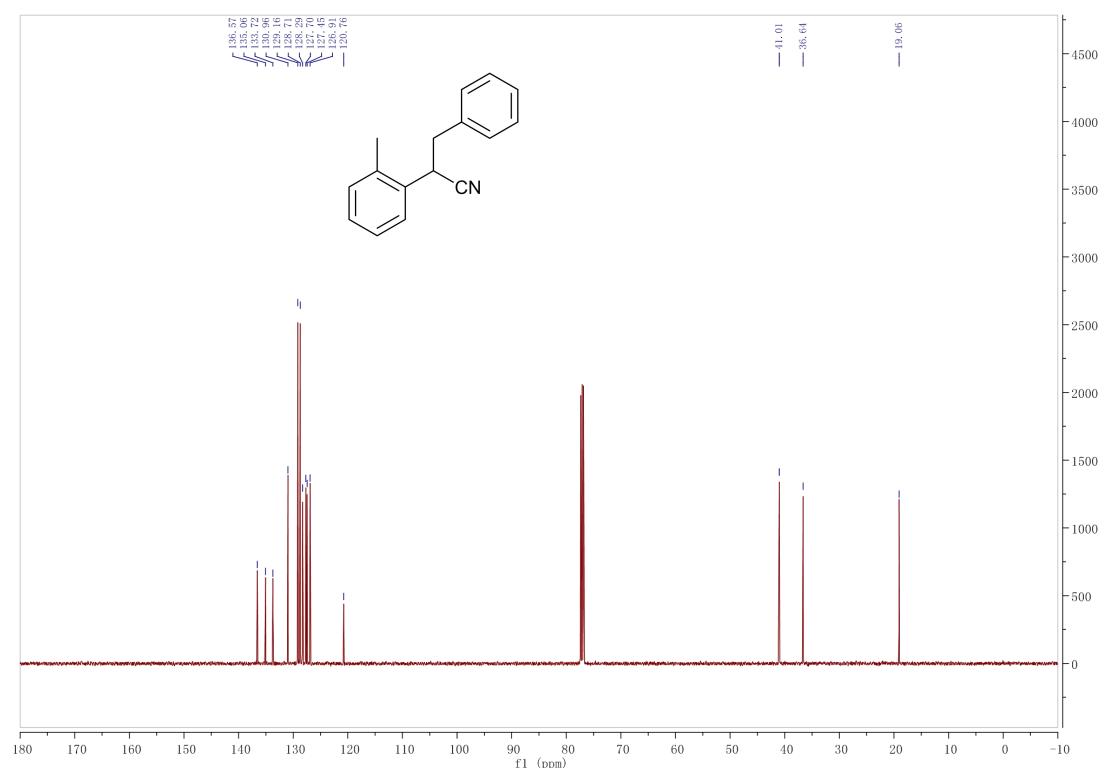
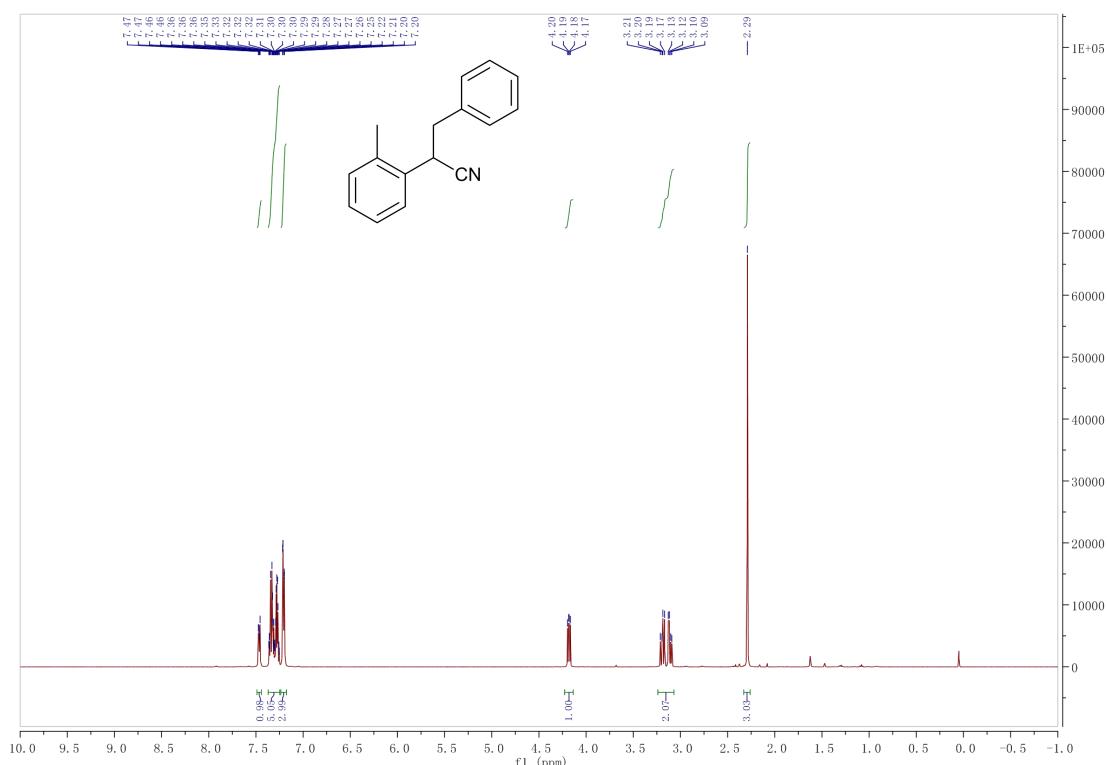
3-cyclopropyl-2-phenylpropanenitrile (**3av**)



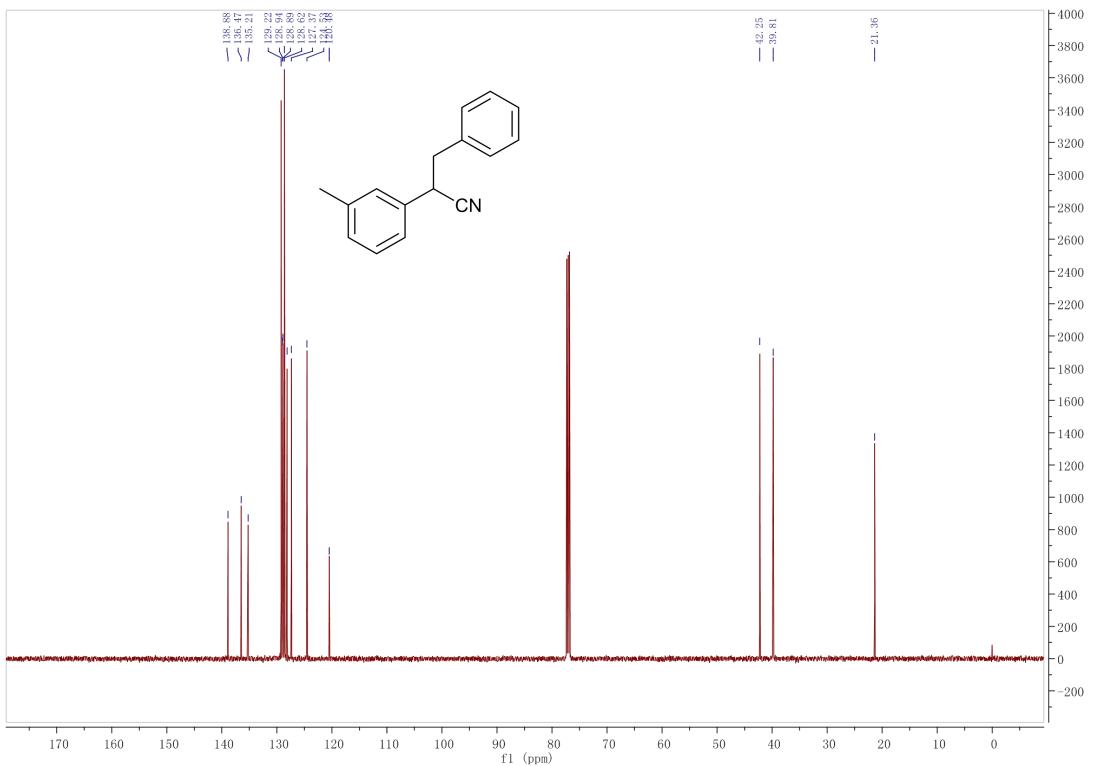
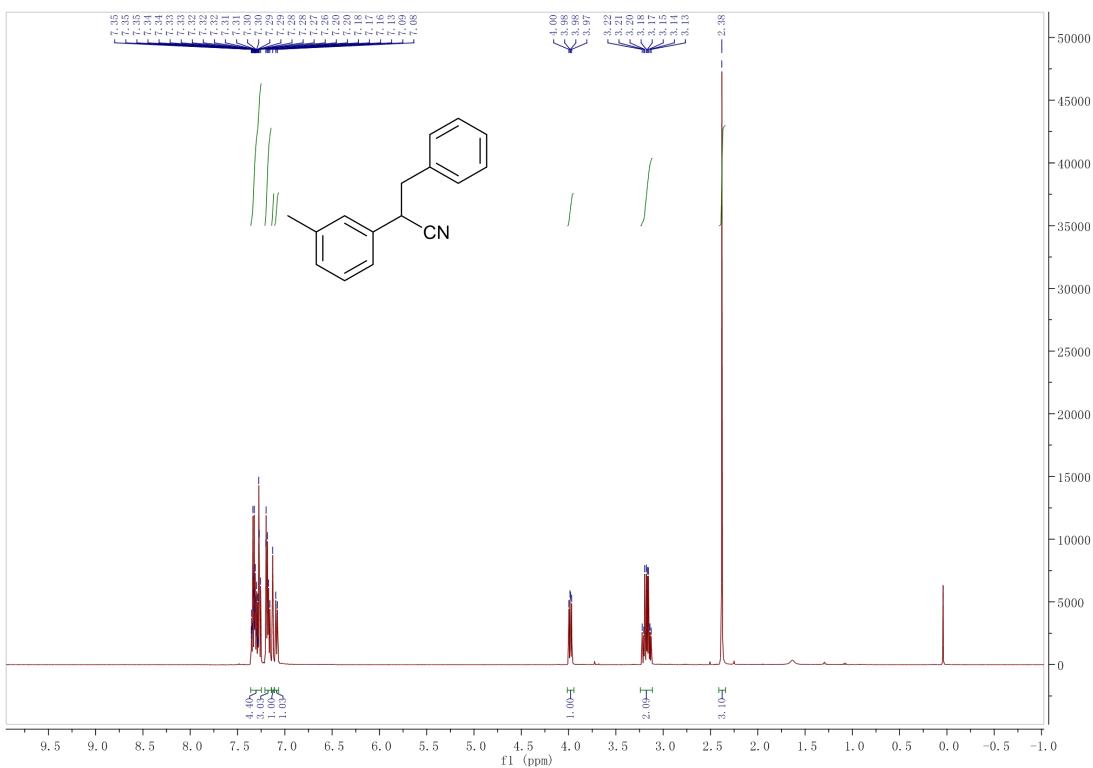
3-(adamantan-1-yl)-2-phenylpropanenitrile (3aw**)**



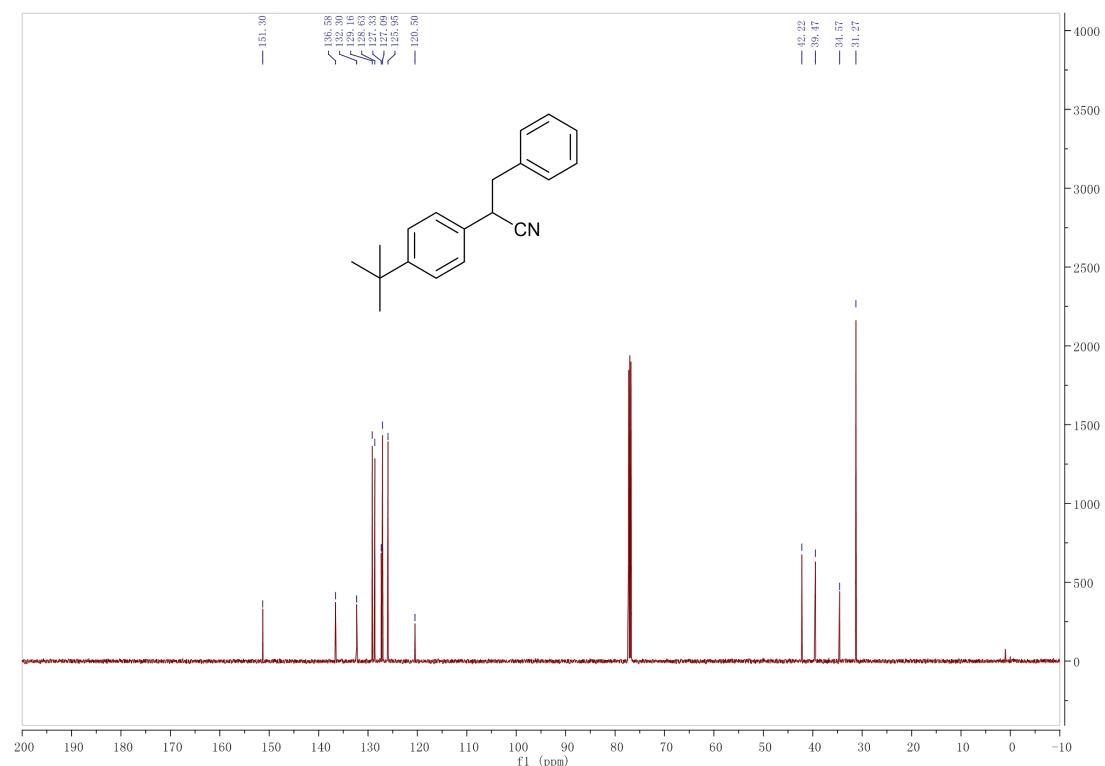
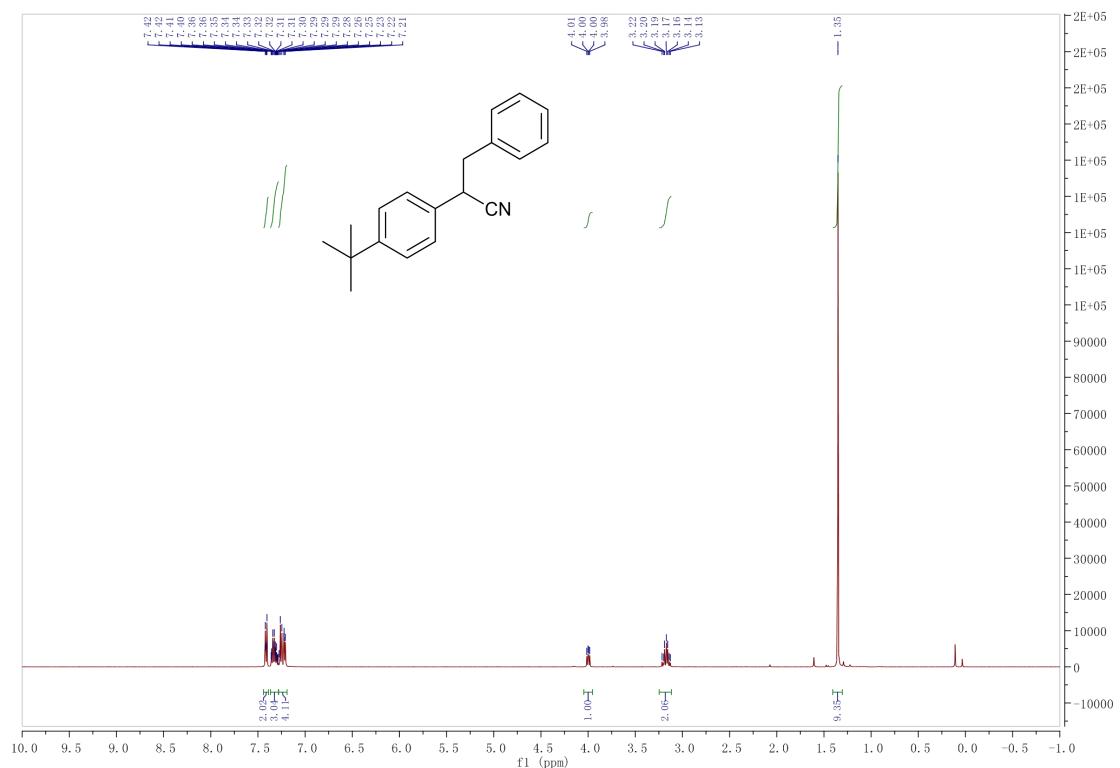
3-phenyl-2-(o-tolyl)propanenitrile (3ba**)**



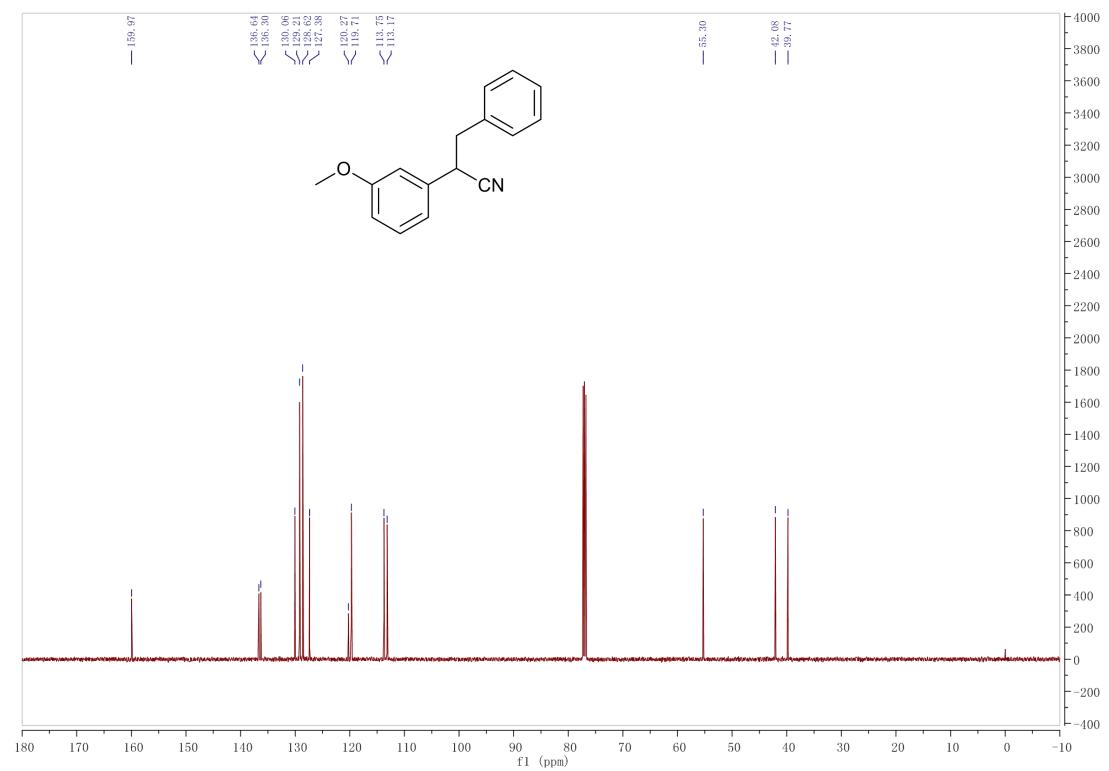
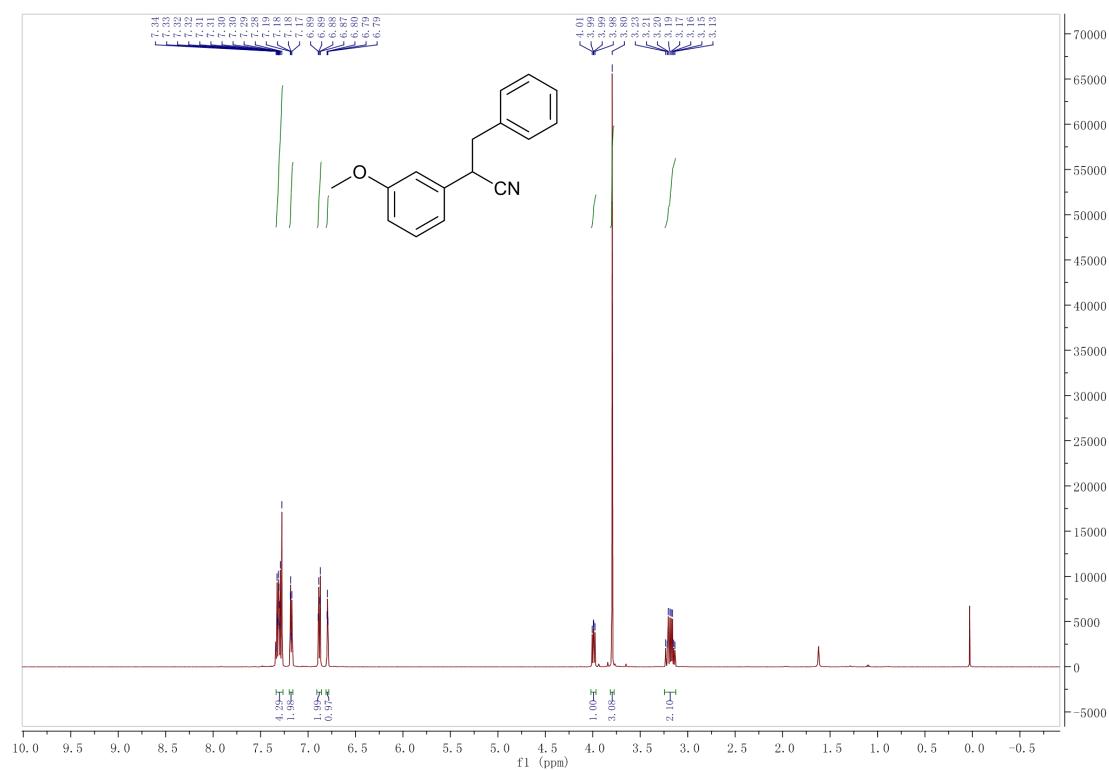
3-phenyl-2-(m-tolyl)propanenitrile (3bb**)**



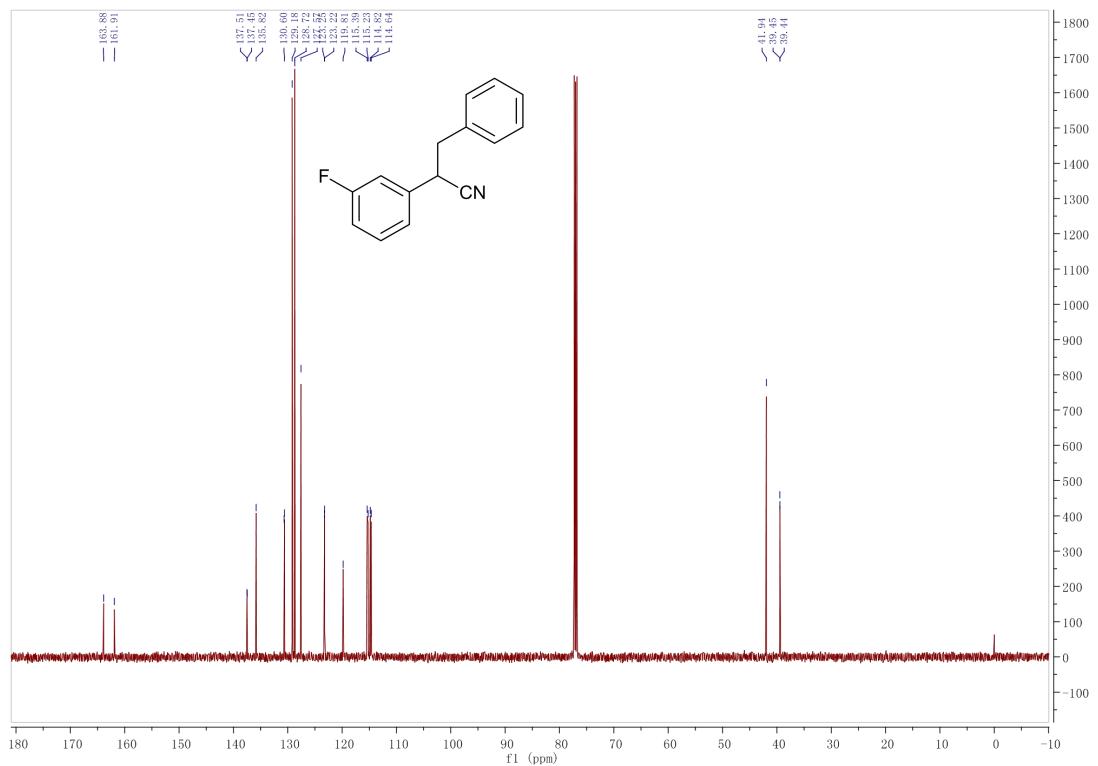
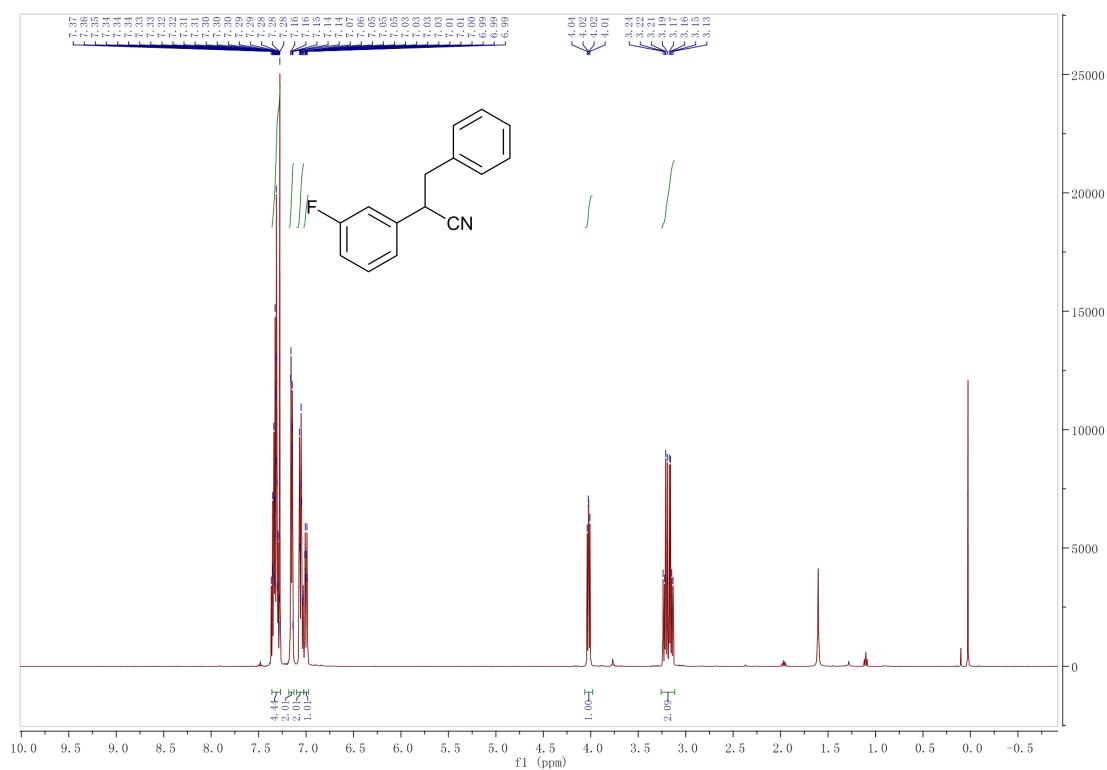
2-(4-(tert-butyl)phenyl)-3-phenylpropanenitrile (3bc**)**



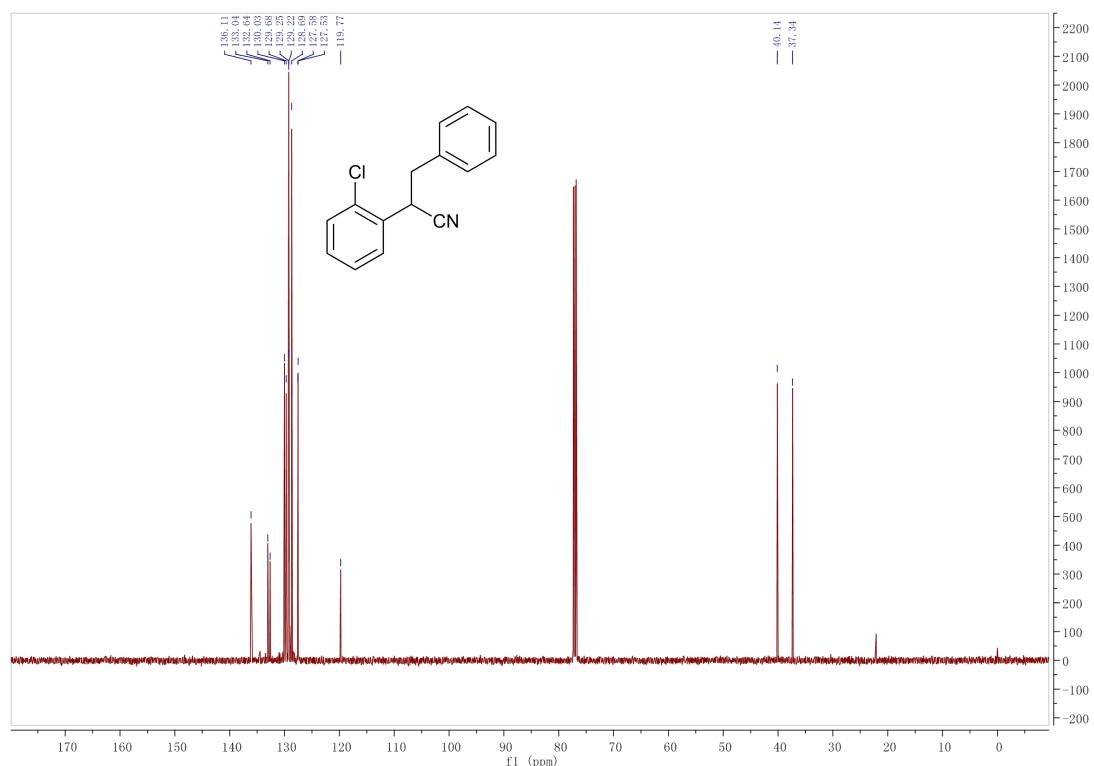
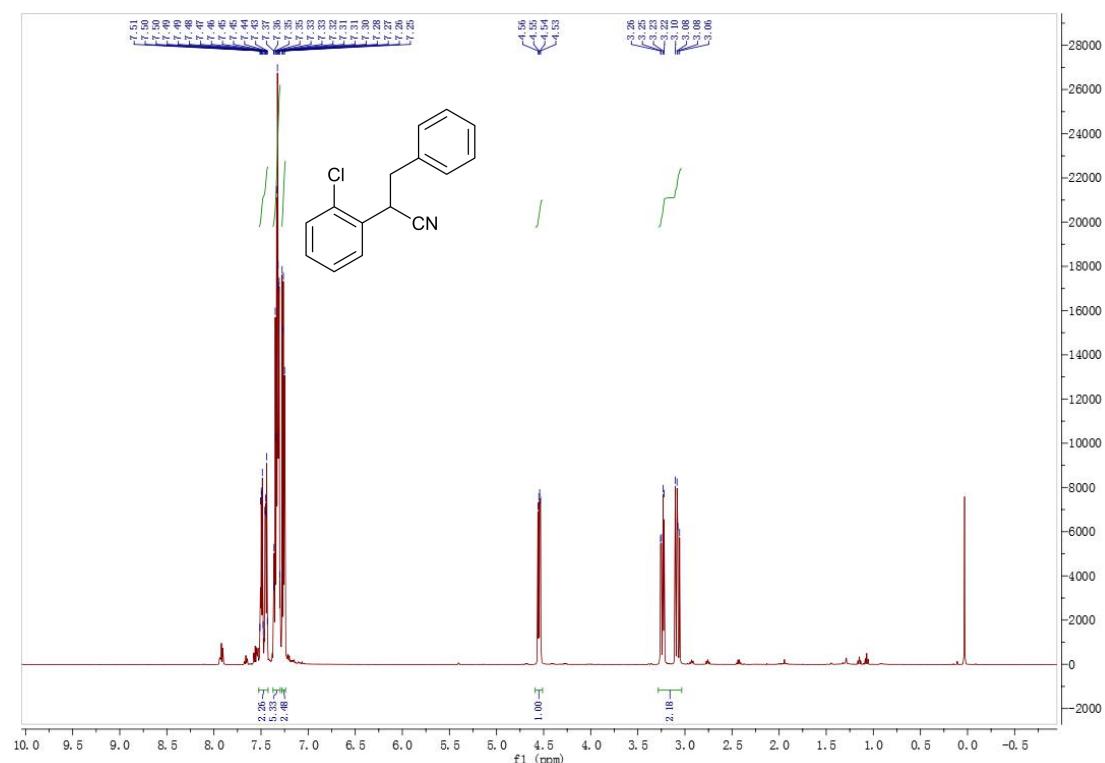
2-(3-methoxyphenyl)-3-phenylpropanenitrile (3bd**)**



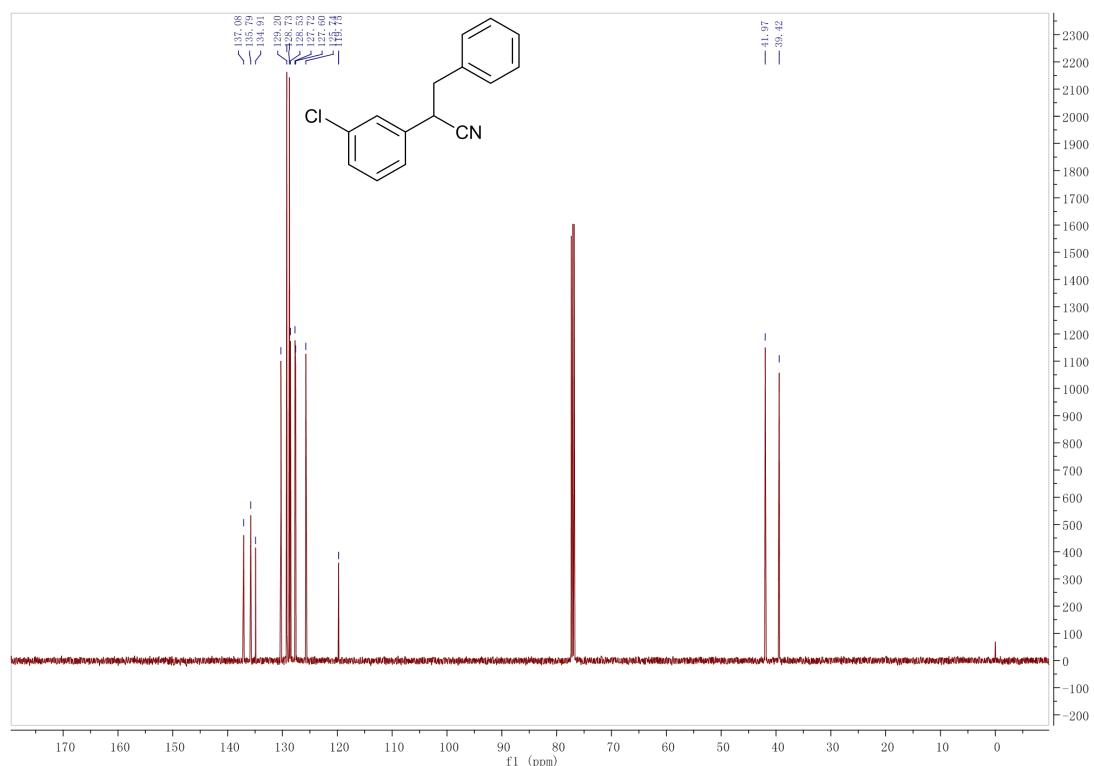
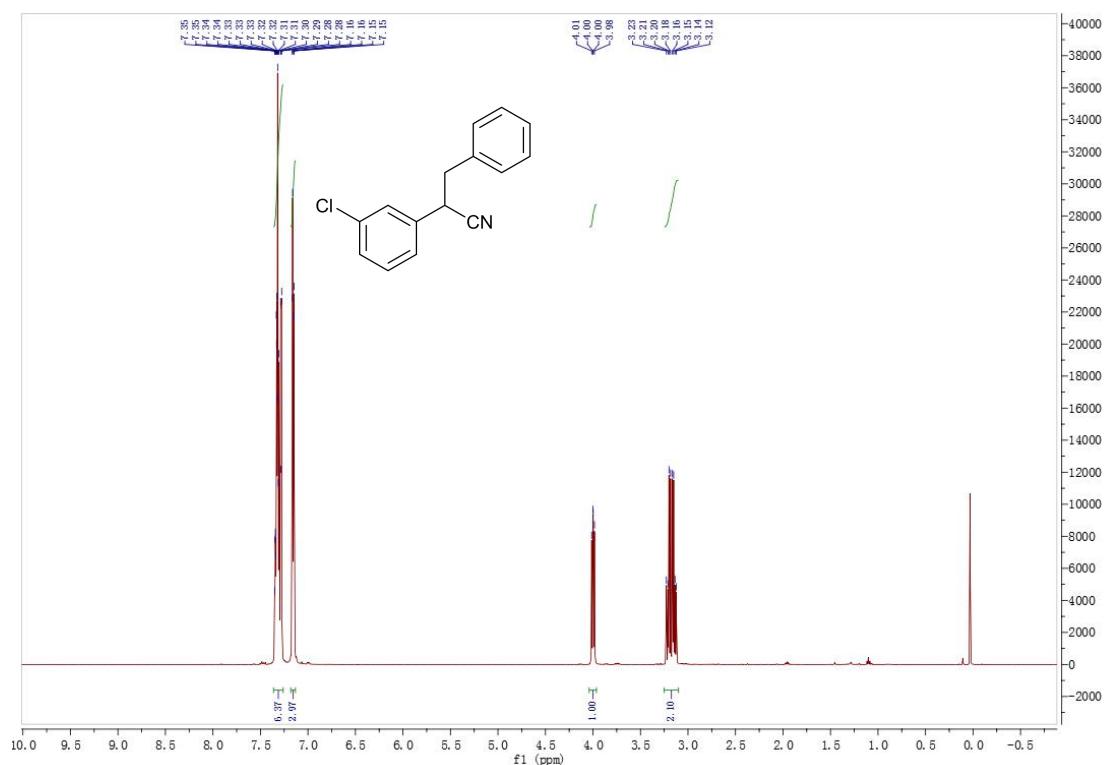
2-(3-fluorophenyl)-3-phenylpropanenitrile (**3be**)



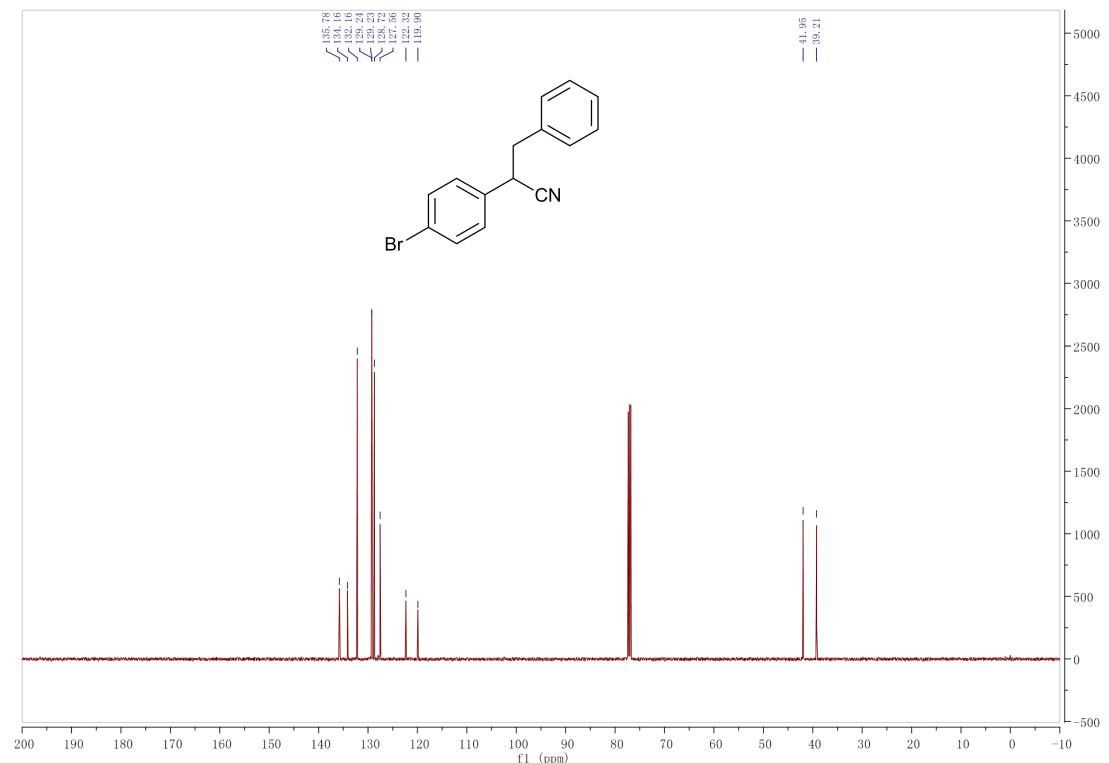
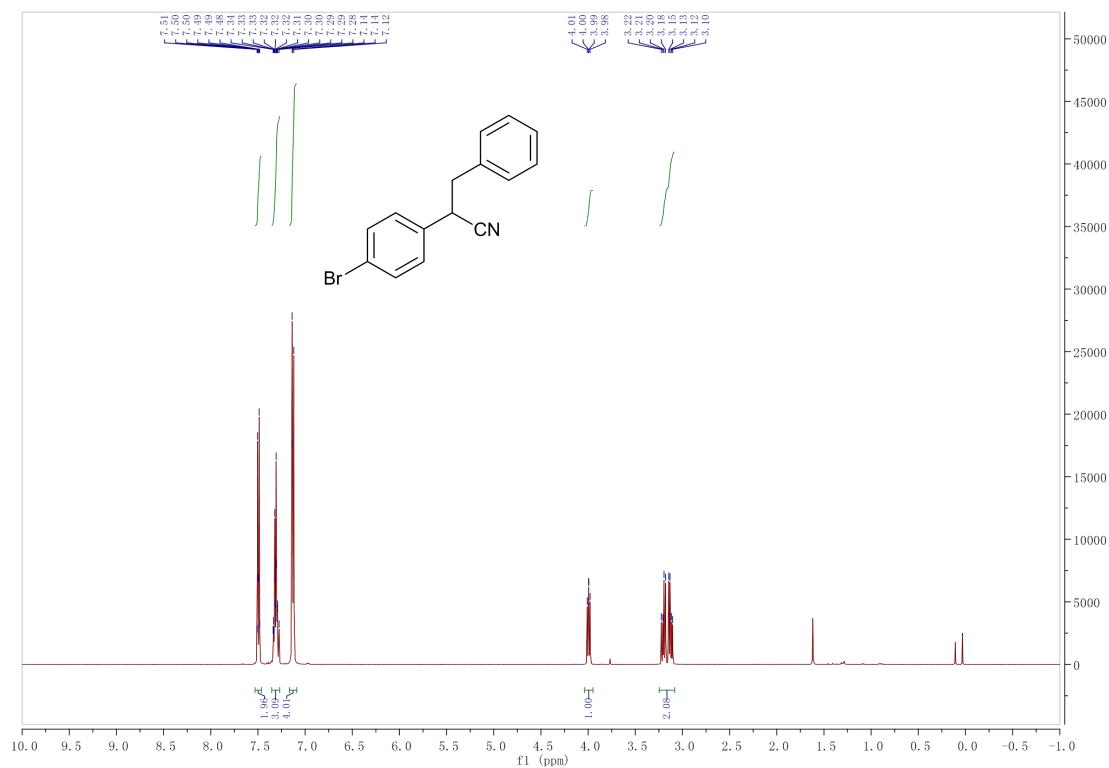
2-(2-chlorophenyl)-3-phenylpropanenitrile (3bf**)**



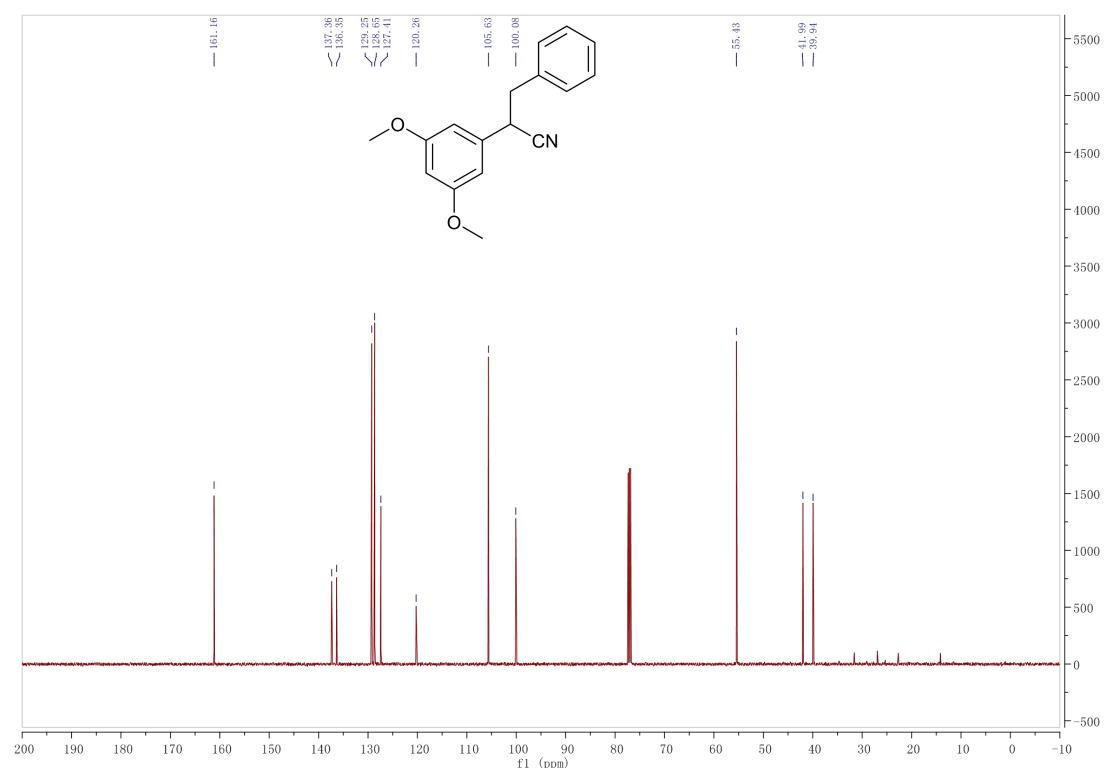
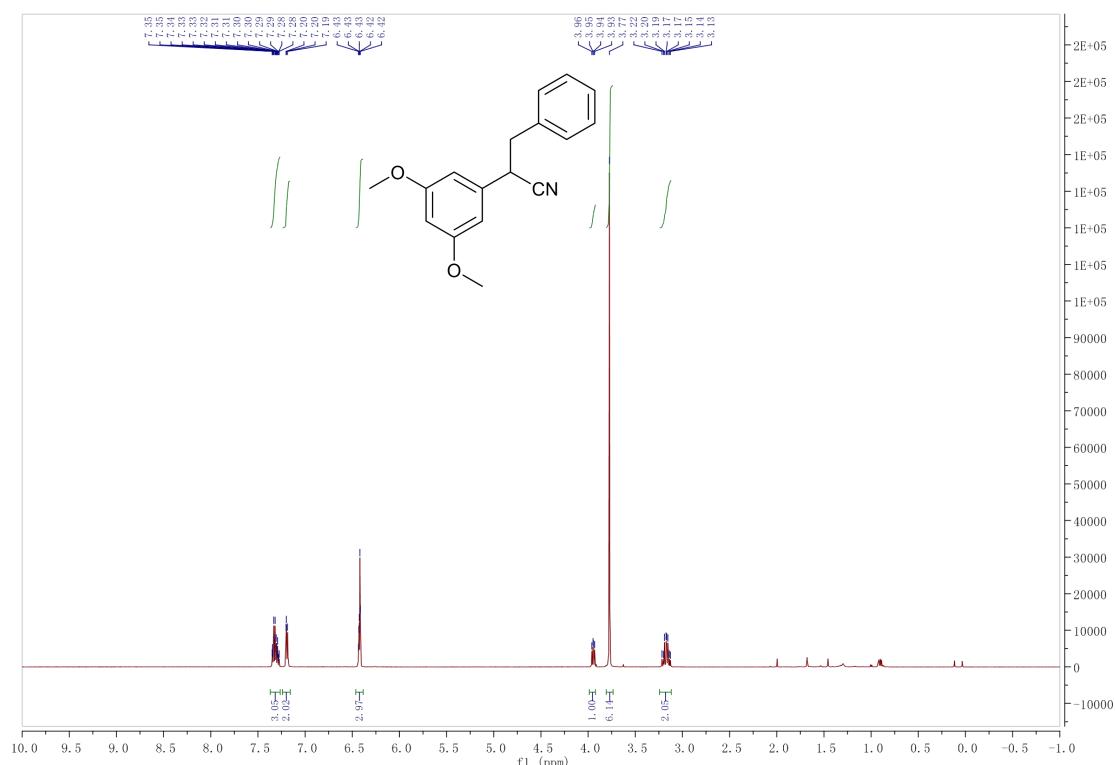
2-(3-chlorophenyl)-3-phenylpropanenitrile (3bg**)**



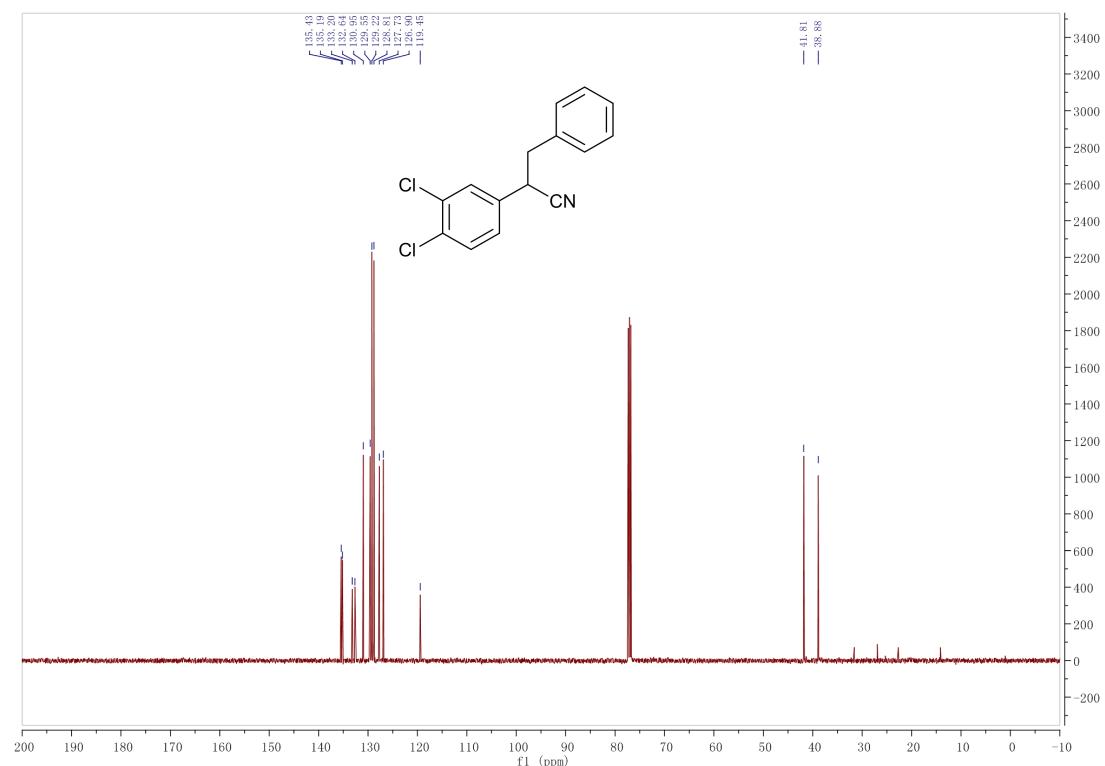
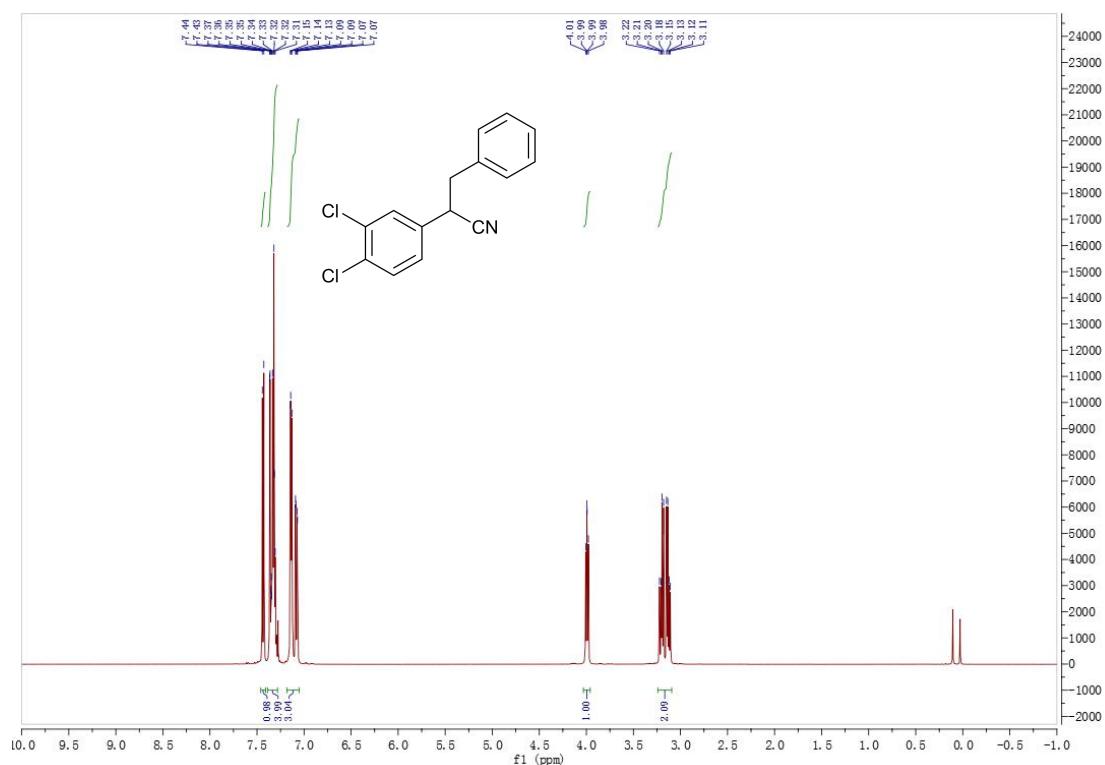
2-(4-bromophenyl)-3-phenylpropanenitrile (3bh**)**



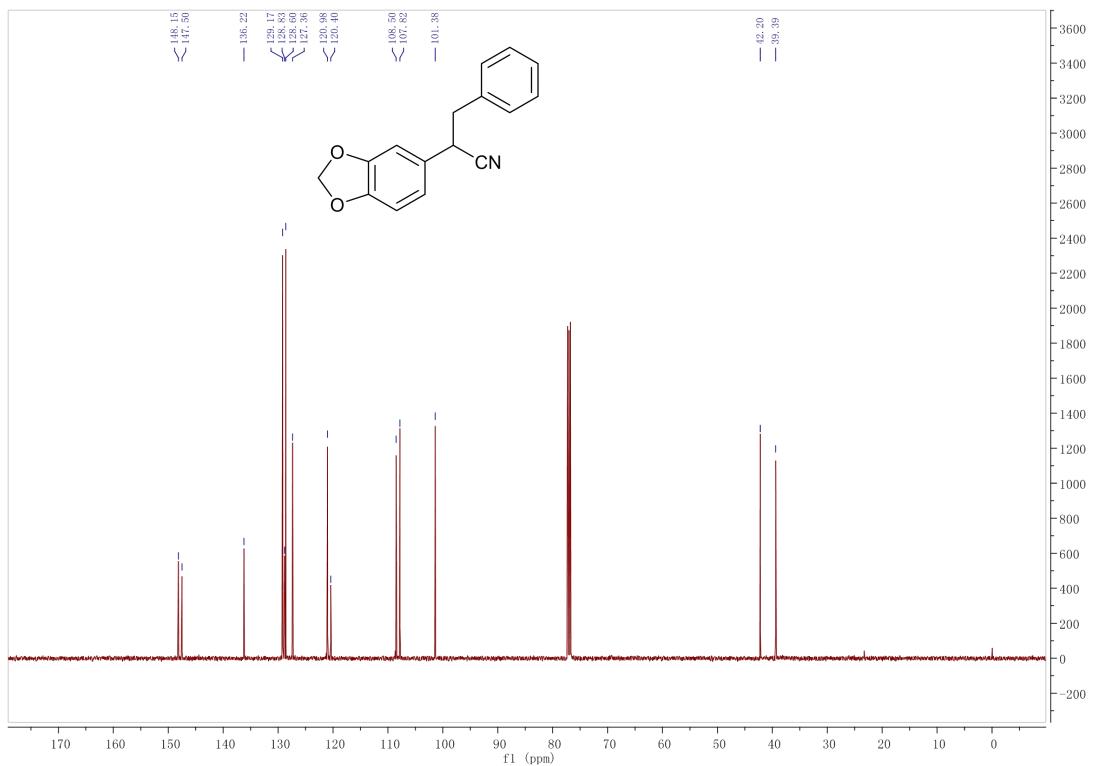
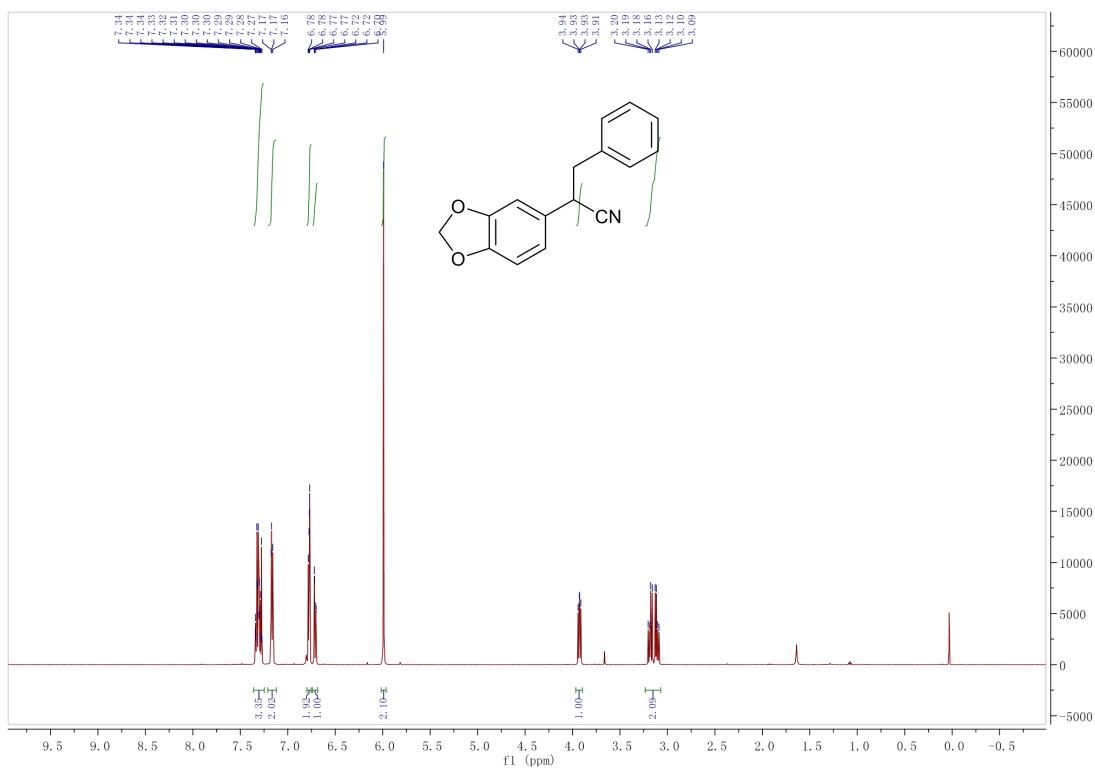
2-(3,5-dimethoxyphenyl)-3-phenylpropanenitrile (3bi**)**



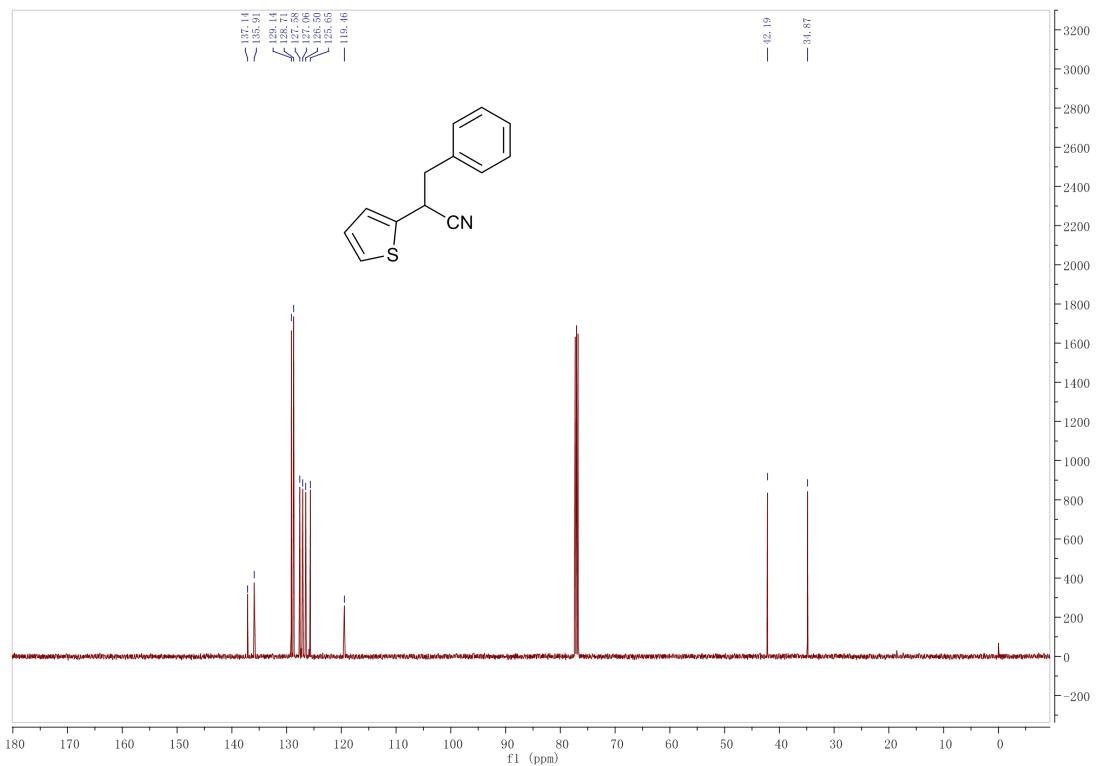
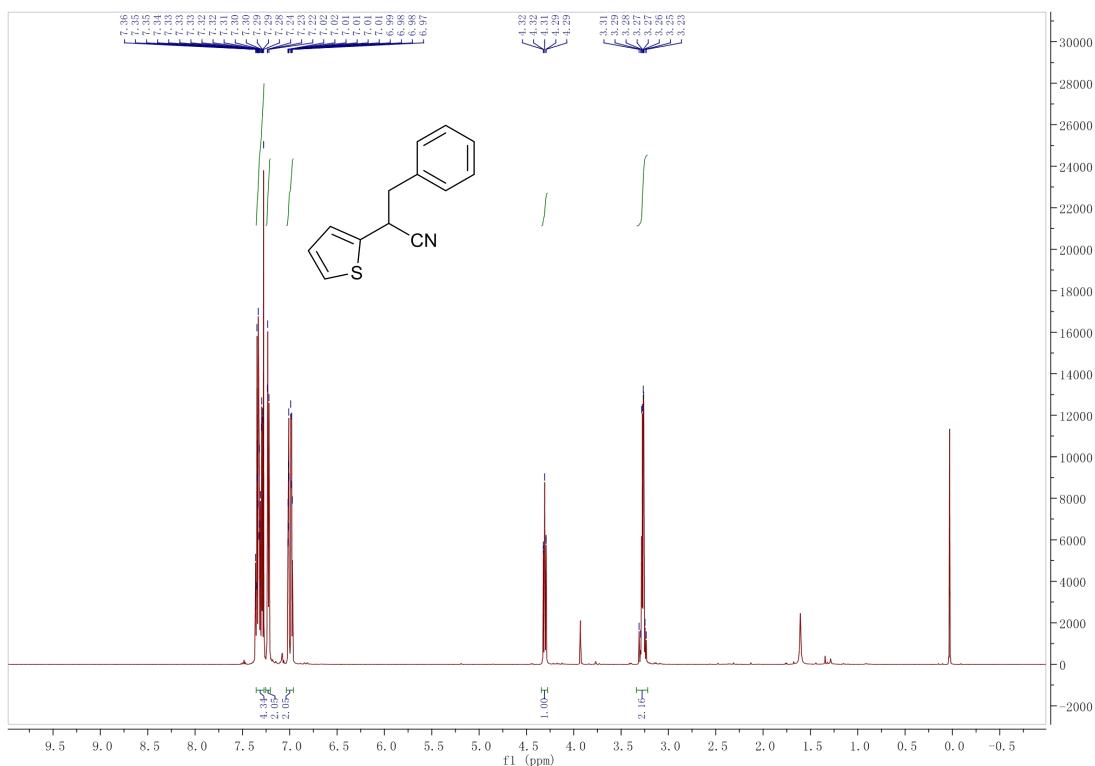
2-(3,4-dichlorophenyl)-3-phenylpropanenitrile (3bj**)**



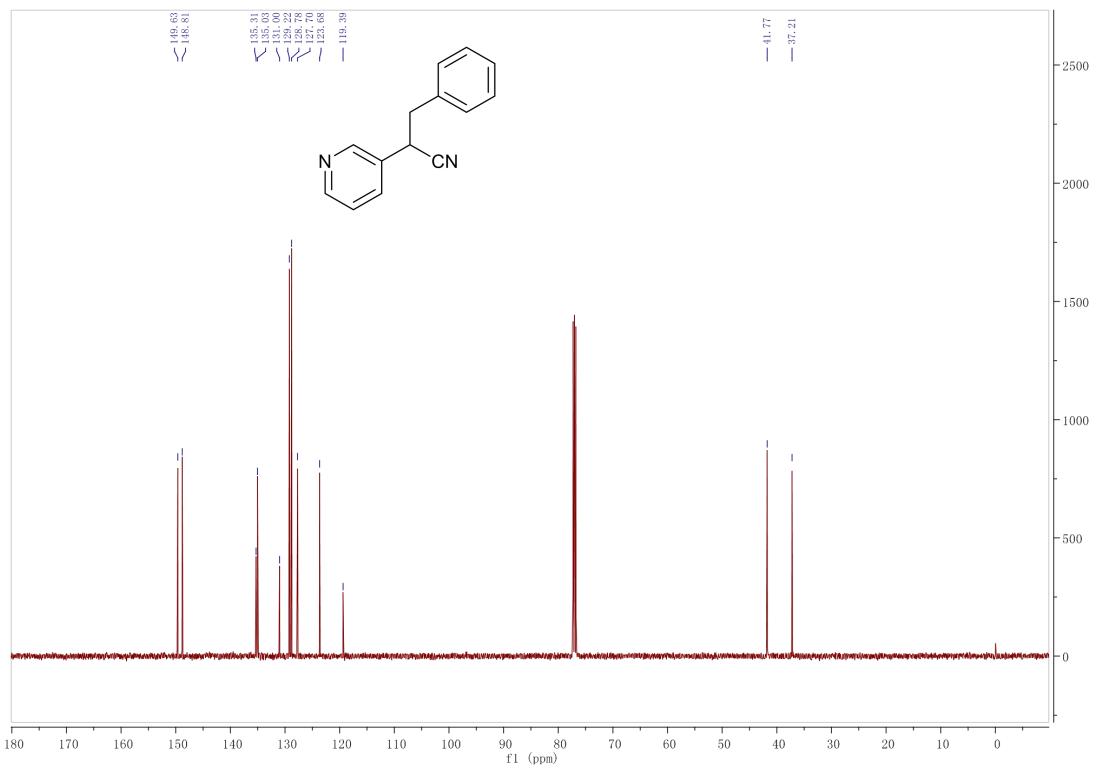
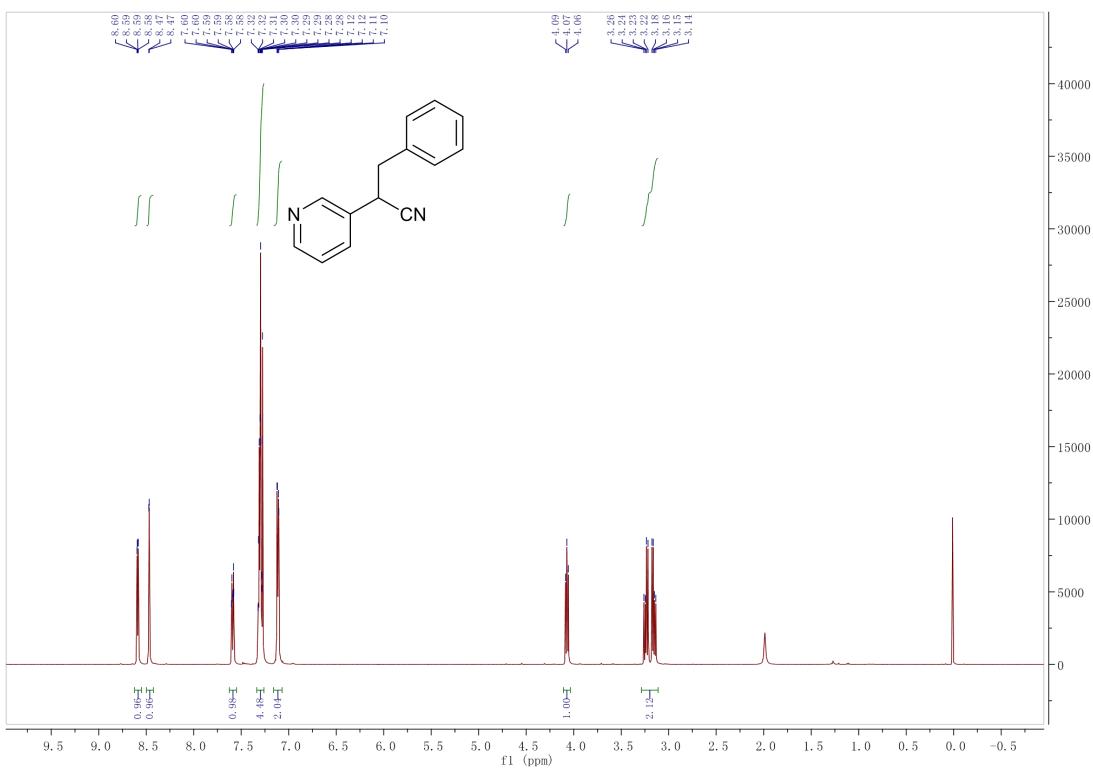
2-(benzo[d][1,3]dioxol-5-yl)-3-phenylpropanenitrile (3bk**)**



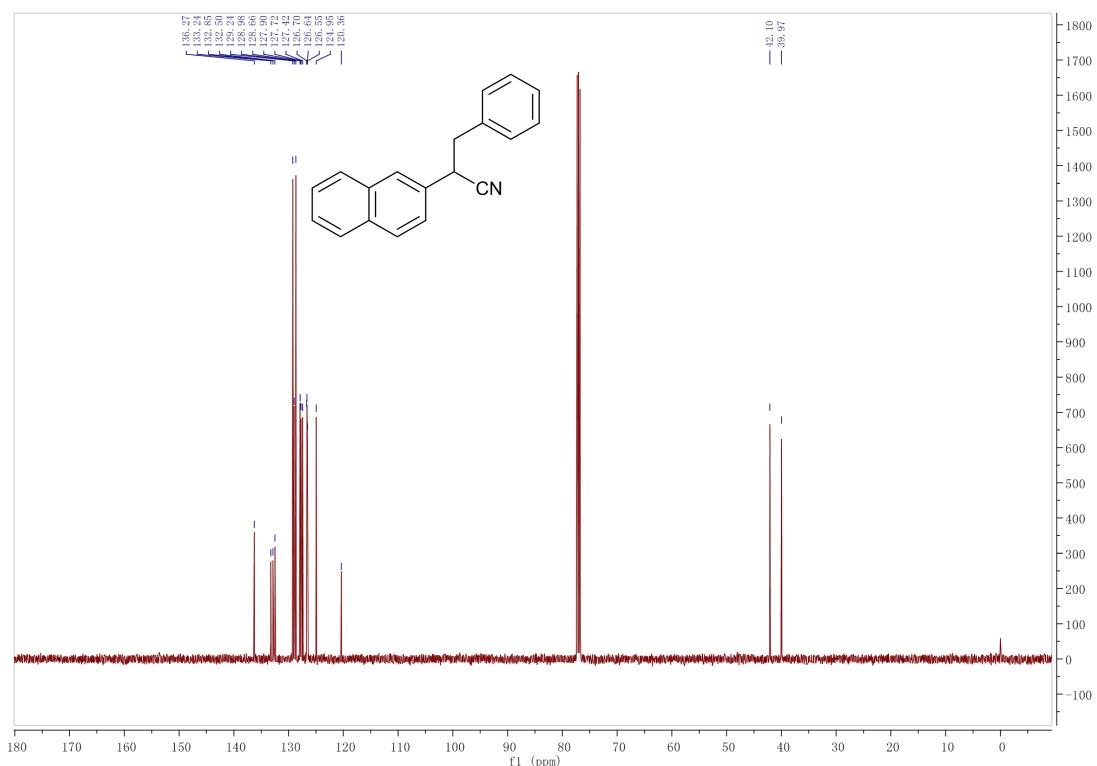
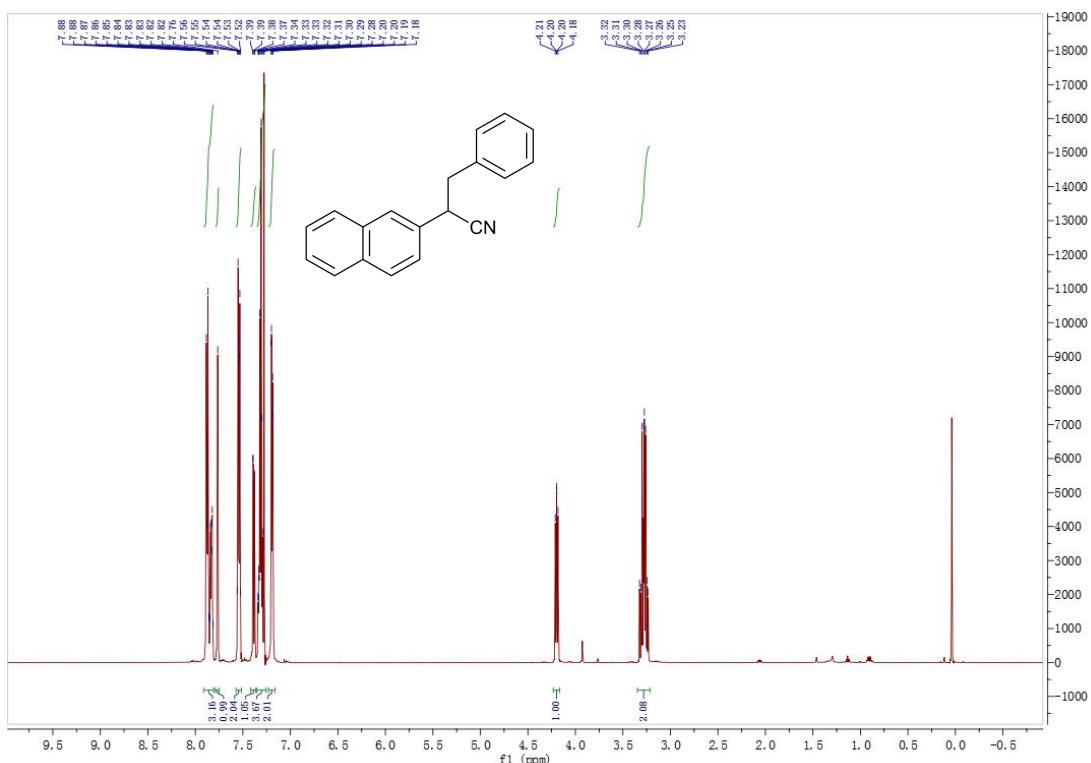
3-phenyl-2-(thiophen-2-yl)propanenitrile (**3bl**)



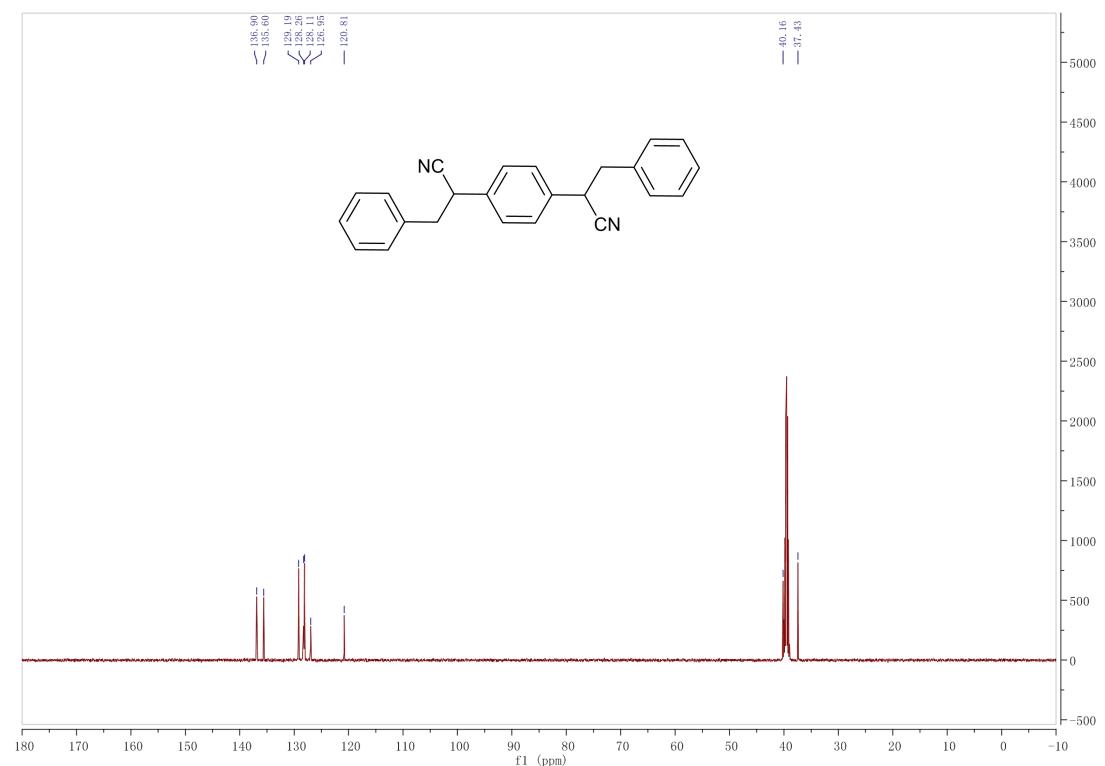
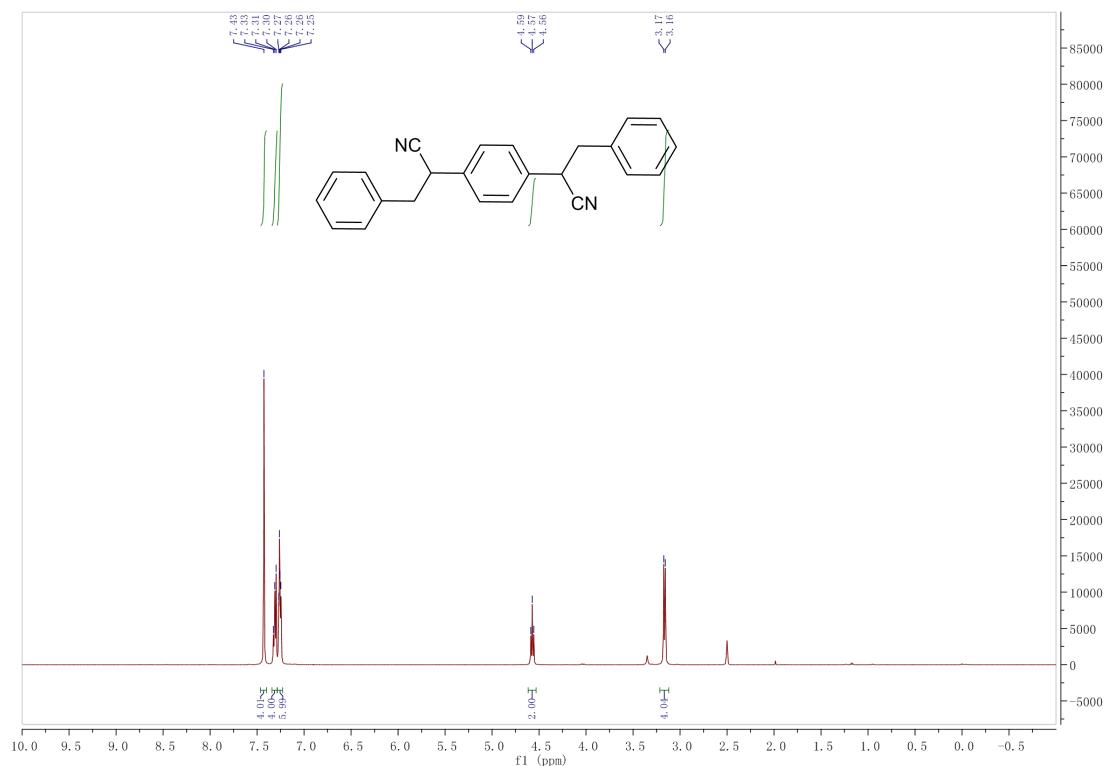
3-phenyl-2-(pyridin-3-yl)propanenitrile (3bm**)**



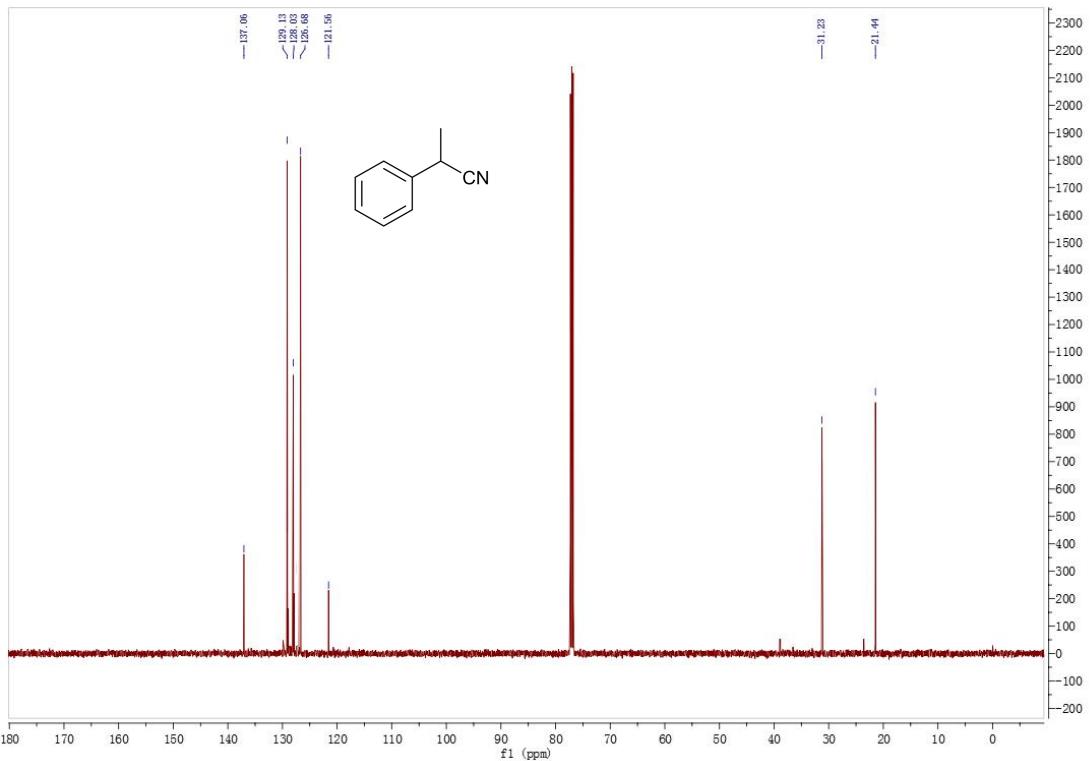
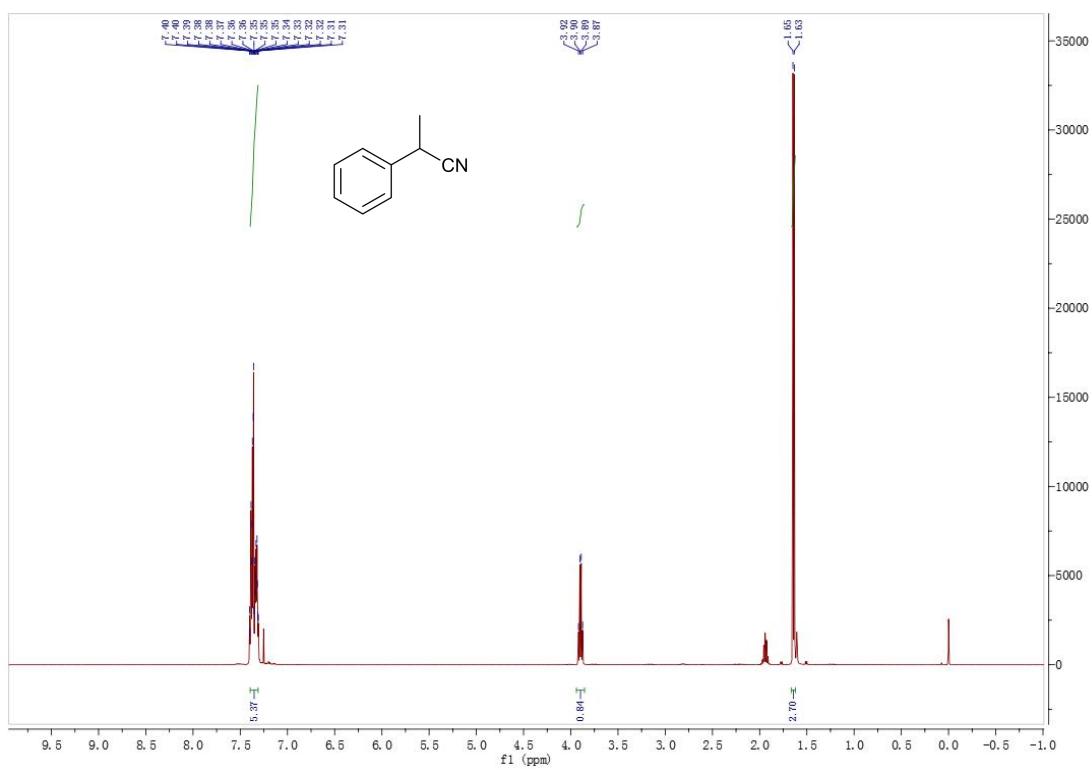
2-(naphthalen-2-yl)-3-phenylpropanenitrile (3bn**)**



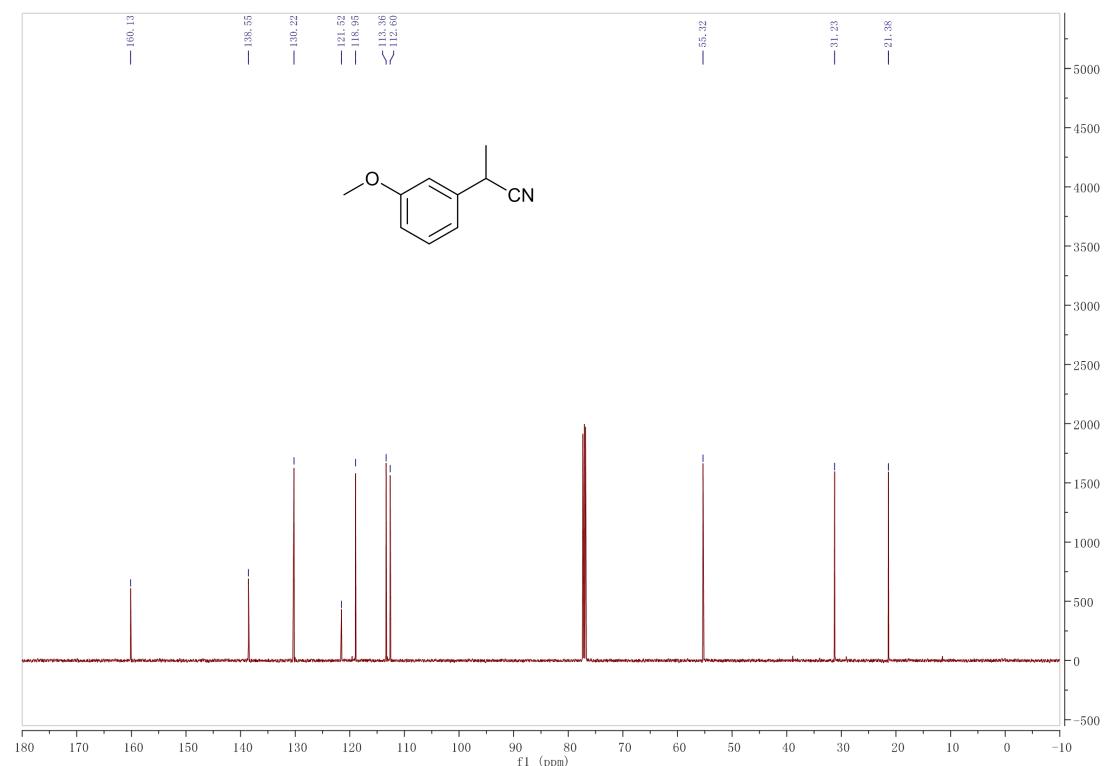
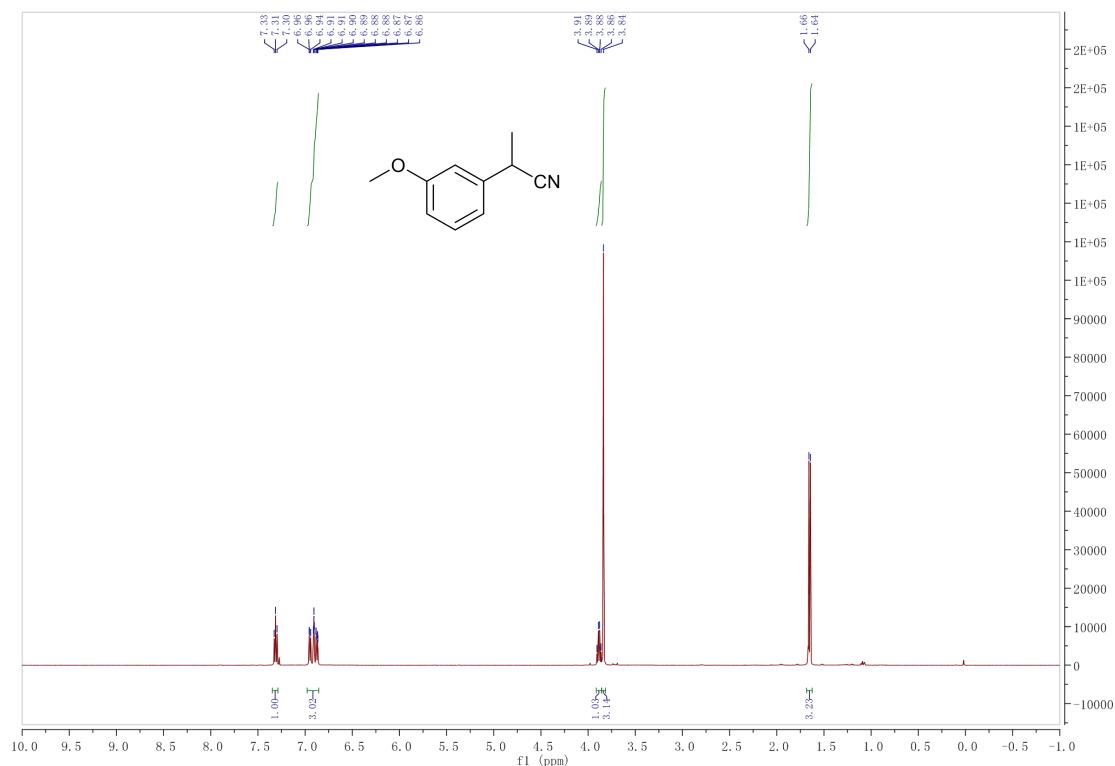
2,2'-(1,4-phenylene)bis(3-phenylpropanenitrile) (3bo**)**



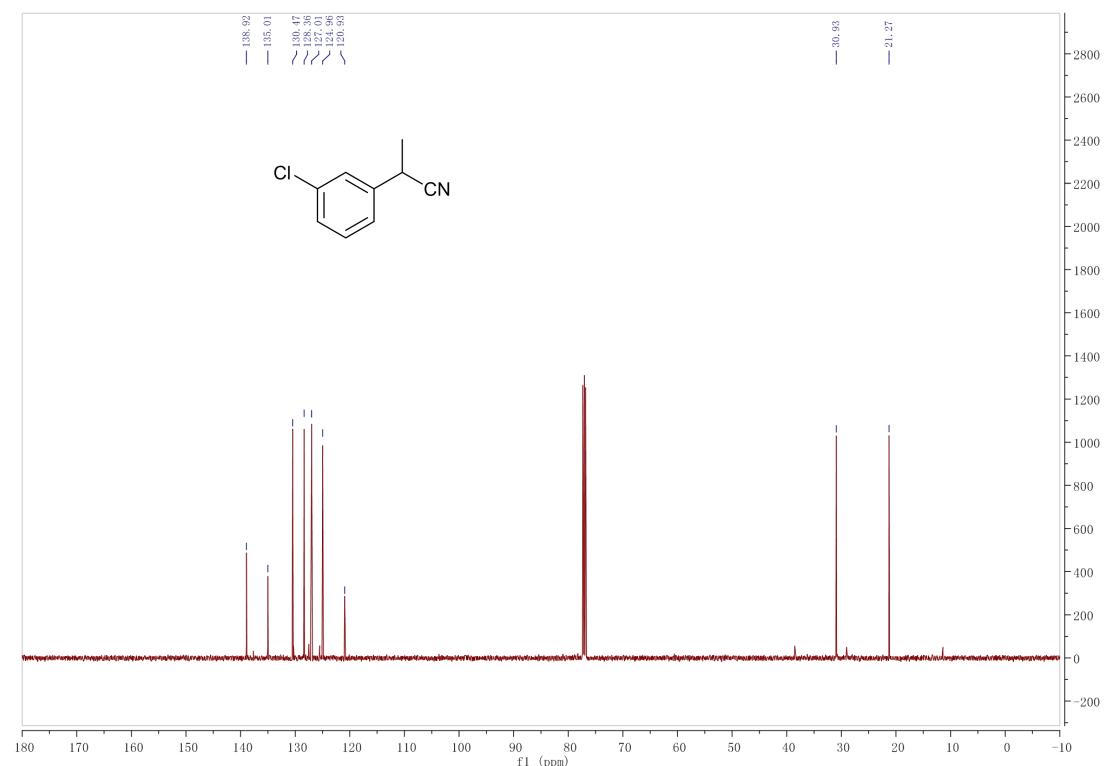
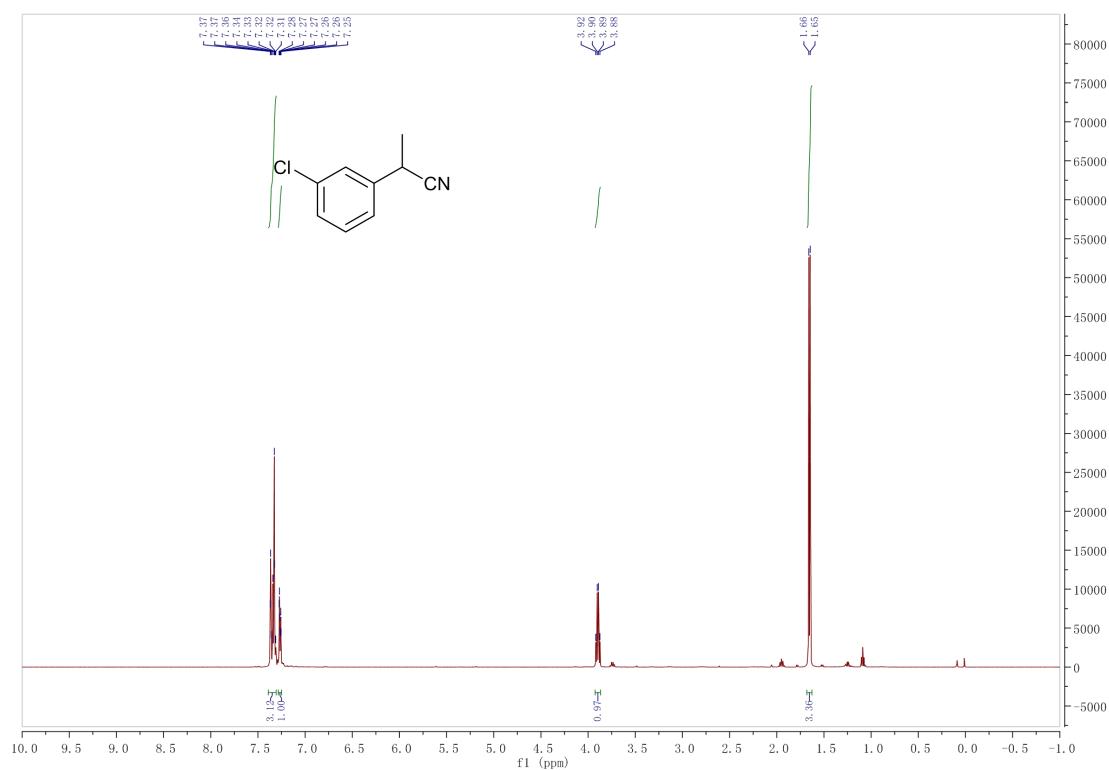
2-phenylpropanenitrile (3ca**)**



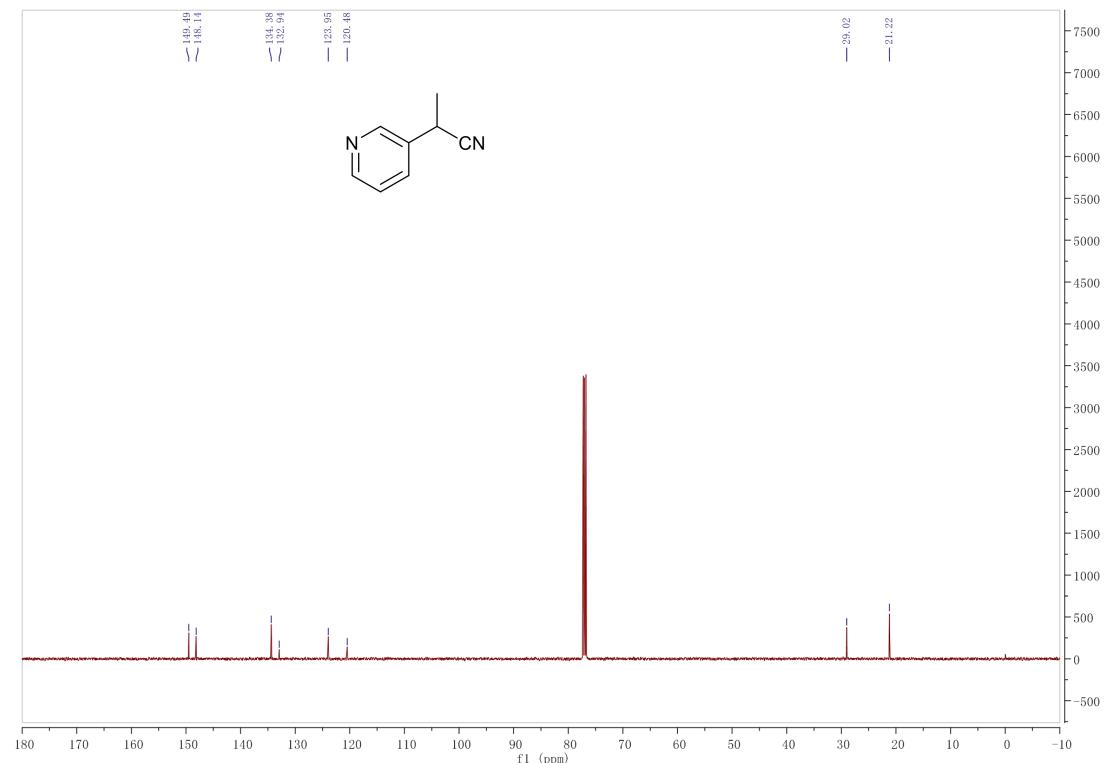
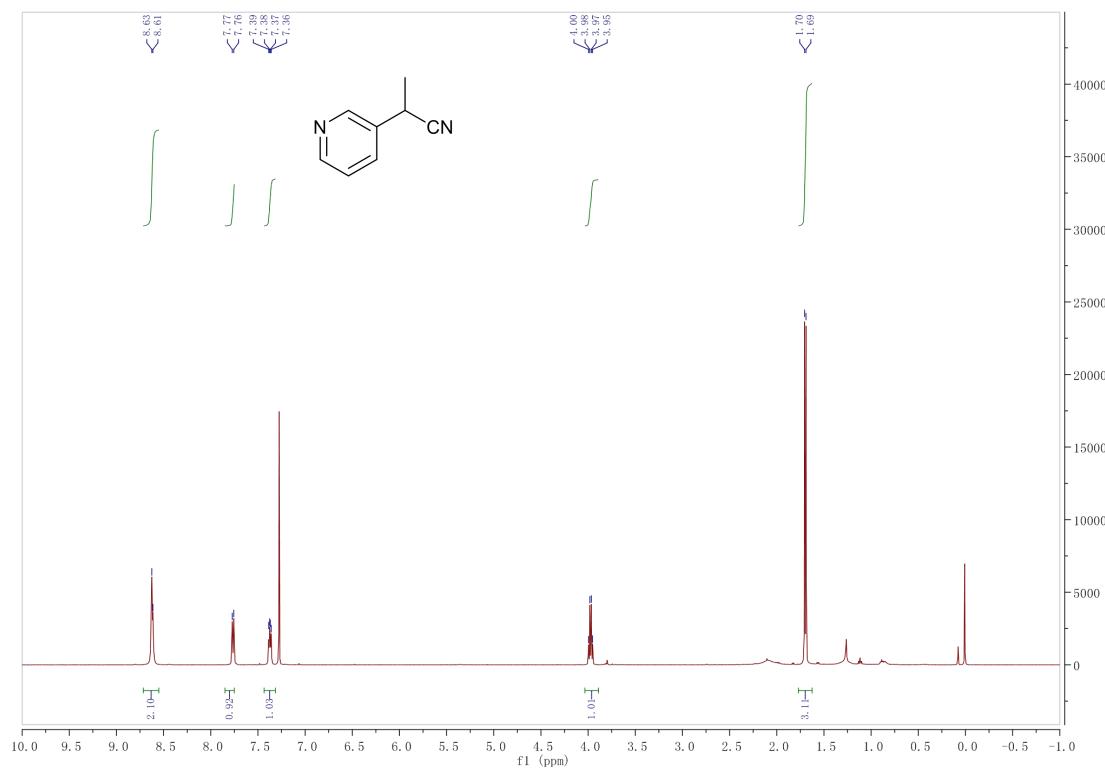
2-(3-methoxyphenyl)propanenitrile (3cb**)**



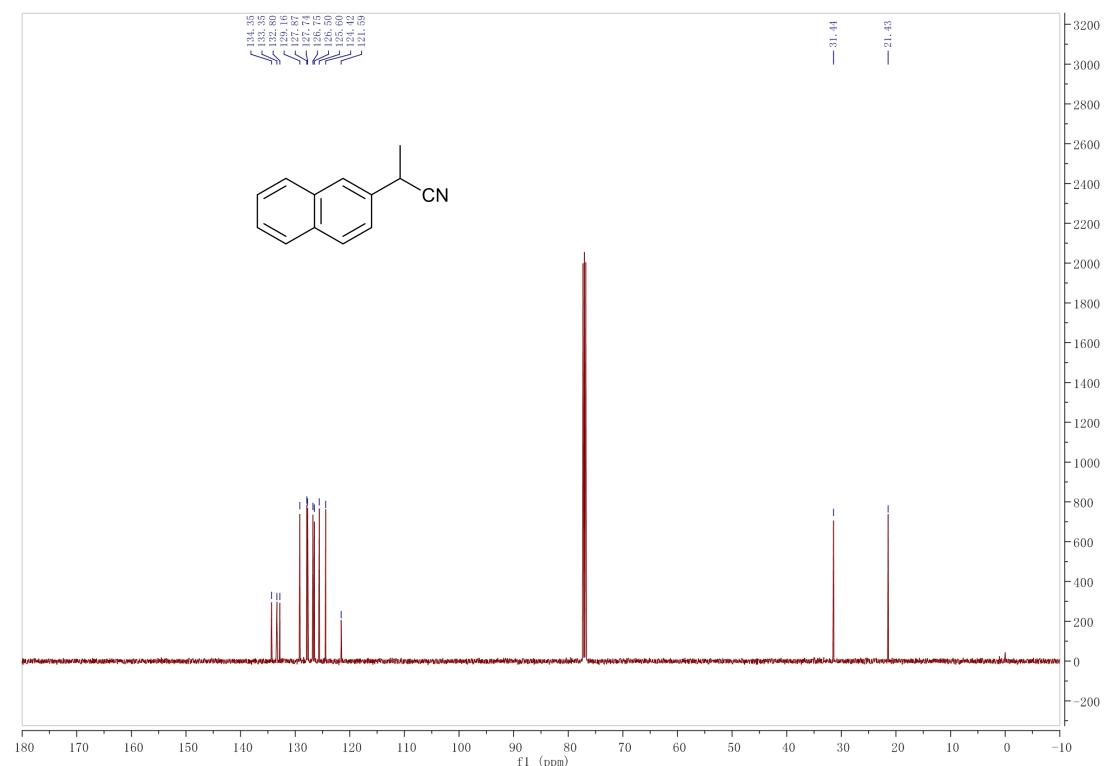
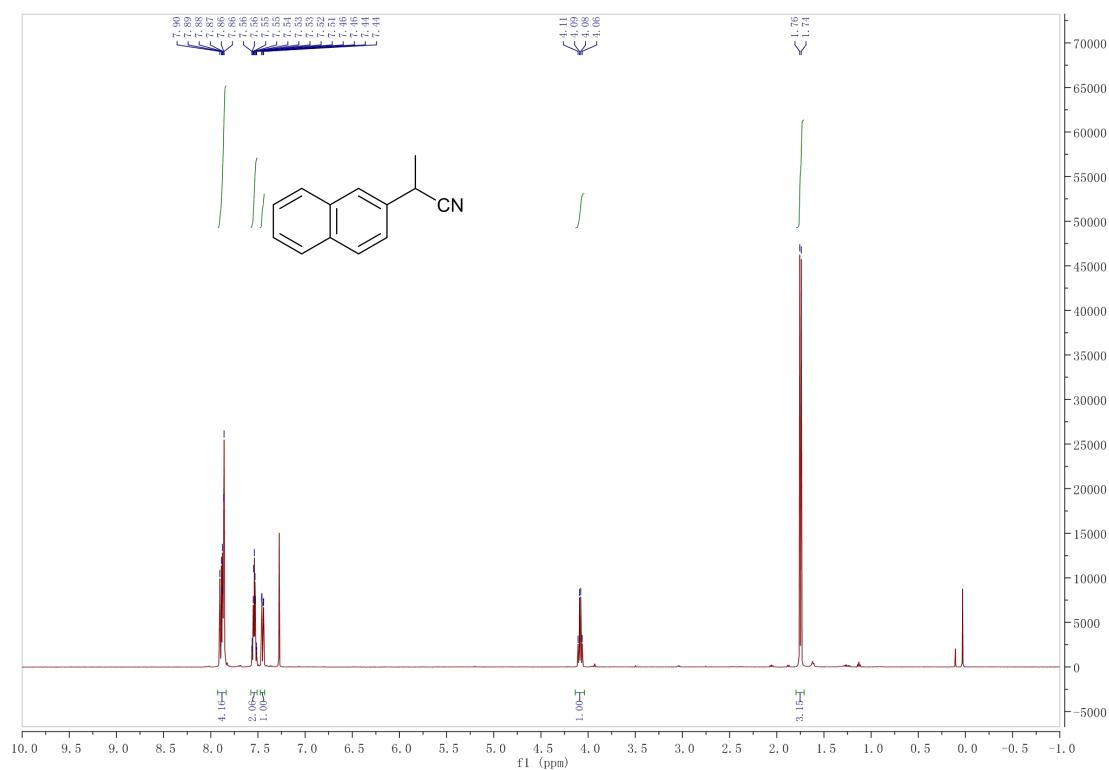
2-(3-chlorophenyl)propanenitrile (3cc**)**



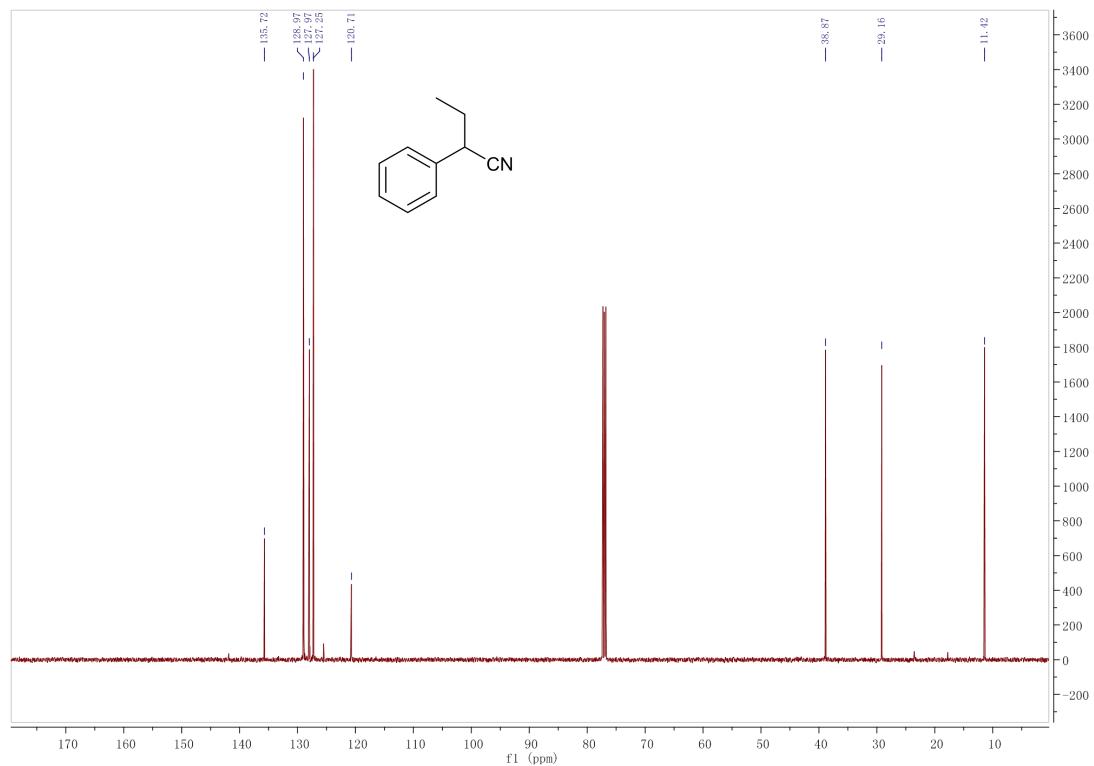
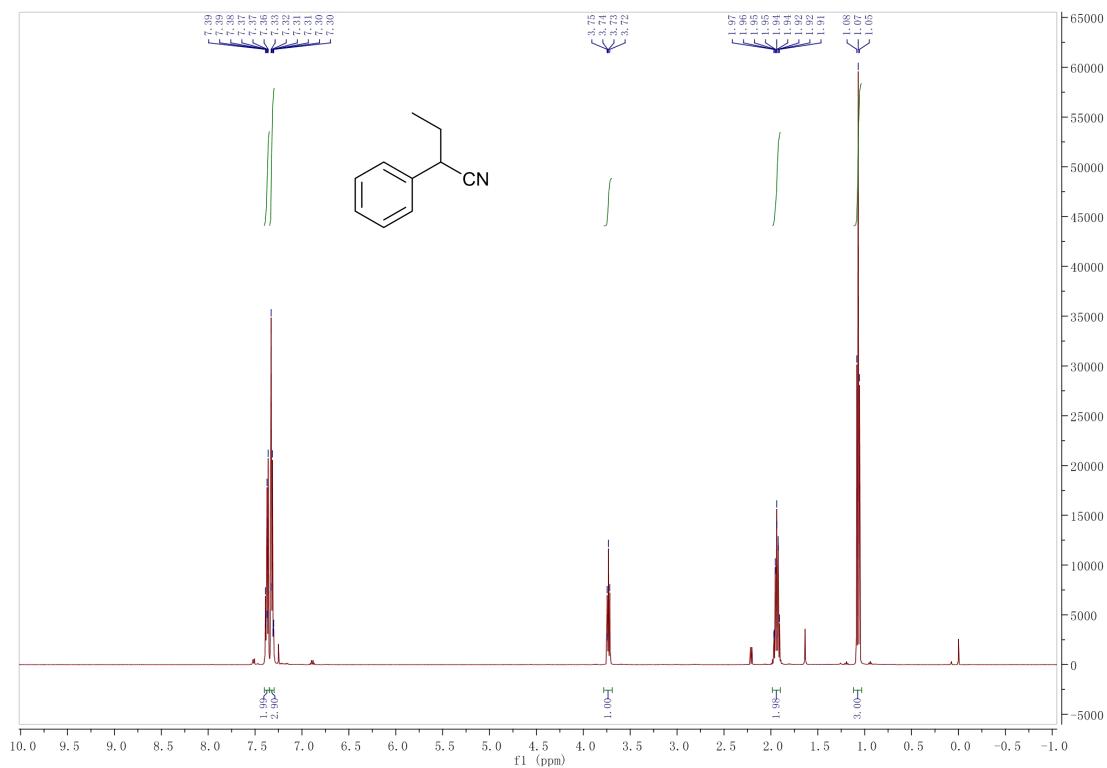
2-(pyridin-3-yl)propanenitrile (3cd**)**



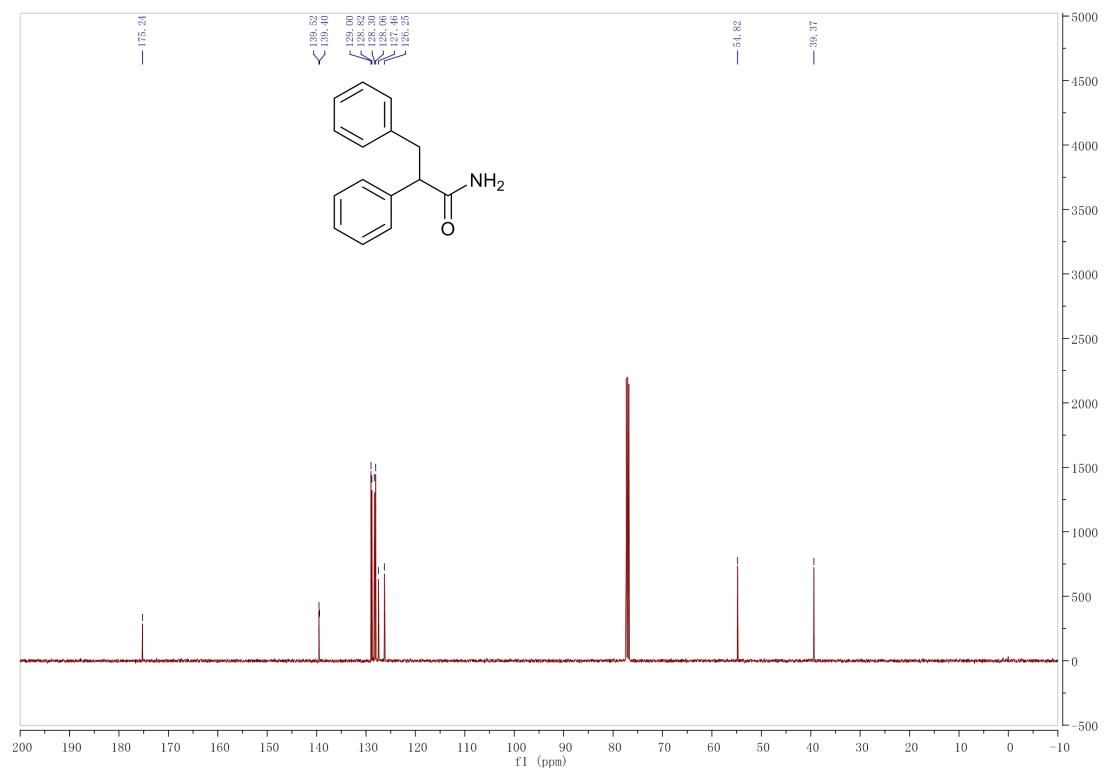
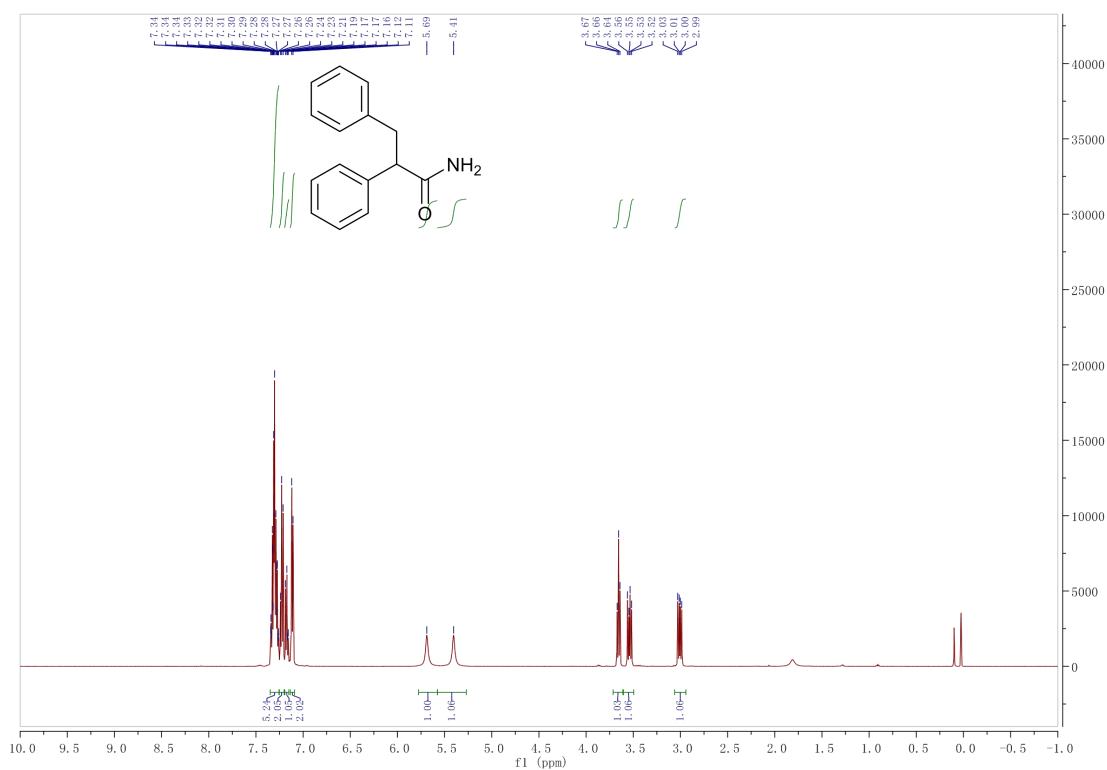
2-(naphthalen-2-yl)propanenitrile (3ce**)**



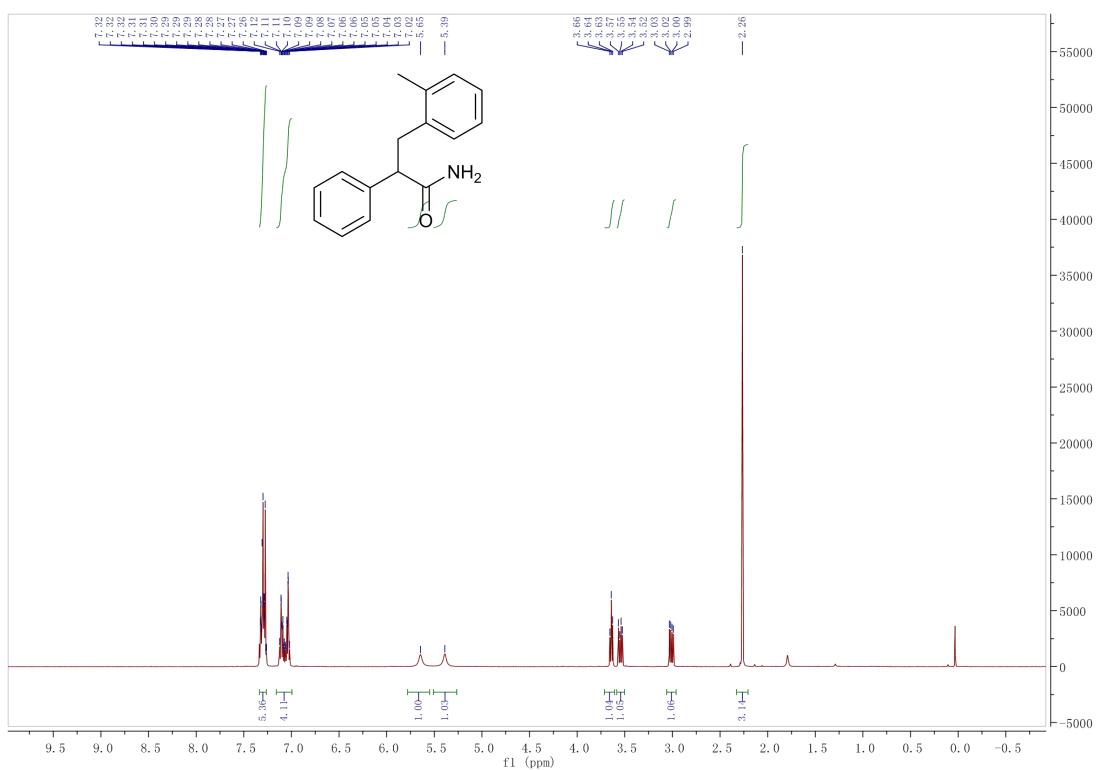
2-phenylbutanenitrile (3cf**)**



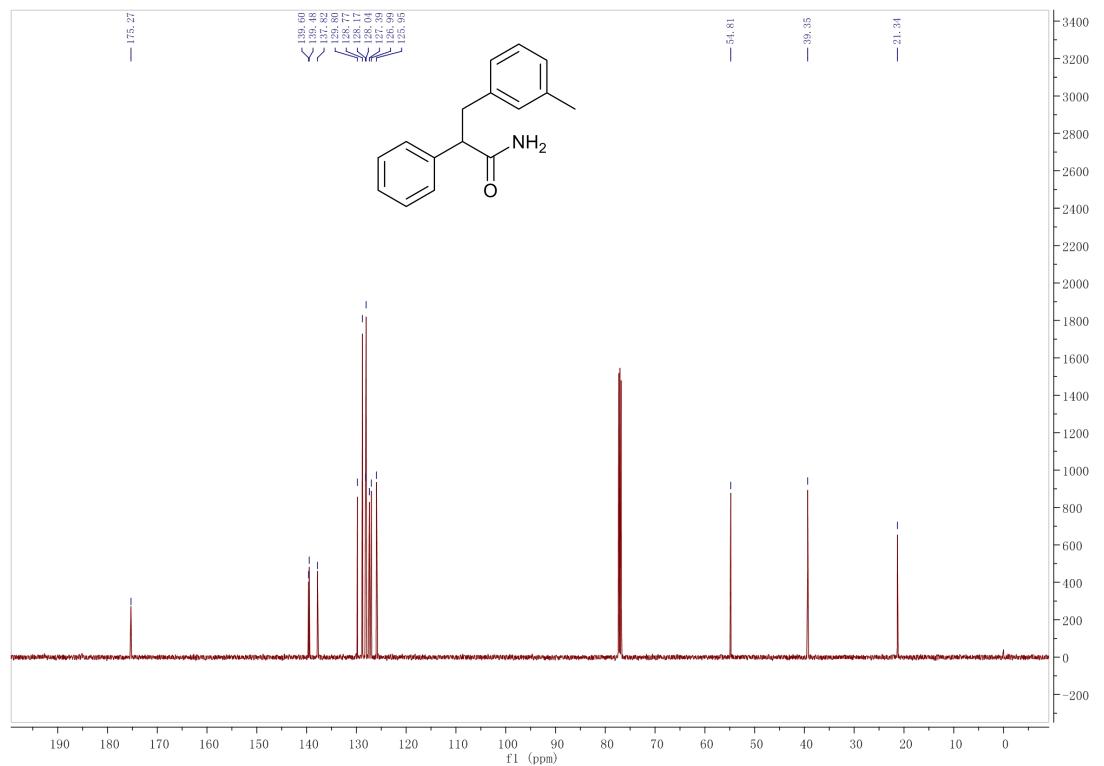
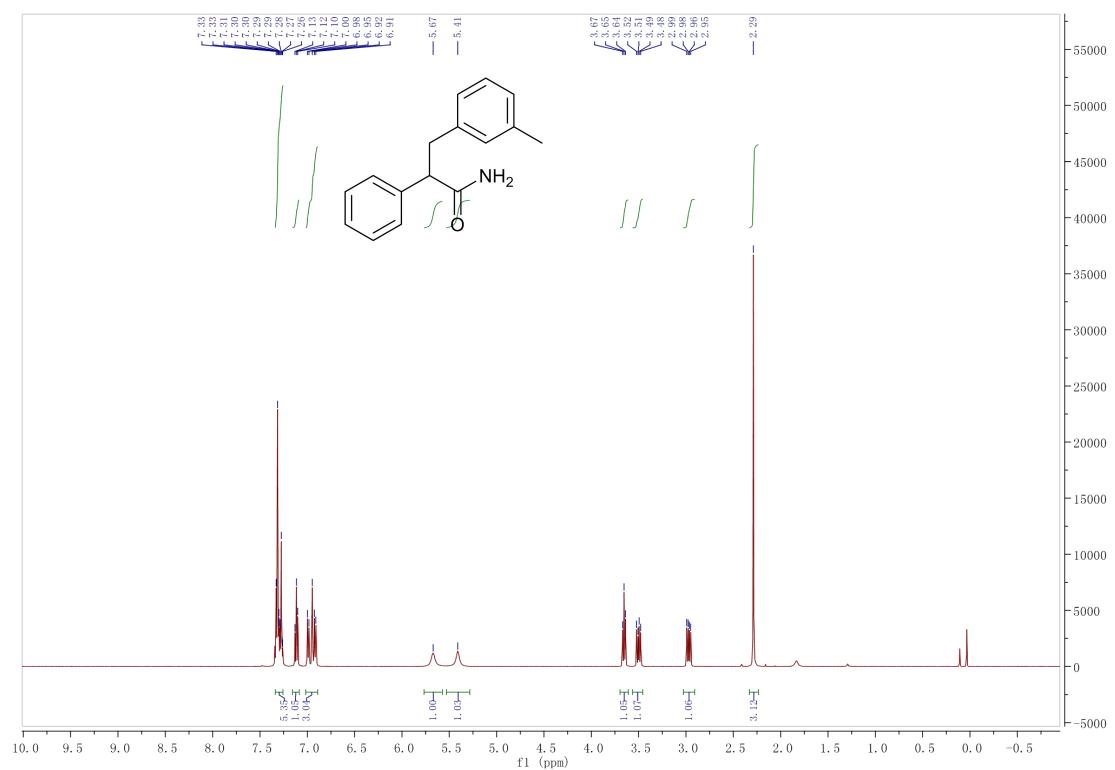
2,3-diphenylpropanamide (**4aa**)



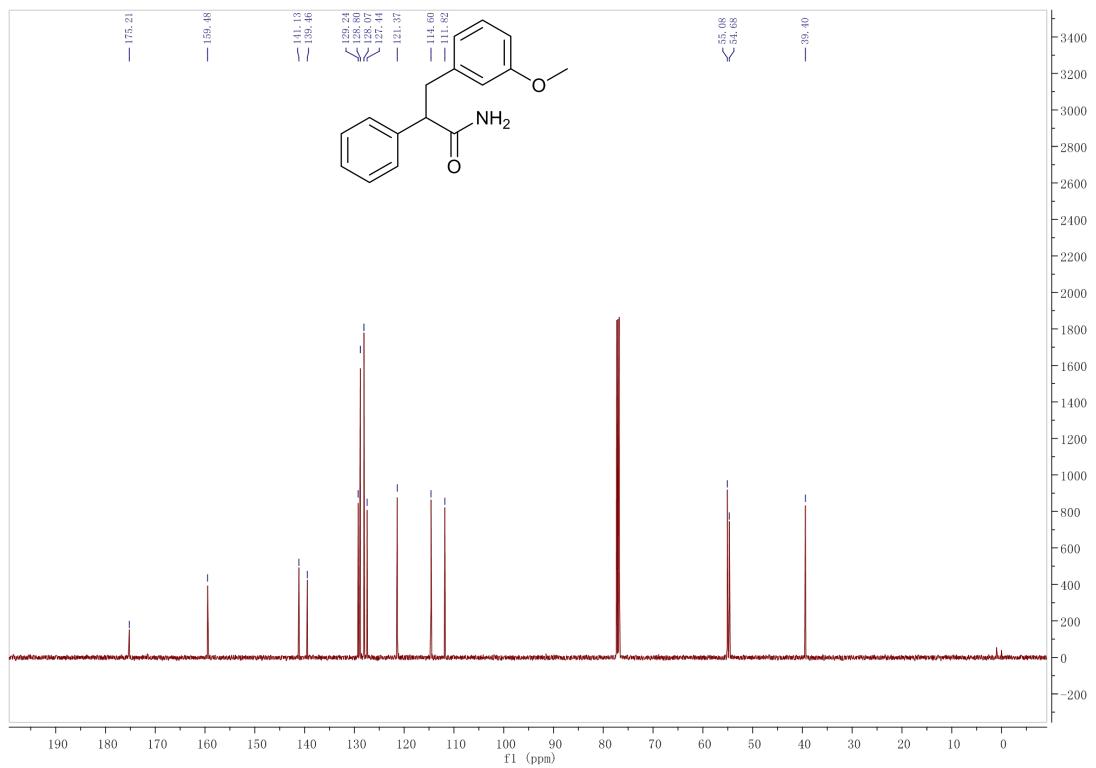
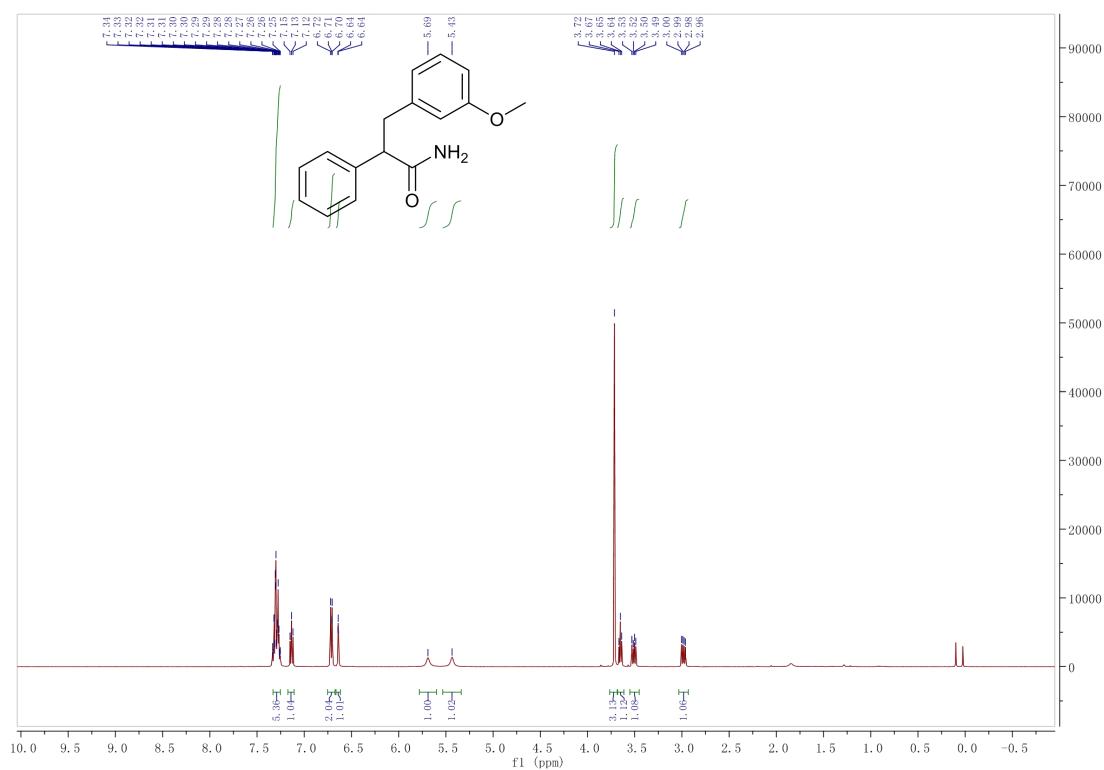
2-phenyl-3-(o-tolyl)propenamide (4ab**)**



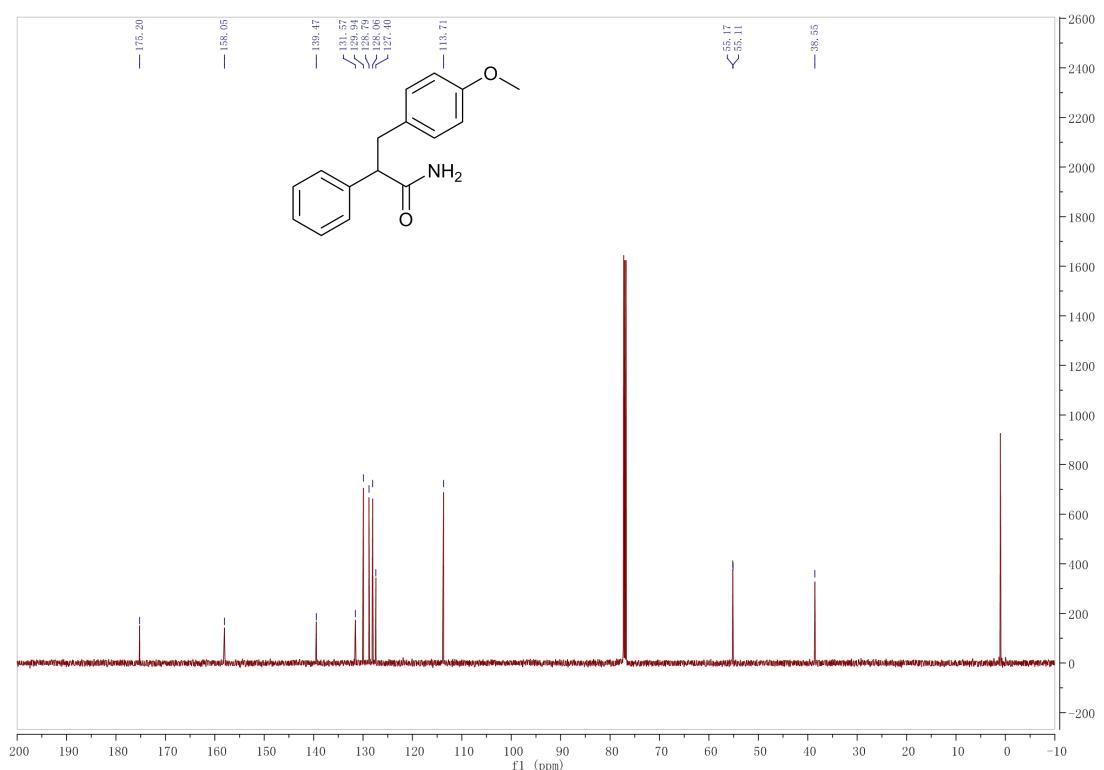
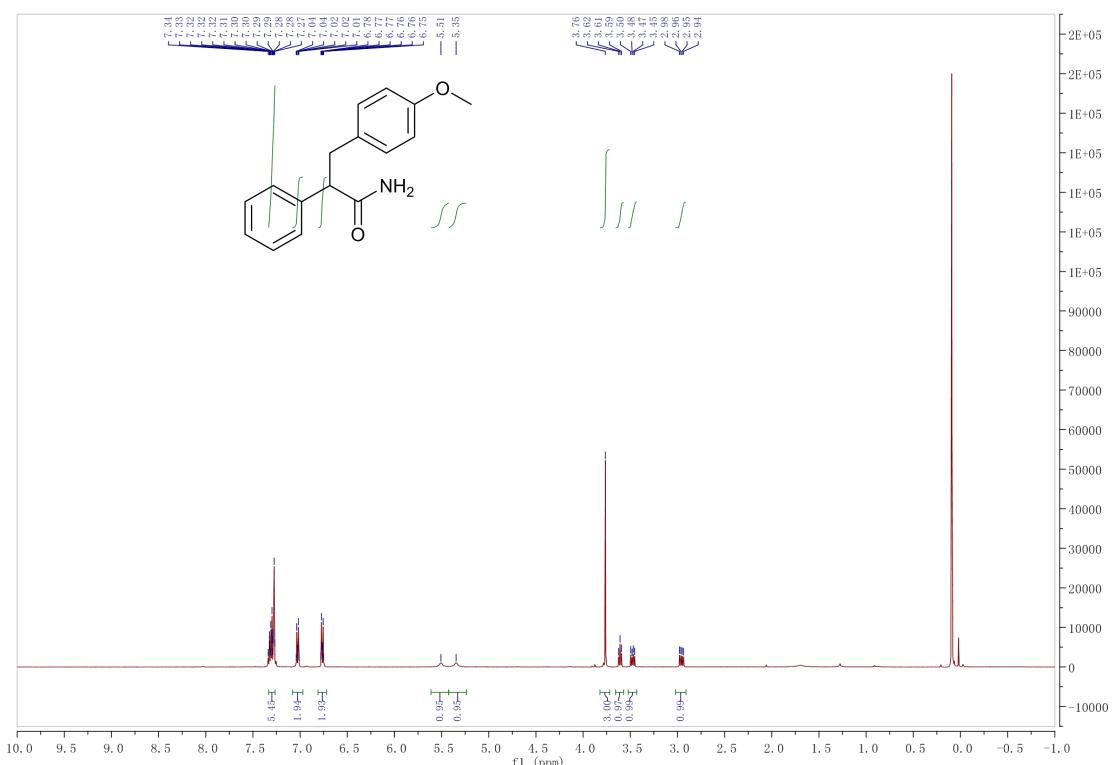
2-phenyl-3-(m-tolyl)propenamide (4ac**)**



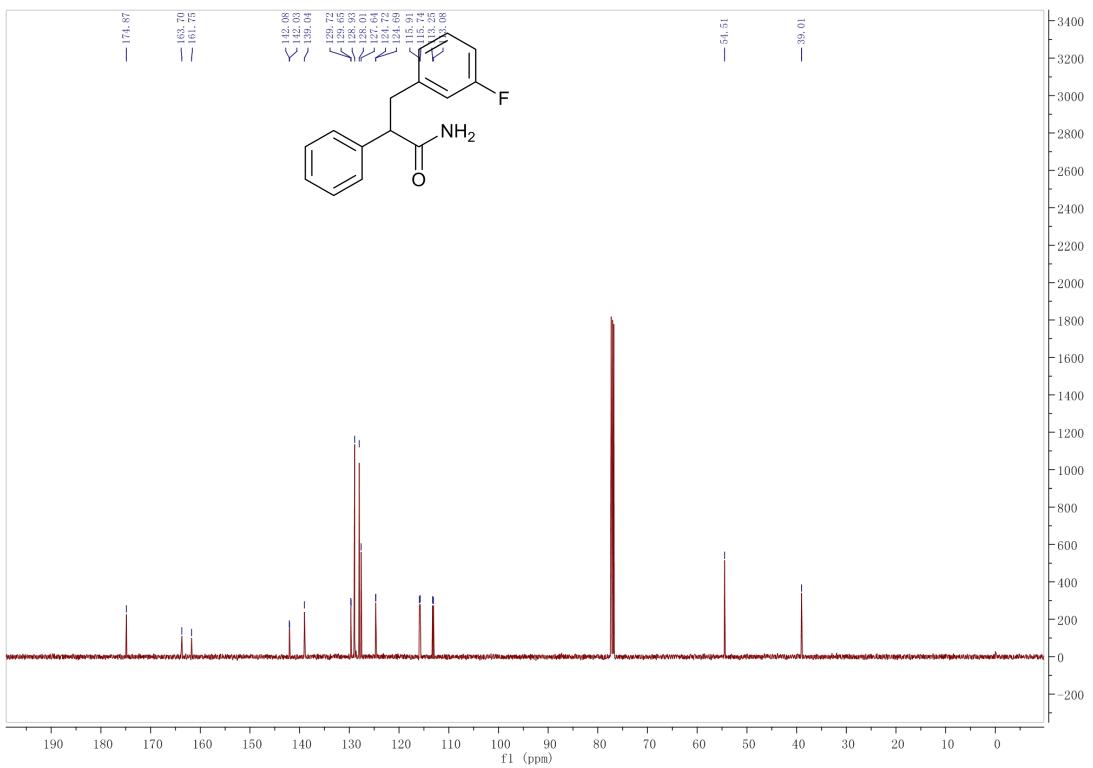
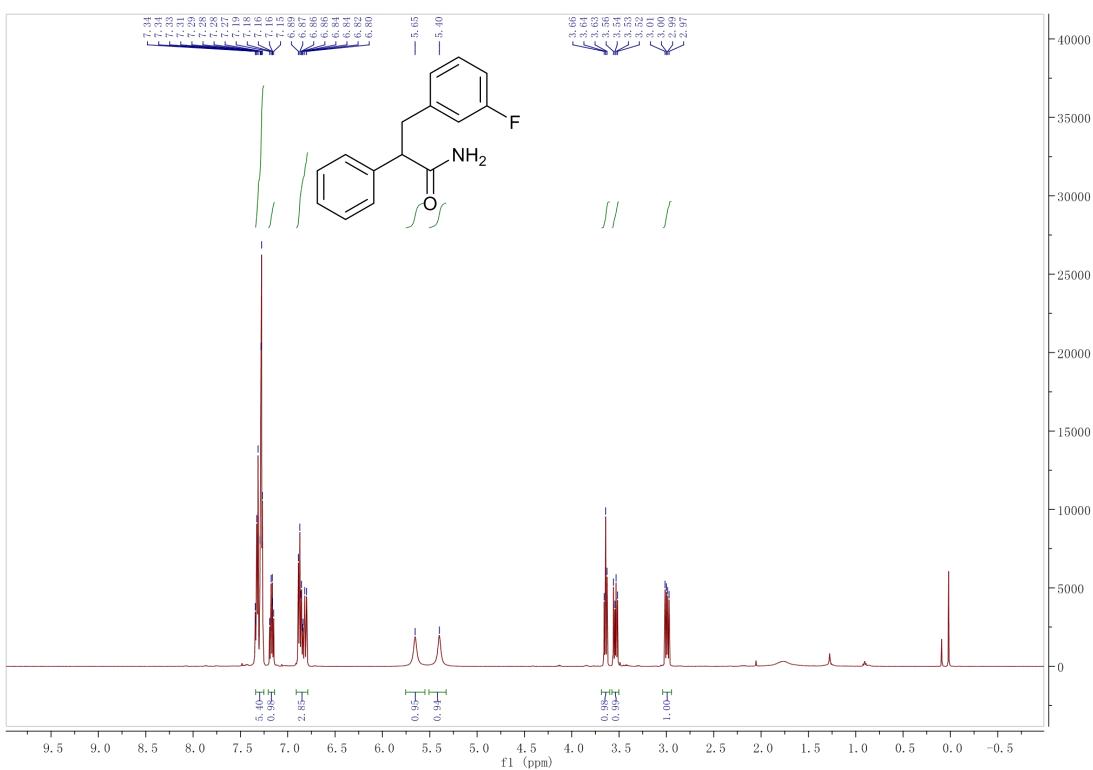
3-(3-methoxyphenyl)-2-phenylpropanamide (4ad**)**



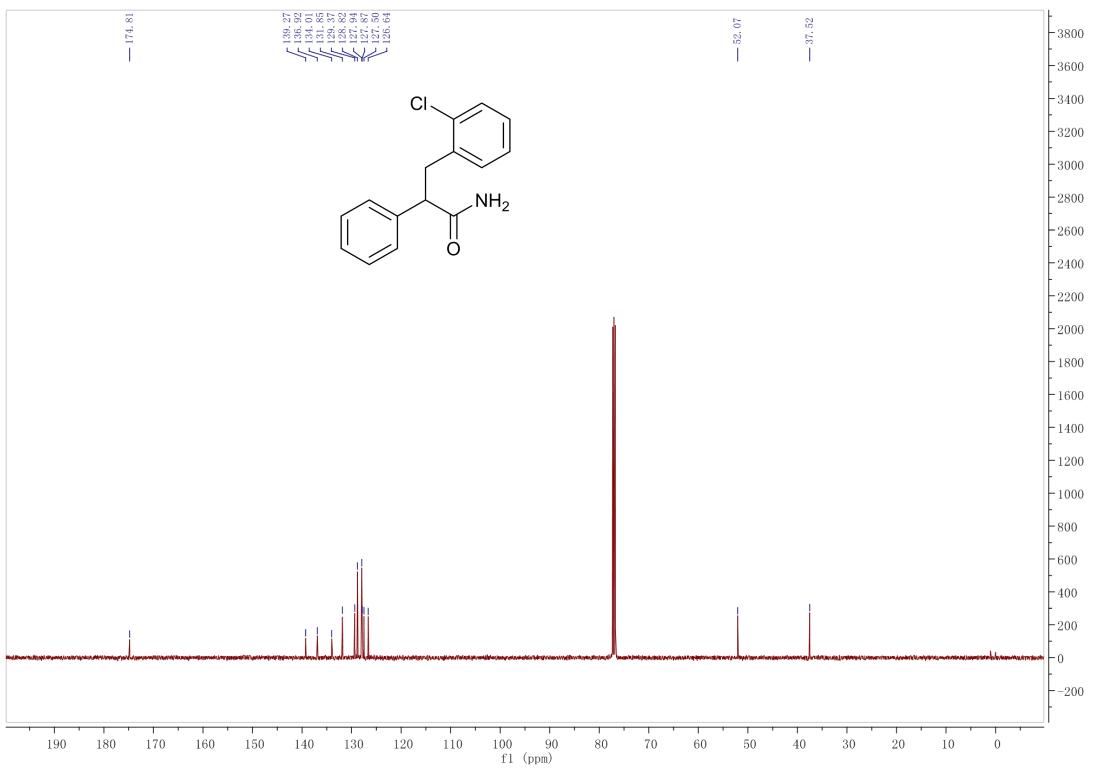
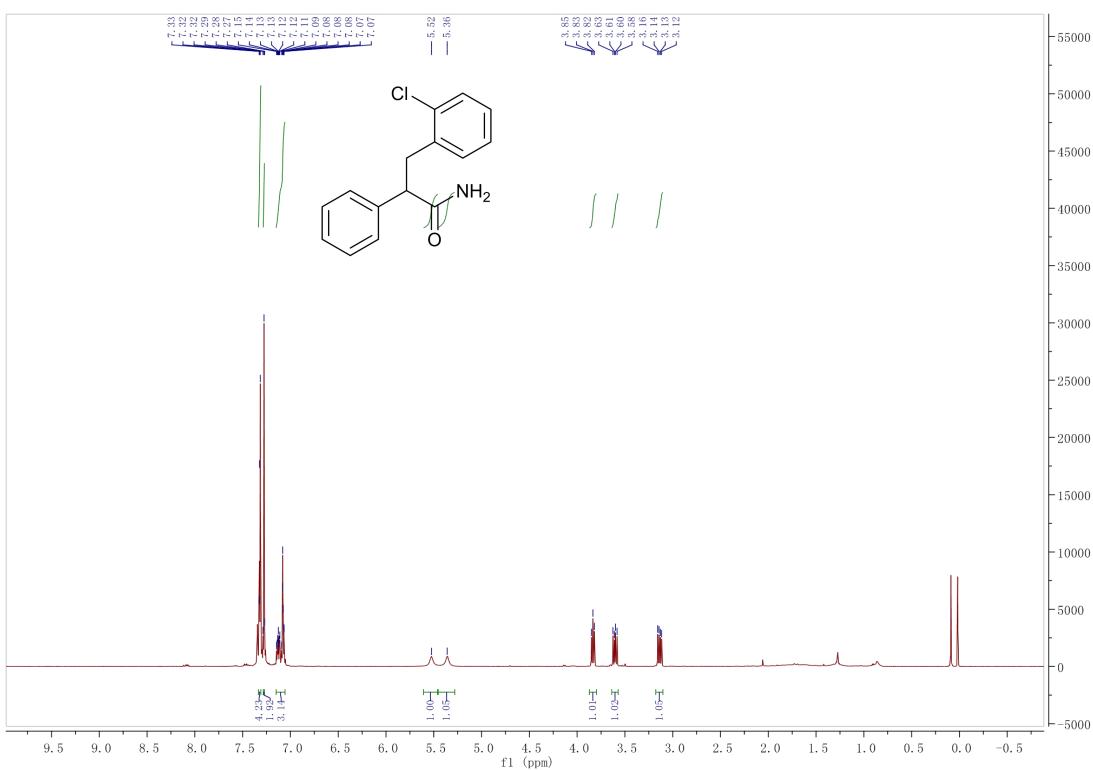
3-(4-methoxyphenyl)-2-phenylpropanamide (**4ae**)



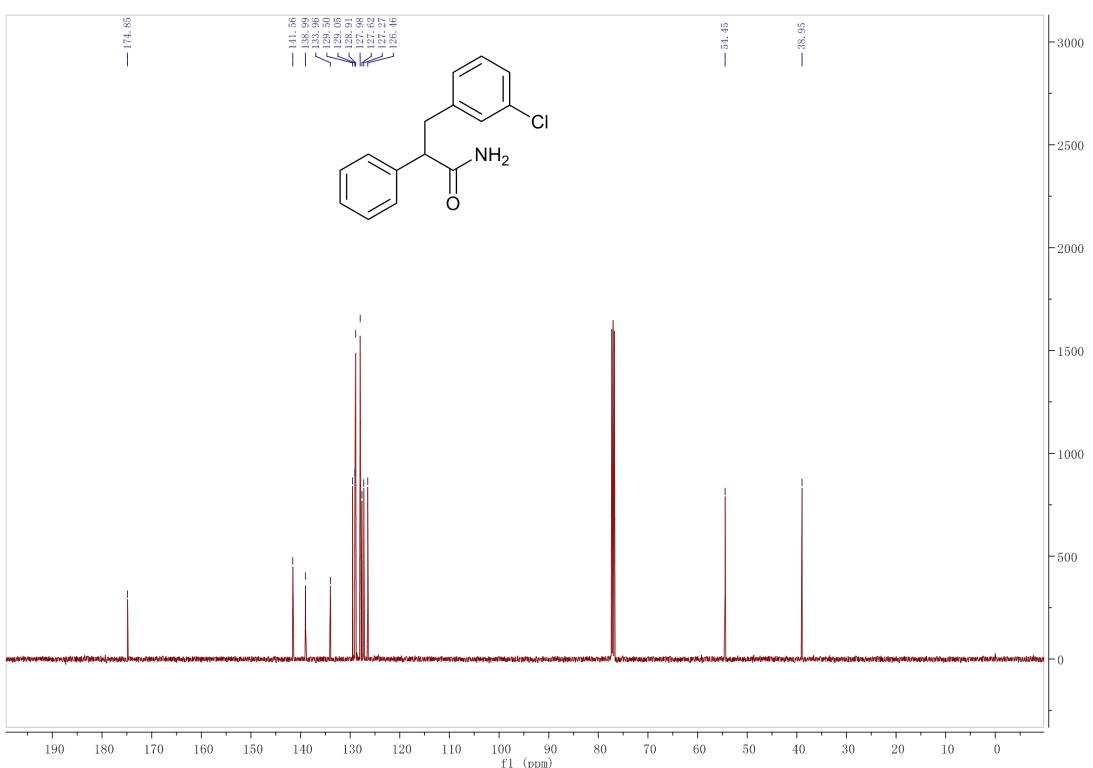
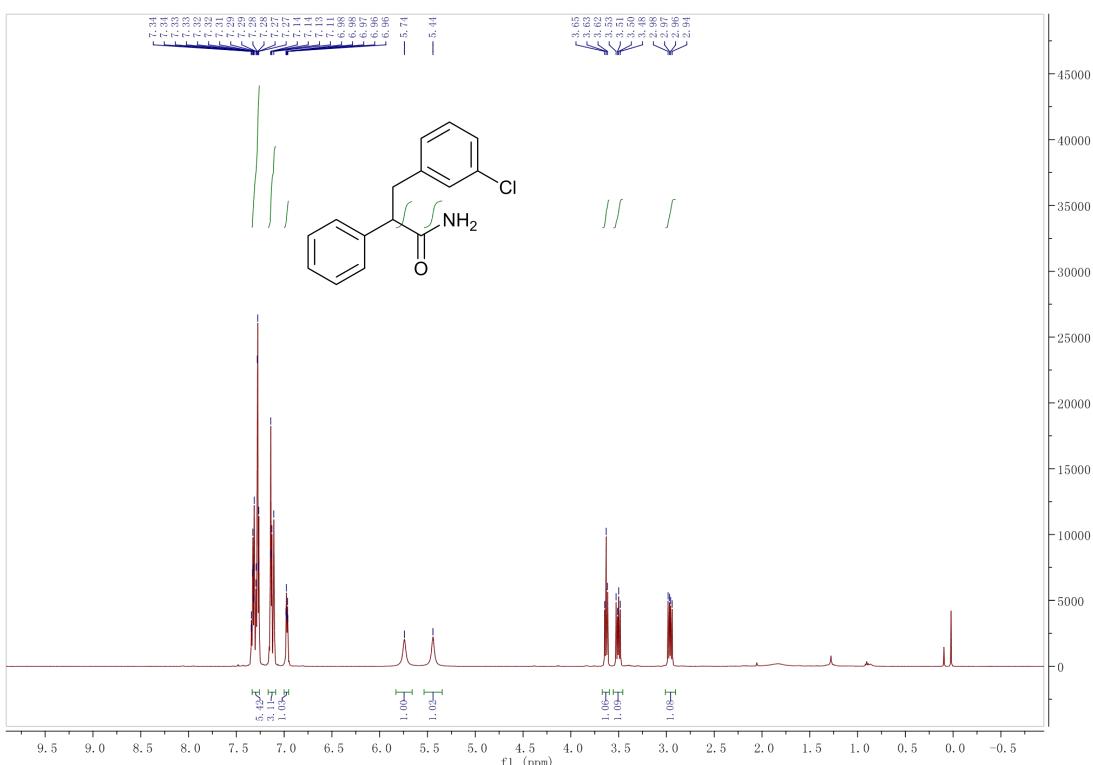
3-(3-fluorophenyl)-2-phenylpropanamide (**4af**)



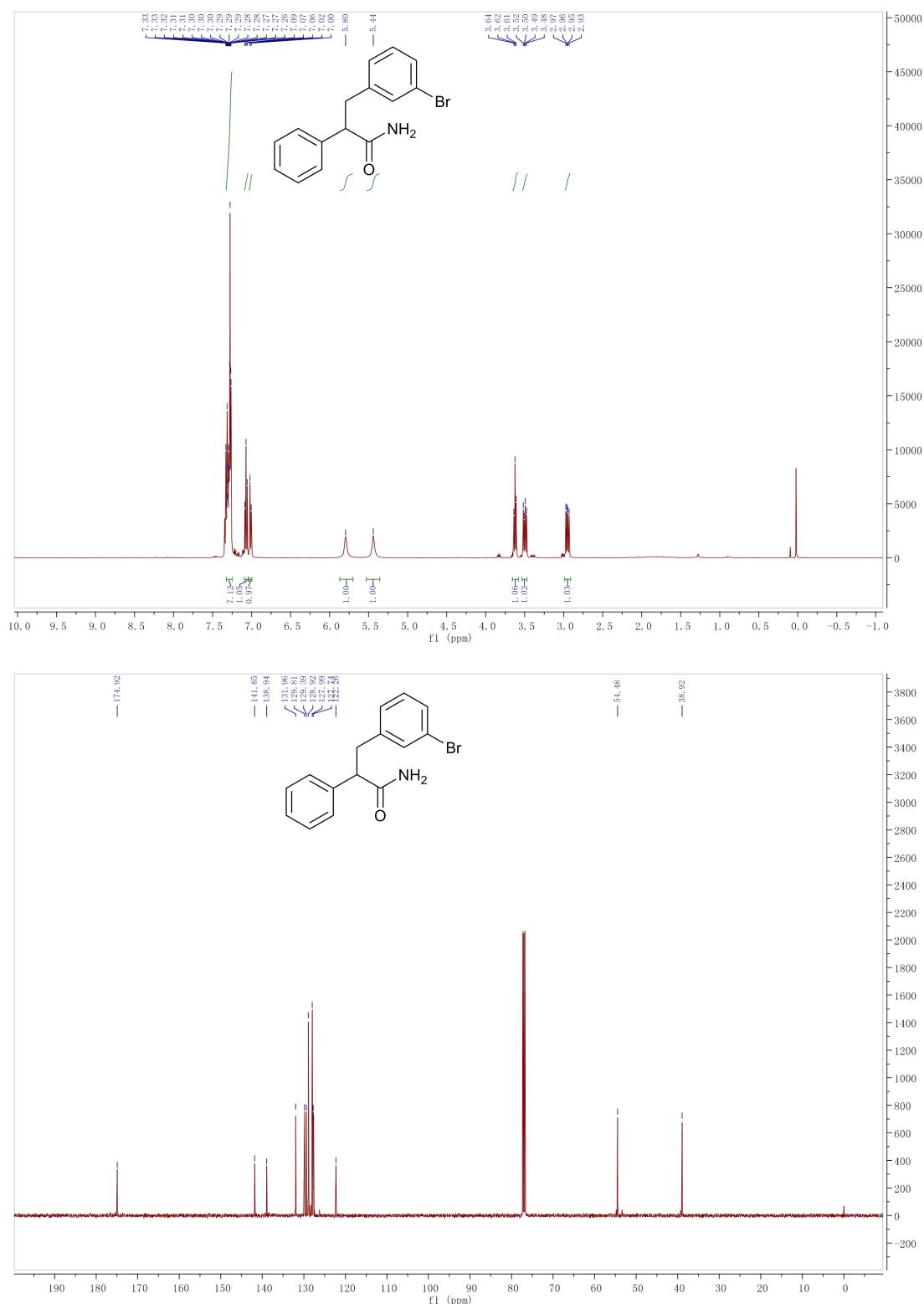
3-(2-chlorophenyl)-2-phenylpropanamide (**4ag**)



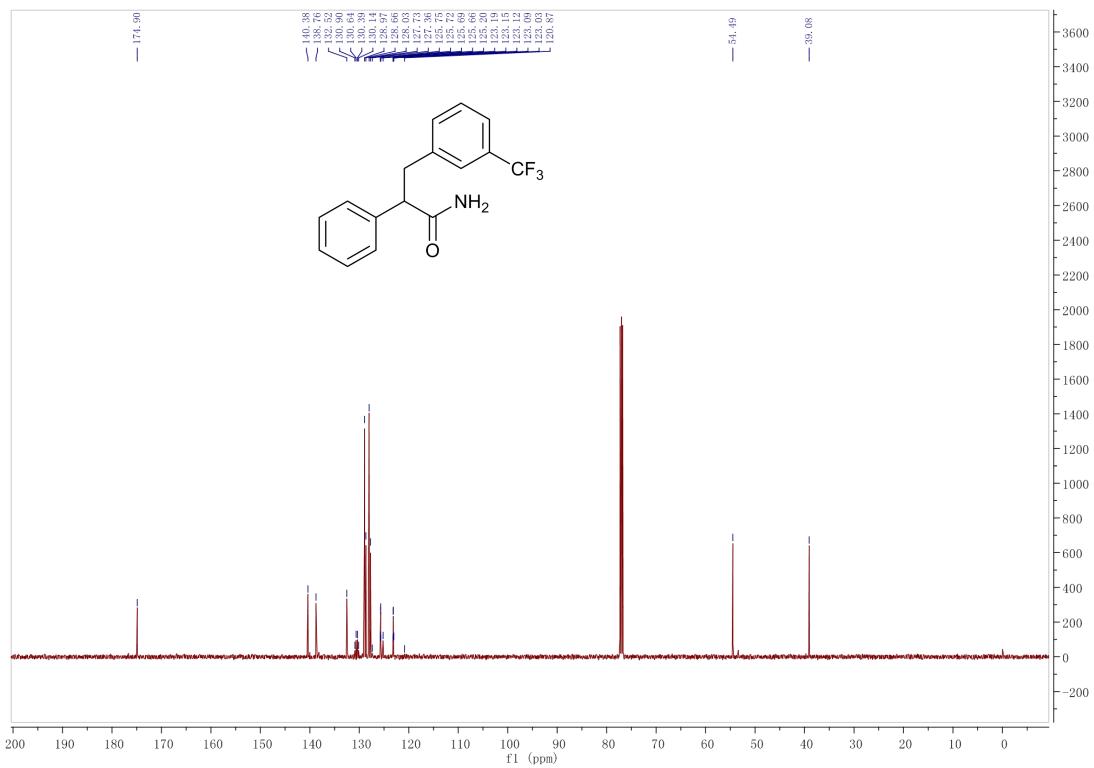
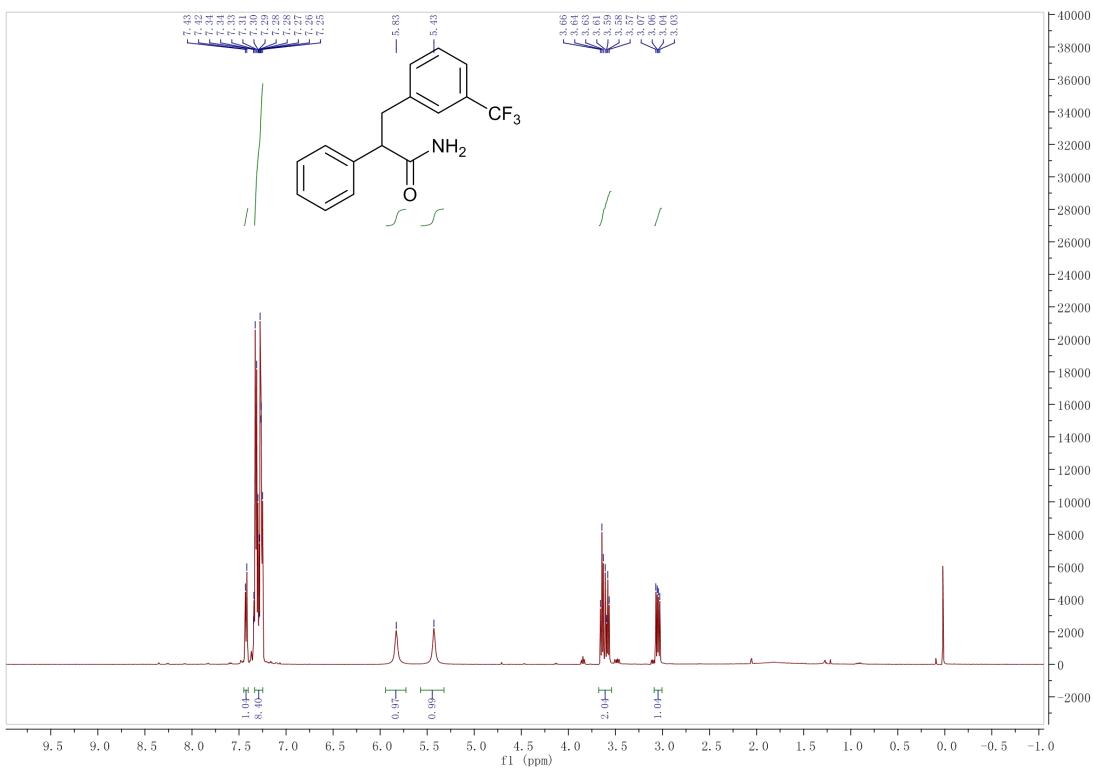
3-(3-chlorophenyl)-2-phenylpropanamide (4ah**)**



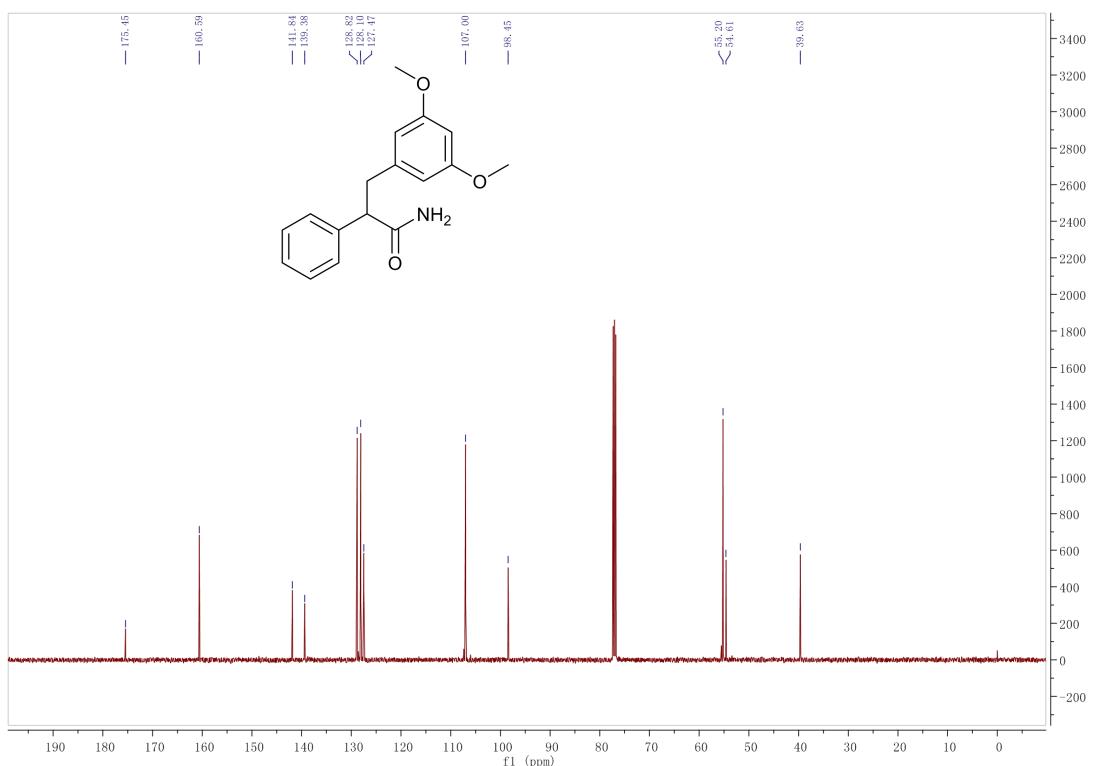
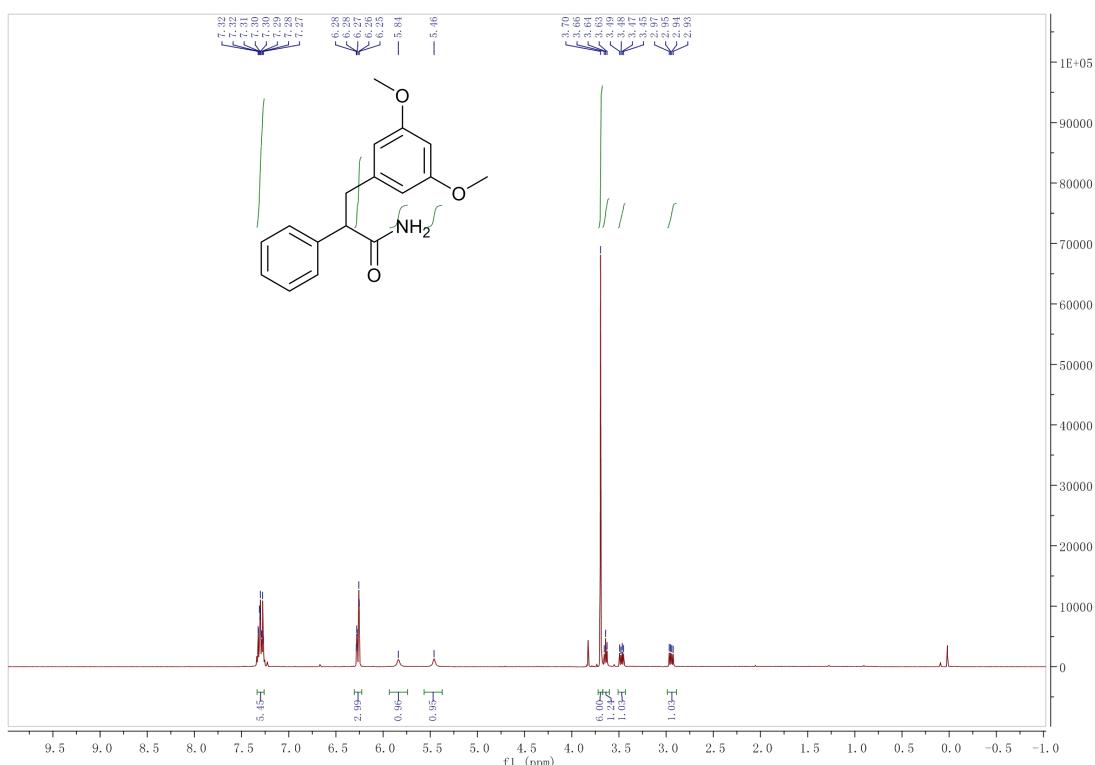
3-(3-bromophenyl)-2-phenylpropanamide (4ai**)**



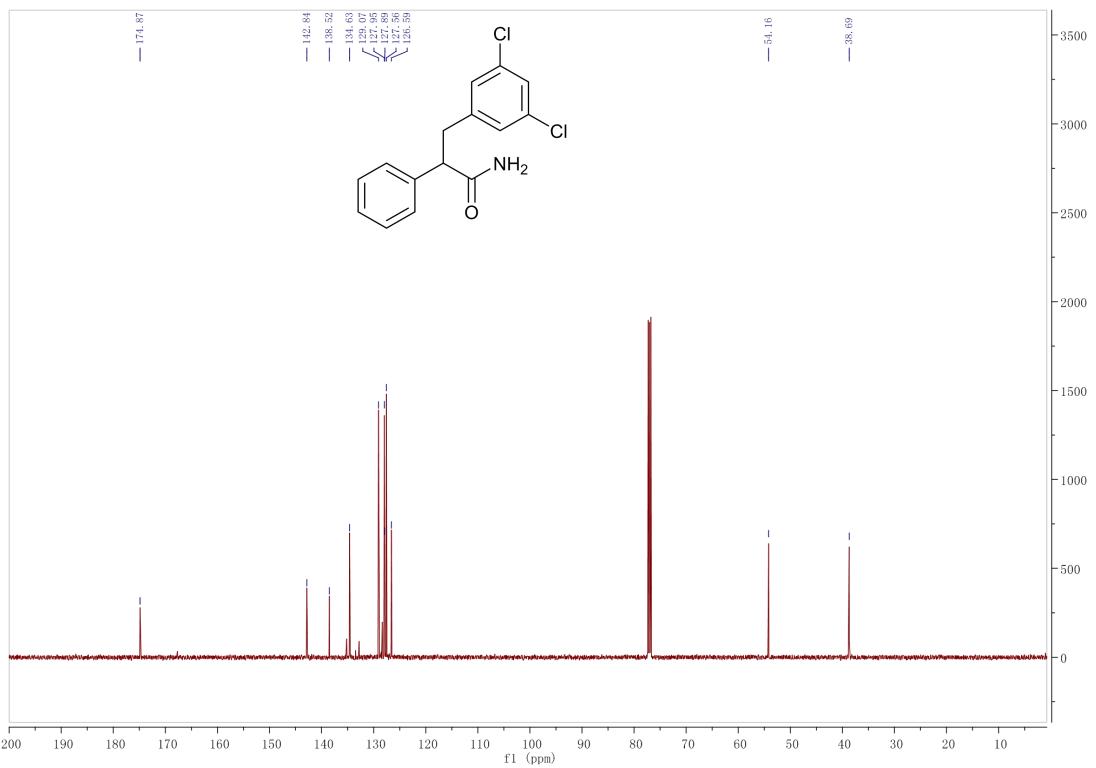
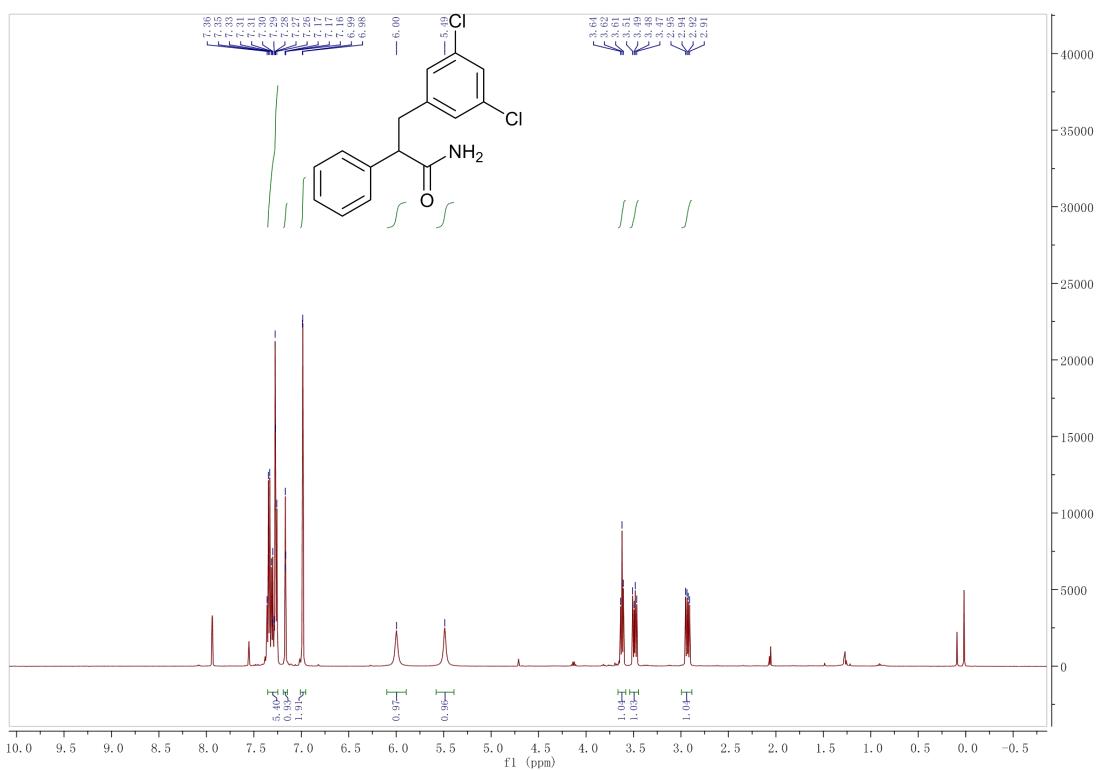
2-phenyl-3-(3-(trifluoromethyl)phenyl)propenamide (4aj**)**



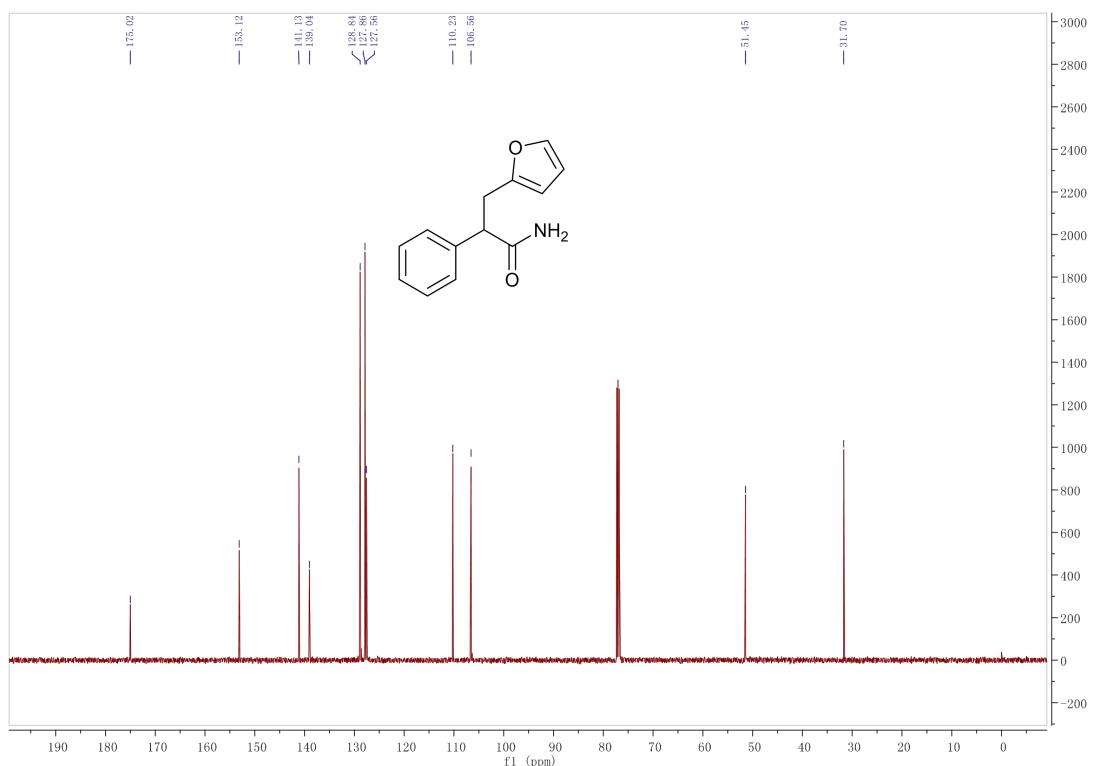
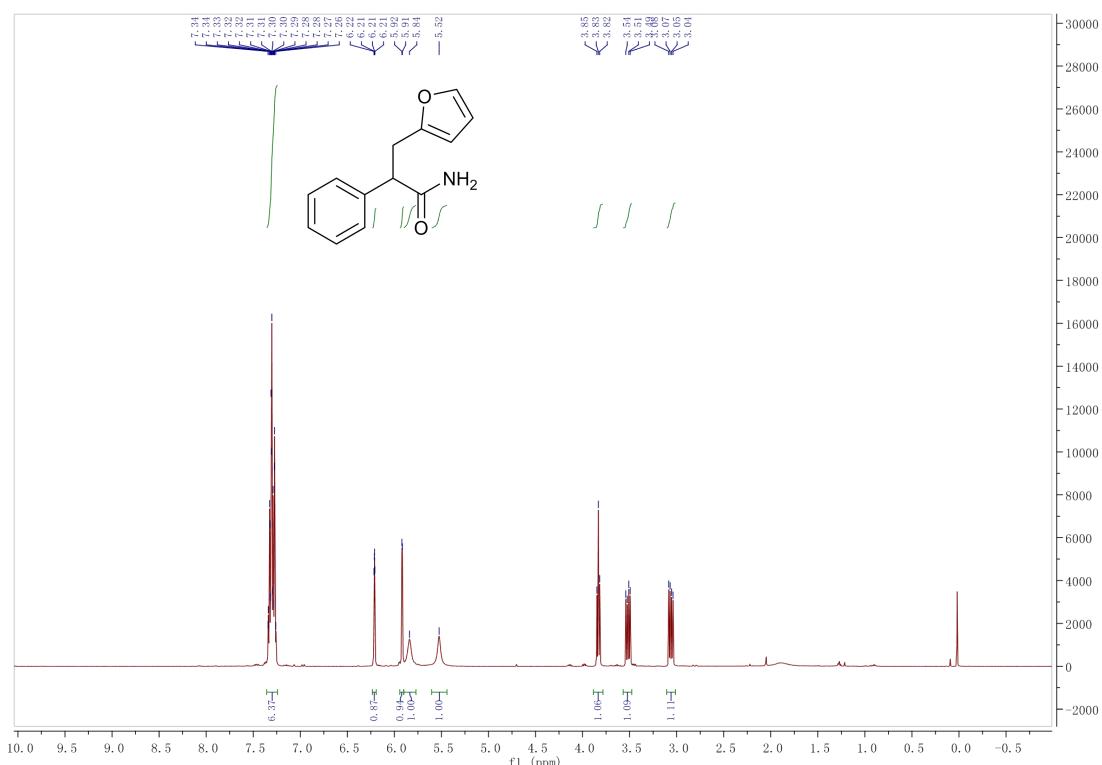
3-(3,5-dimethoxyphenyl)-2-phenylpropanamide (4ak**)**



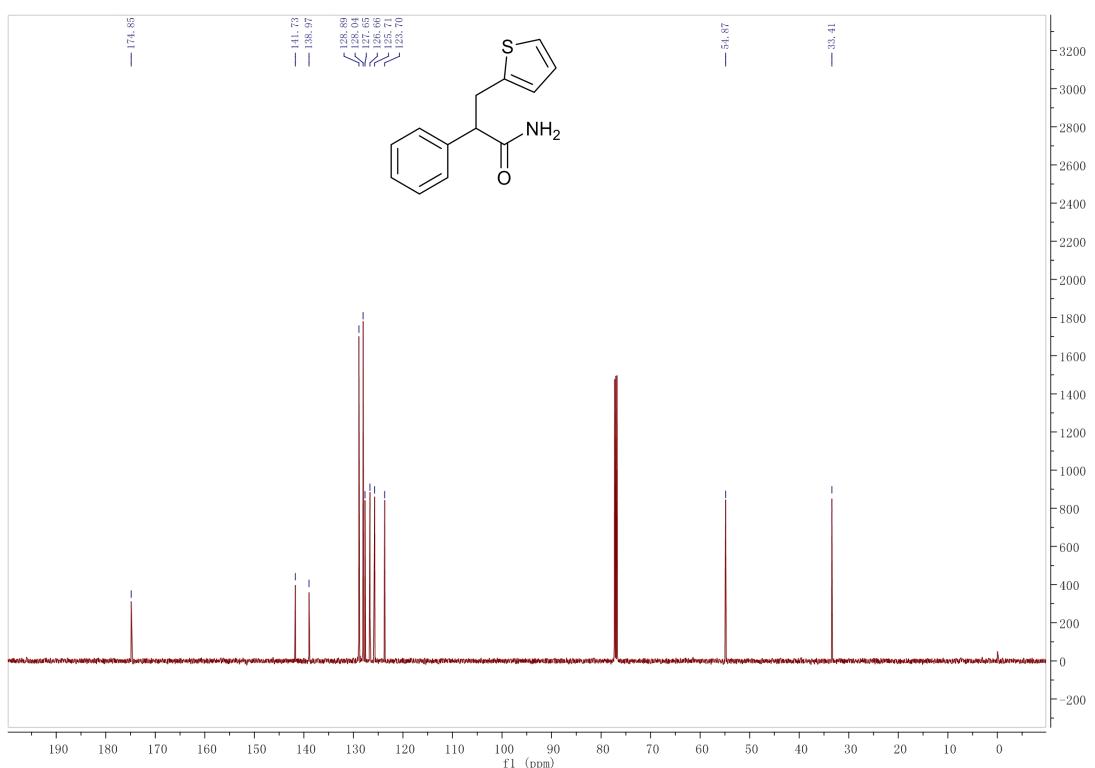
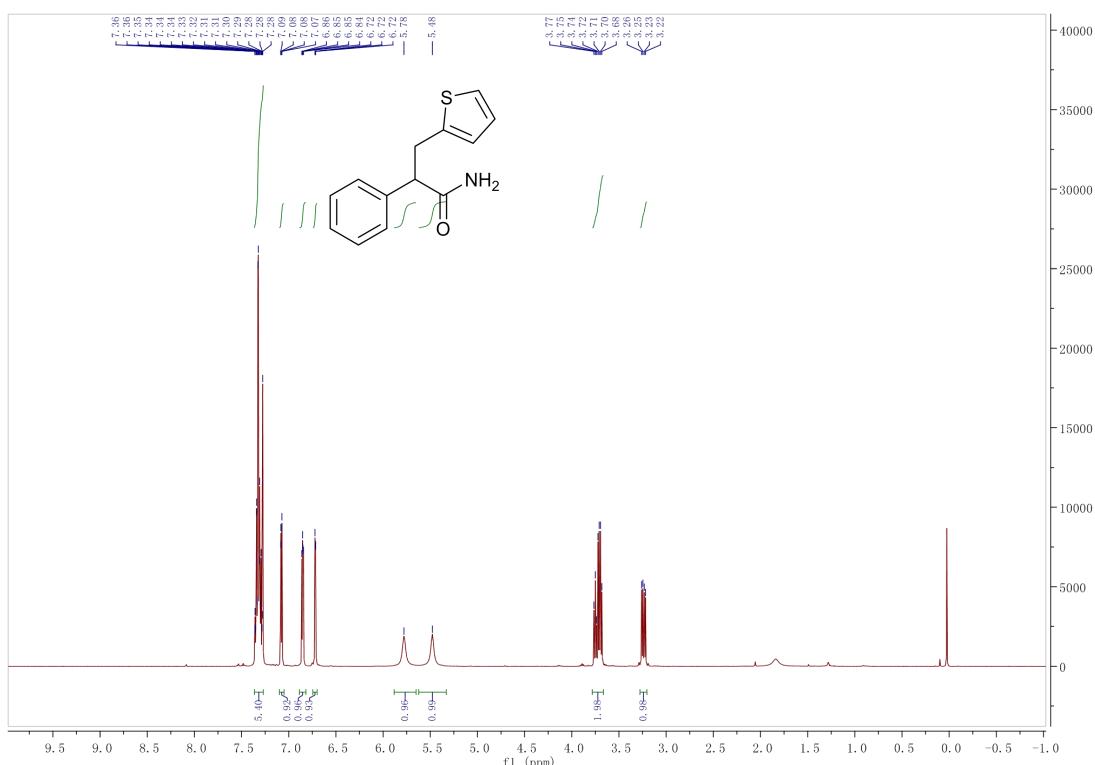
3-(3,5-dichlorophenyl)-2-phenylpropanamide (**4al**)



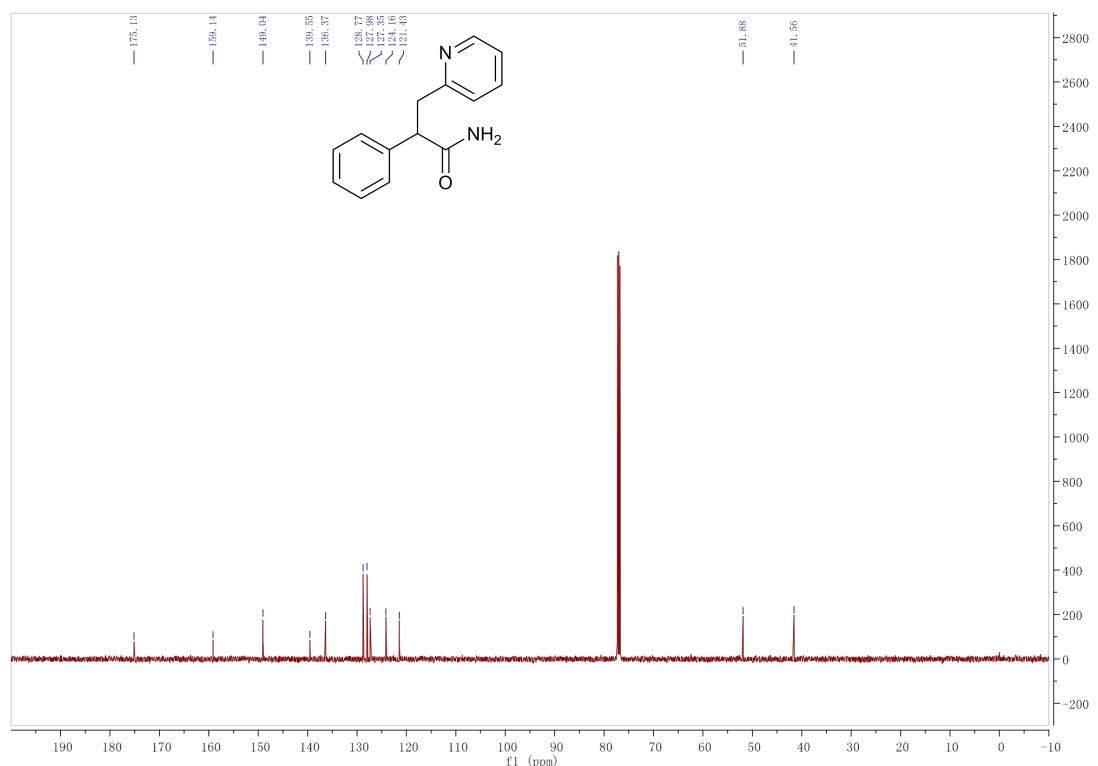
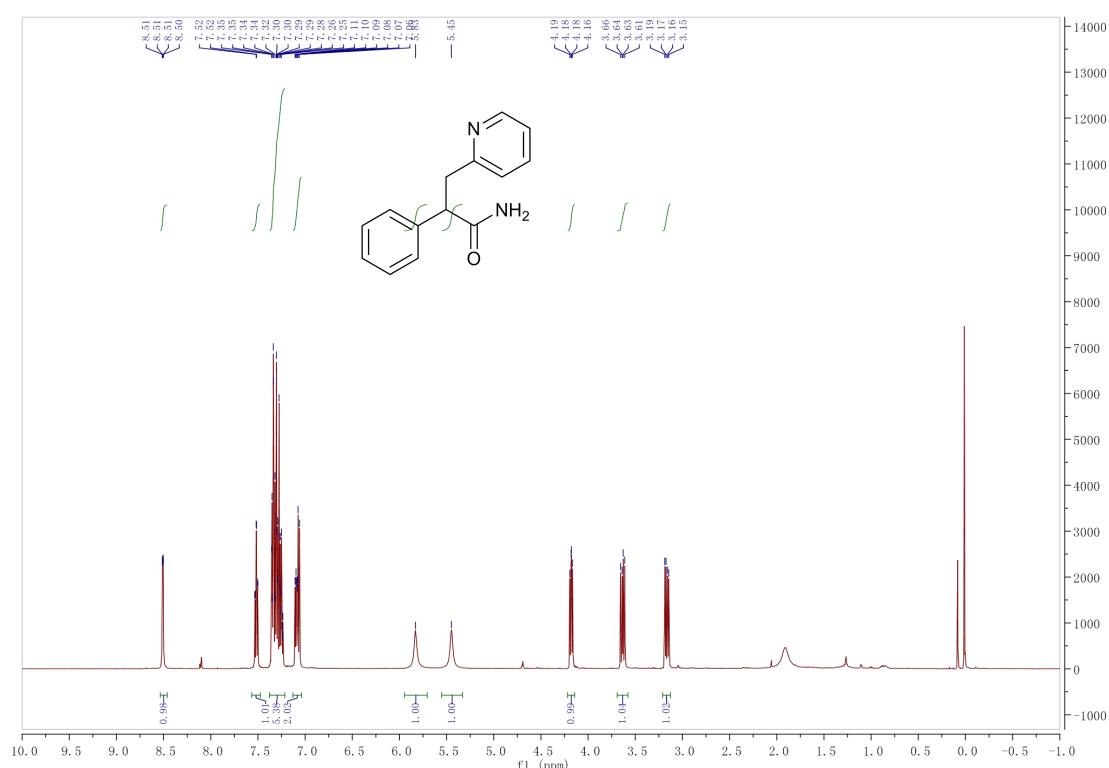
3-(furan-2-yl)-2-phenylpropanamide (4am**)**



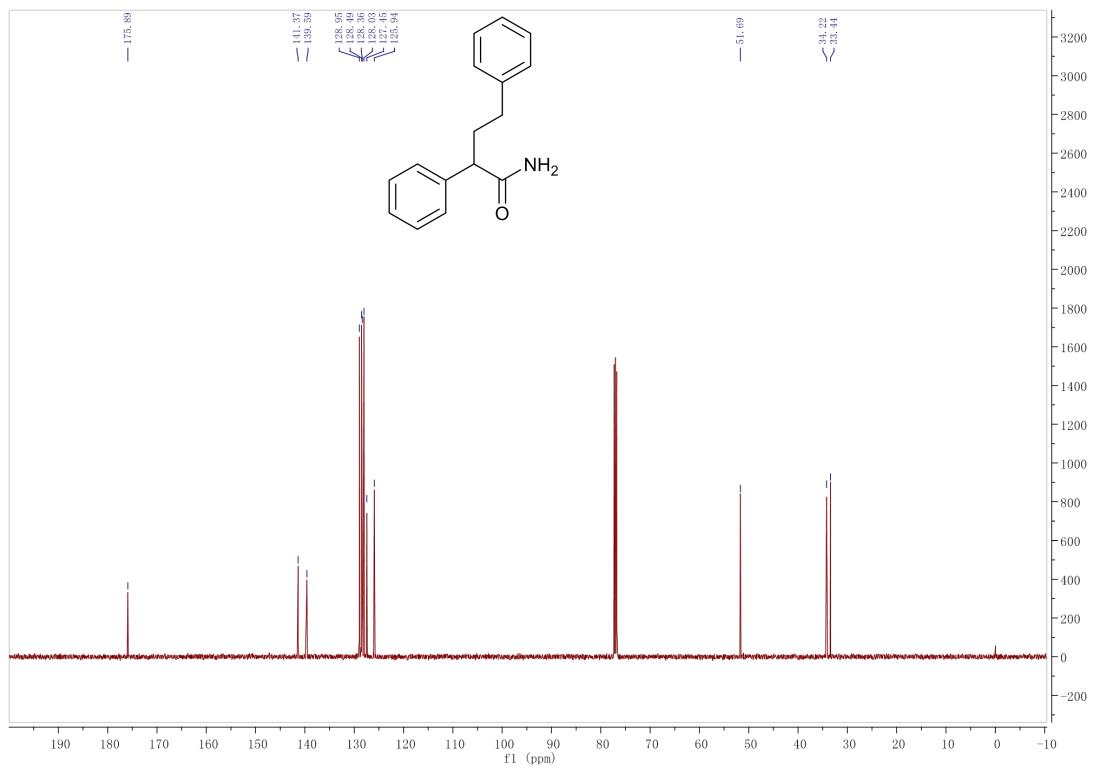
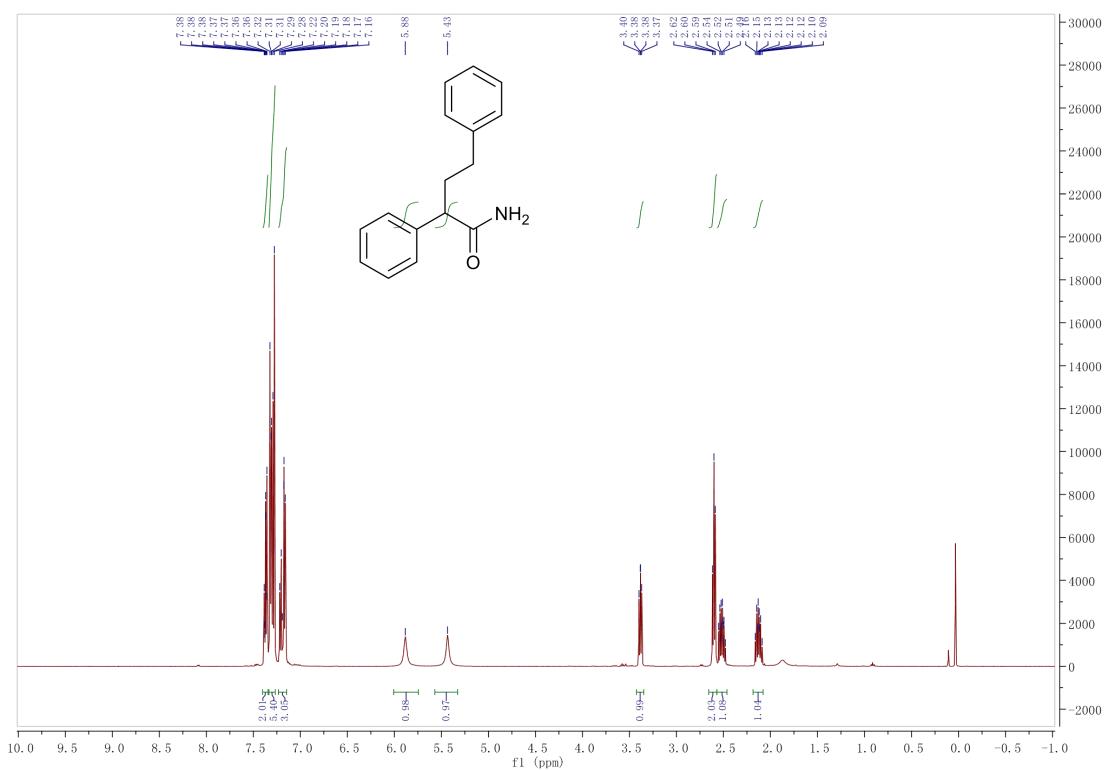
2-phenyl-3-(thiophen-2-yl)propenamide (**4an**)



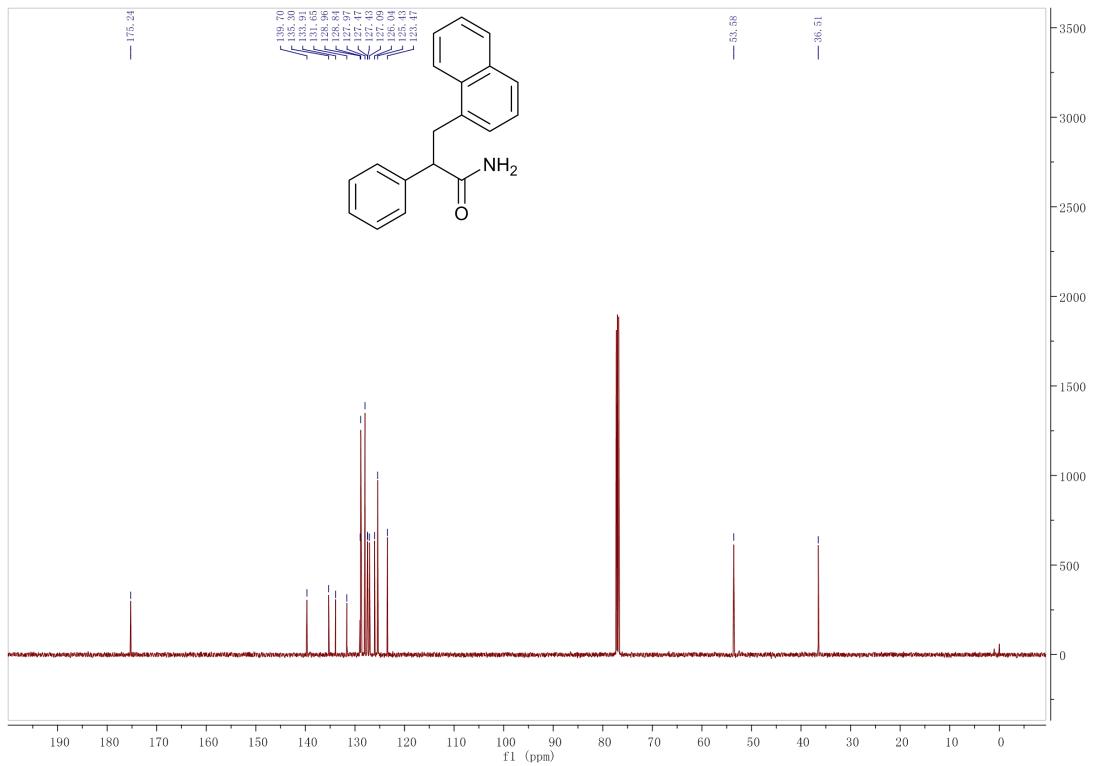
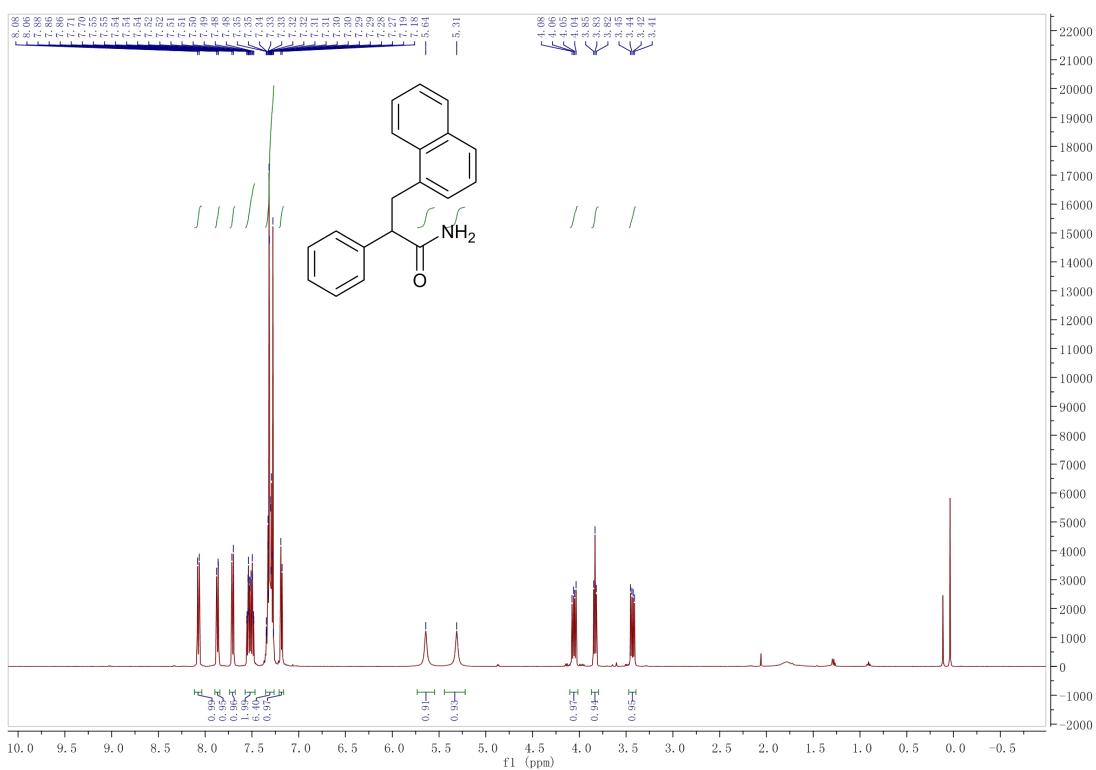
2-phenyl-3-(pyridin-2-yl)propenamide (4ao**)**



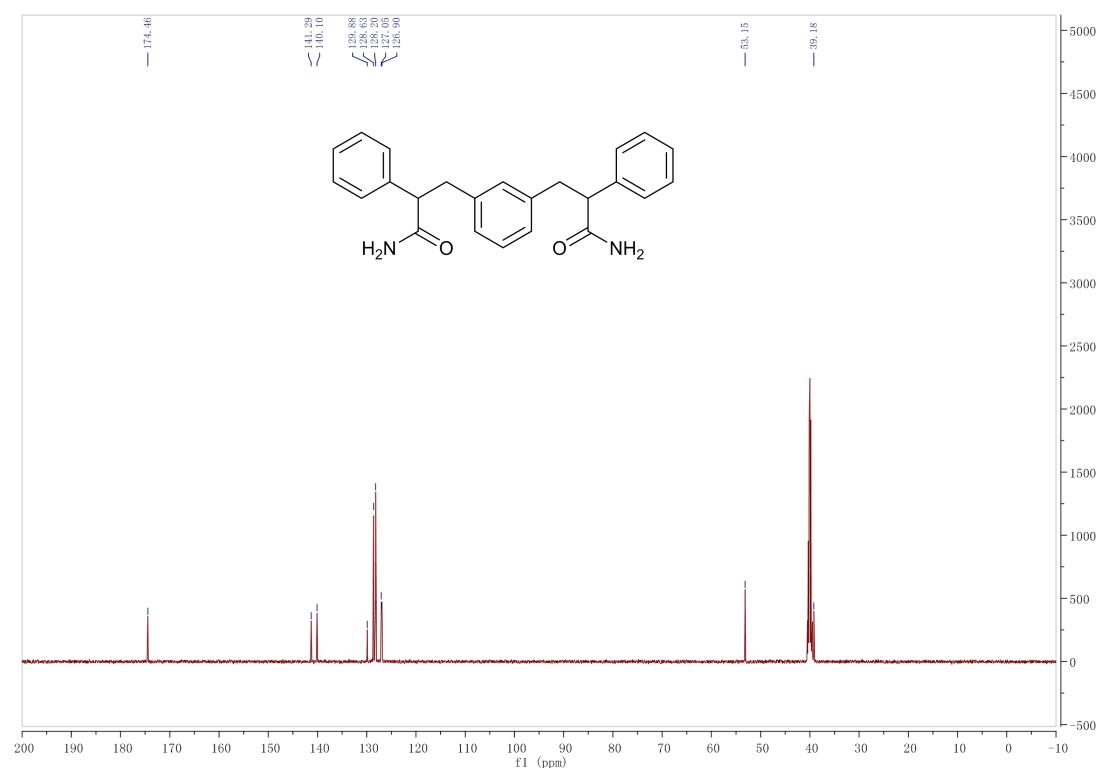
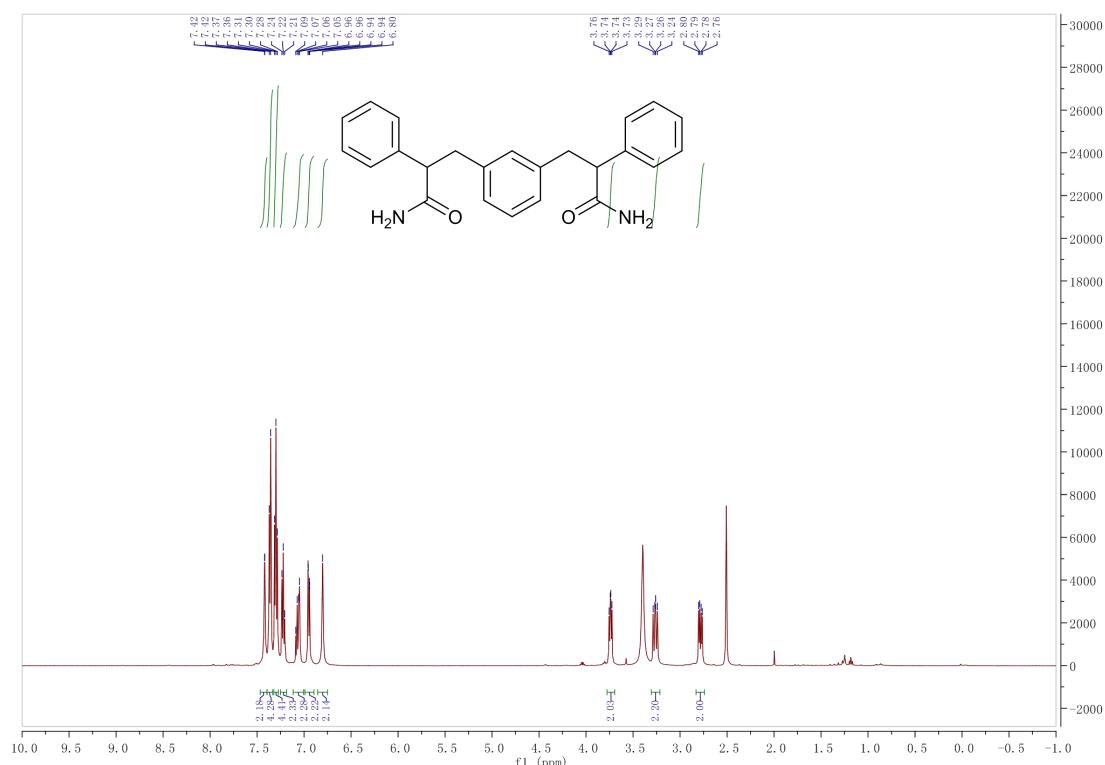
2,4-diphenylbutanamide (4ap)



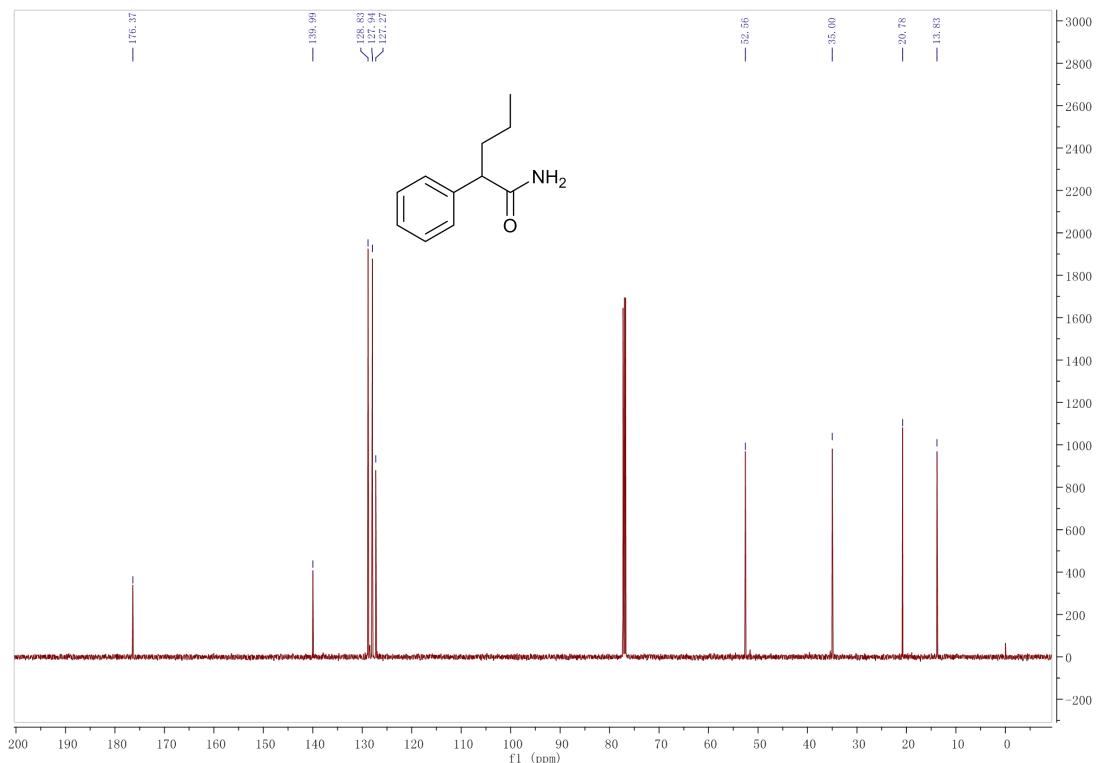
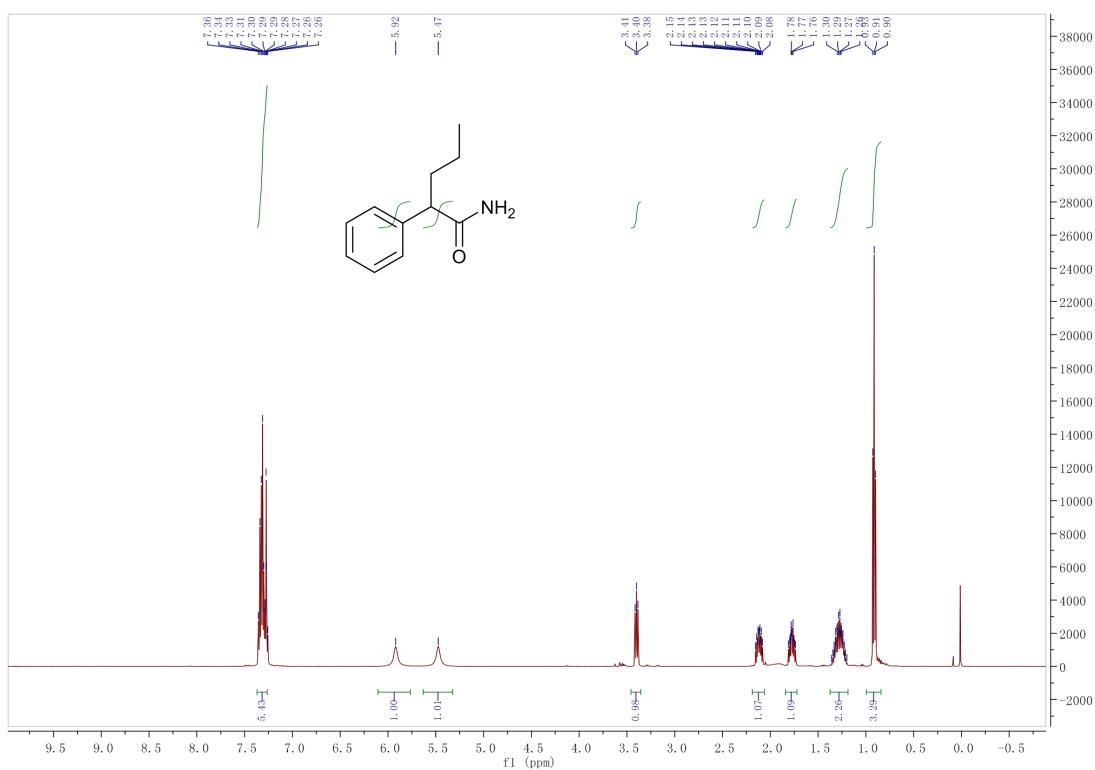
3-(naphthalen-1-yl)-2-phenylpropanamide (4aq**)**



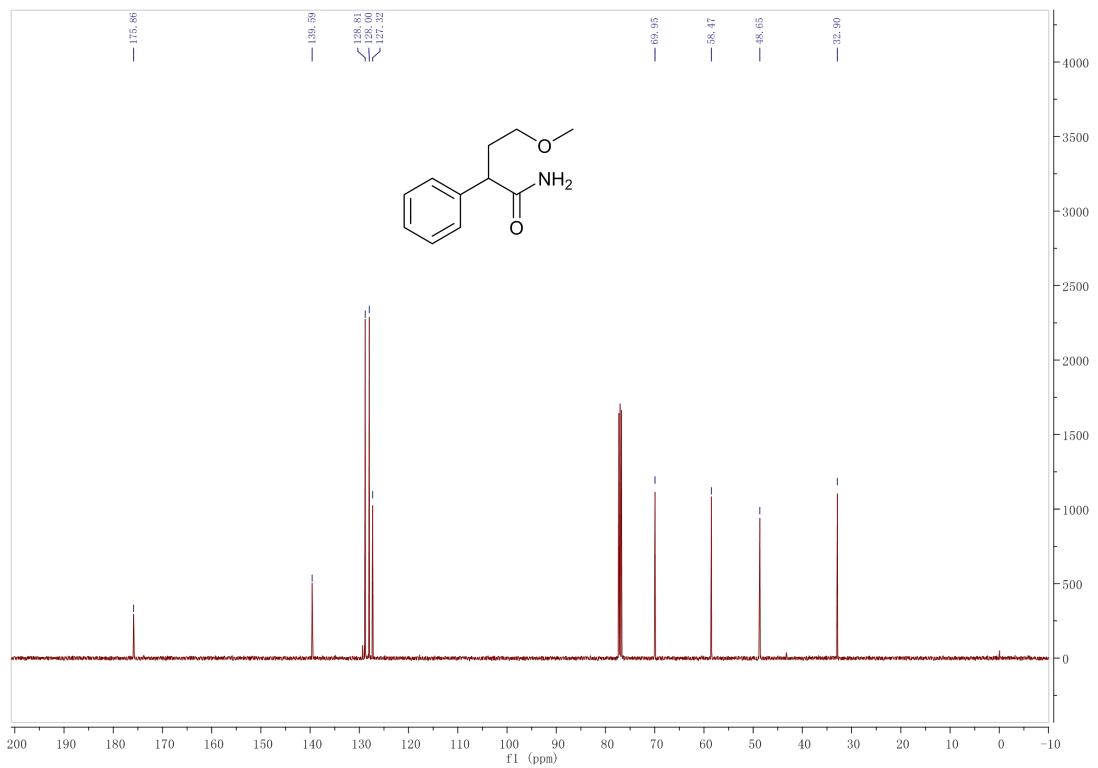
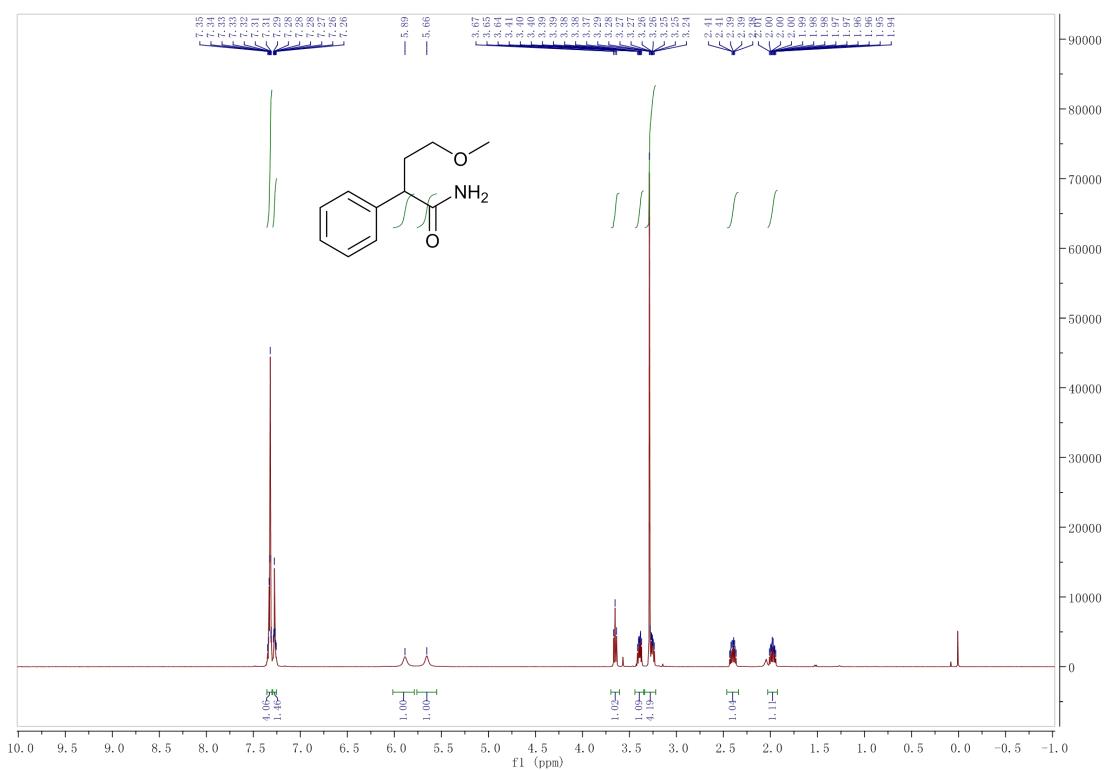
3,3'-(1,3-phenylene)bis(2-phenylpropanamide) (4ar**)**



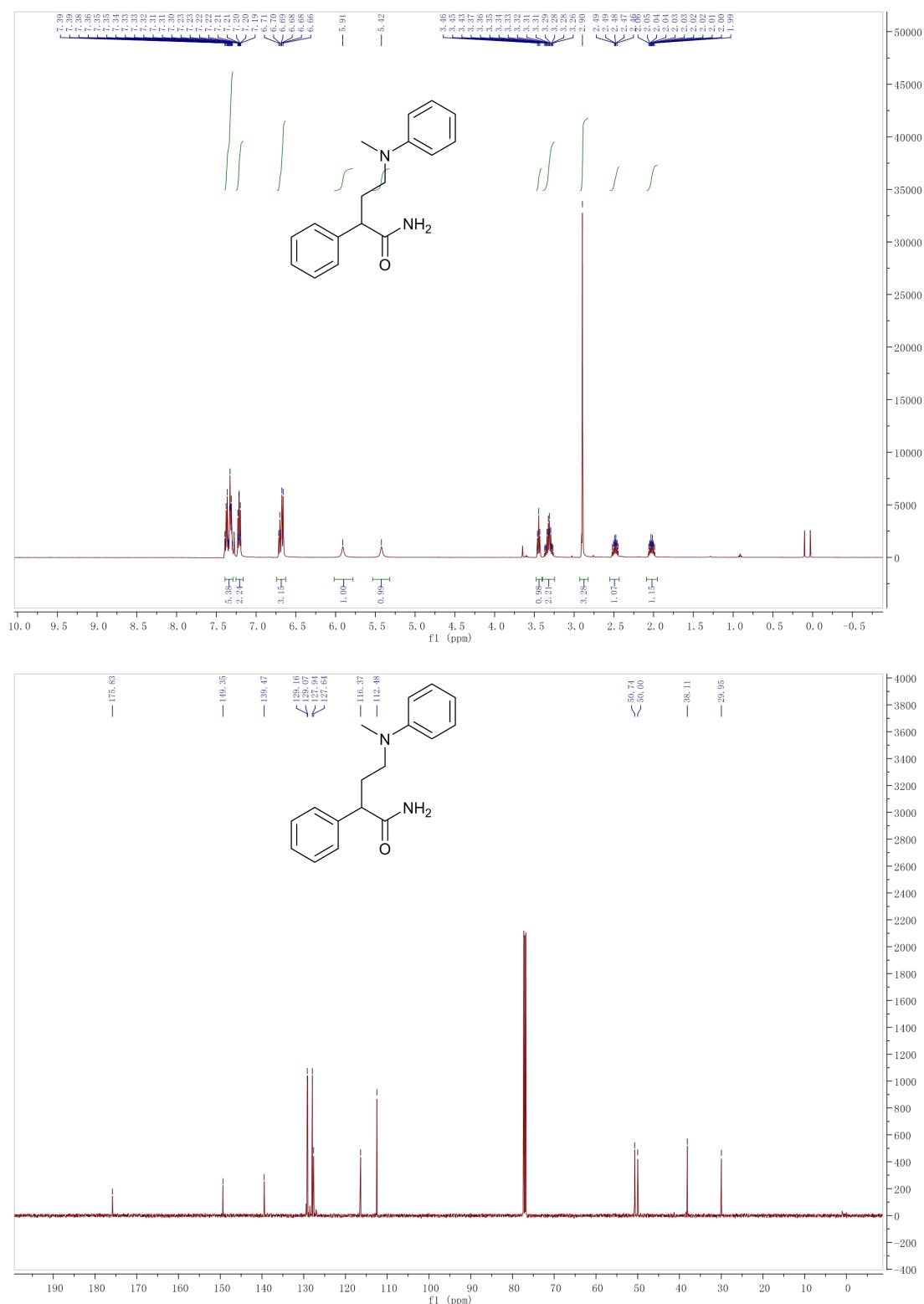
2-phenylpentanamide (4as**)**



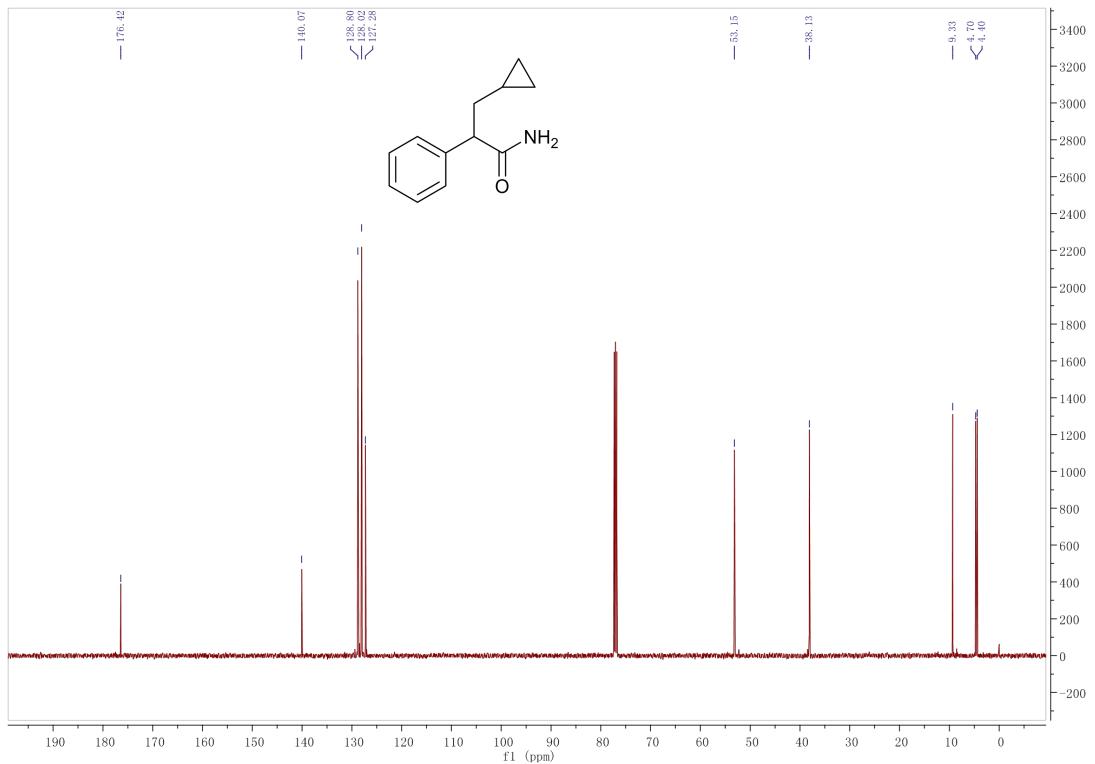
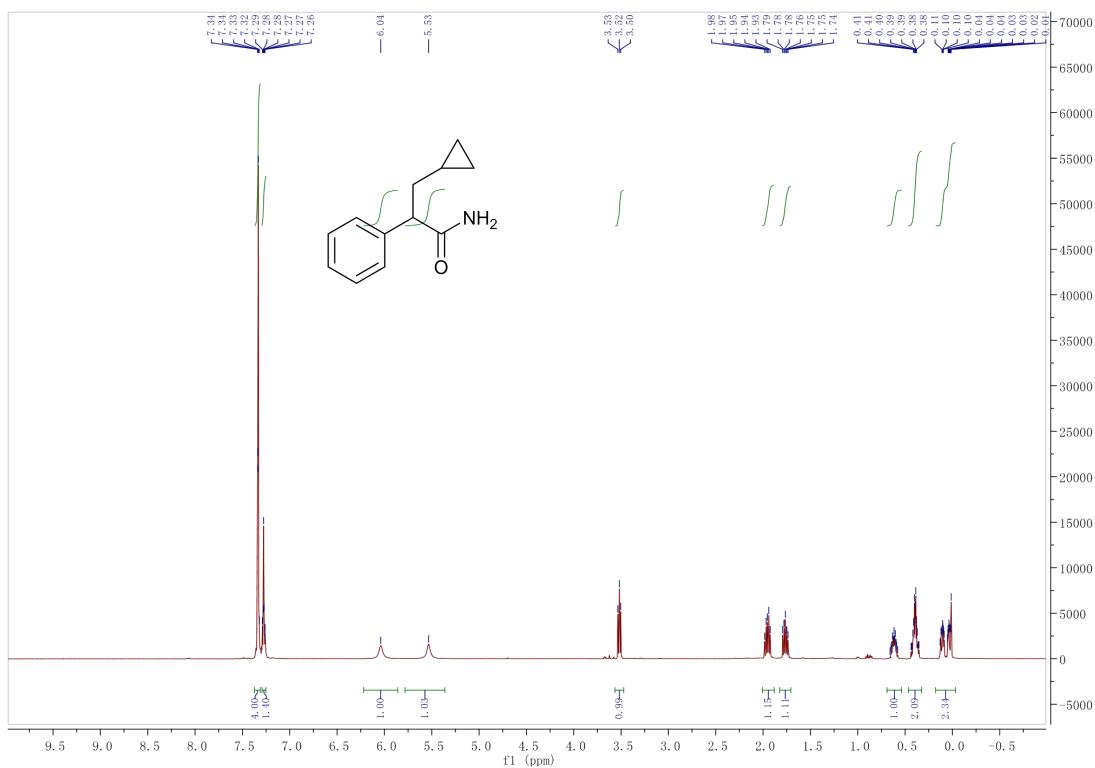
4-methoxy-2-phenylbutanamide (**4at**)



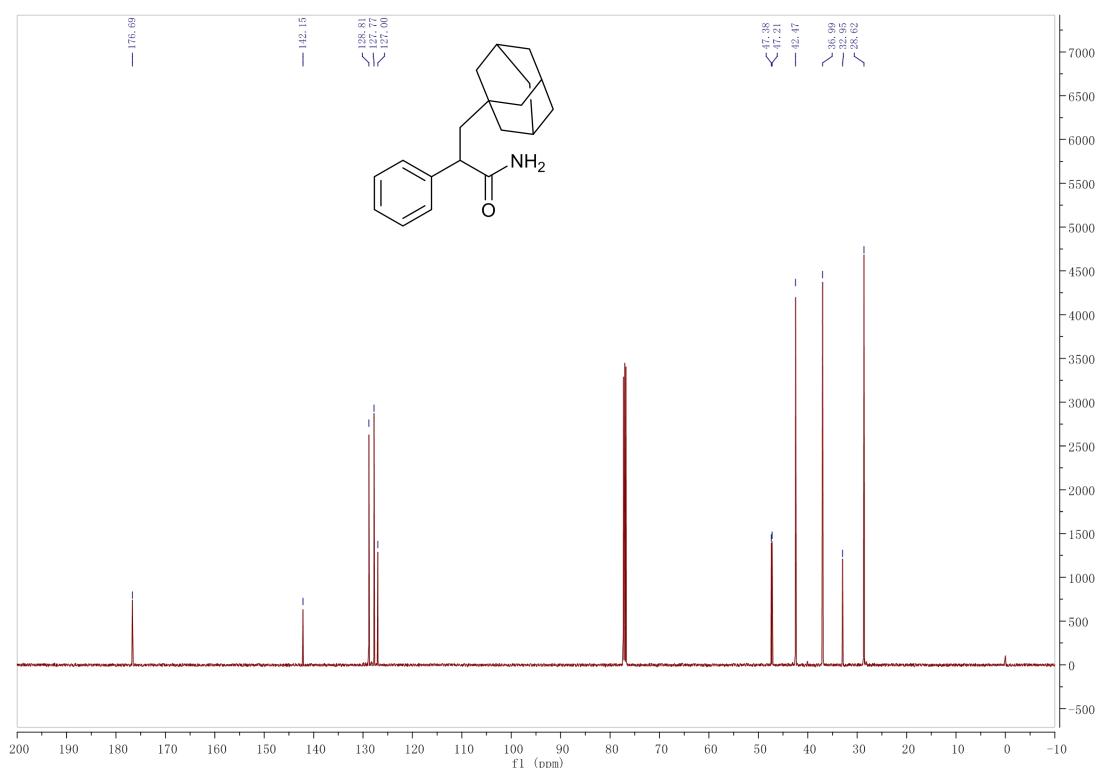
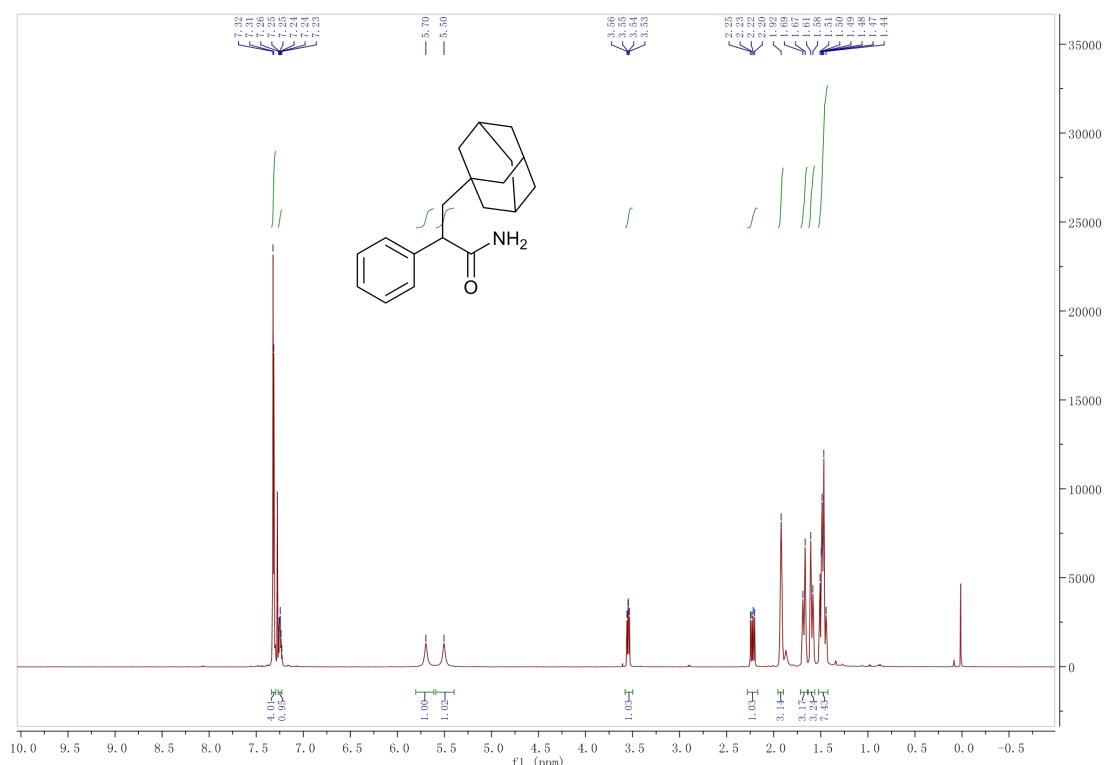
4-(methyl(phenyl)amino)-2-phenylbutanamide (4au**)**



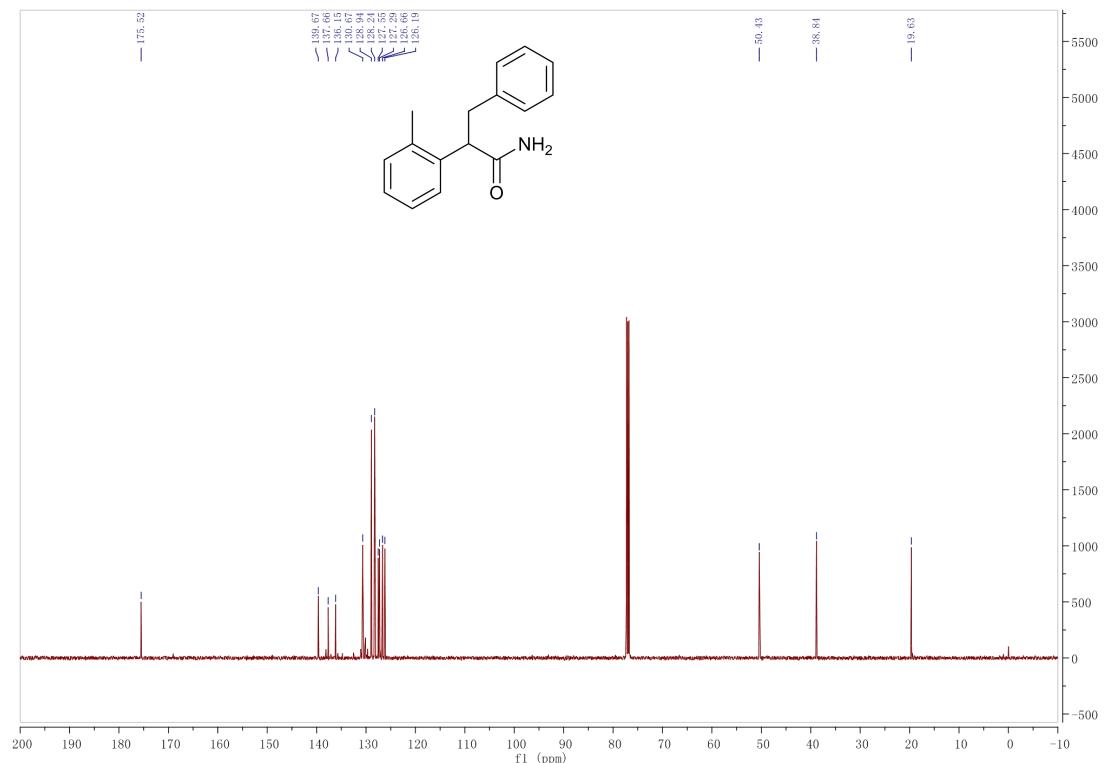
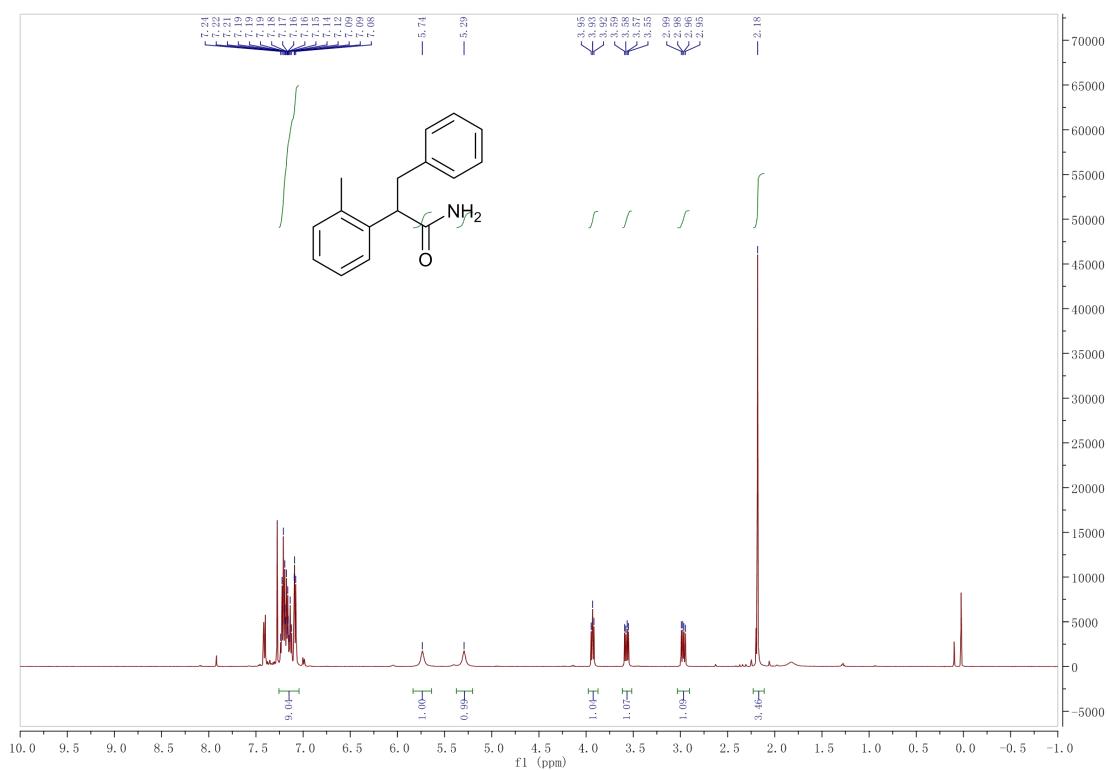
3-cyclopropyl-2-phenylpropanamide (4av**)**



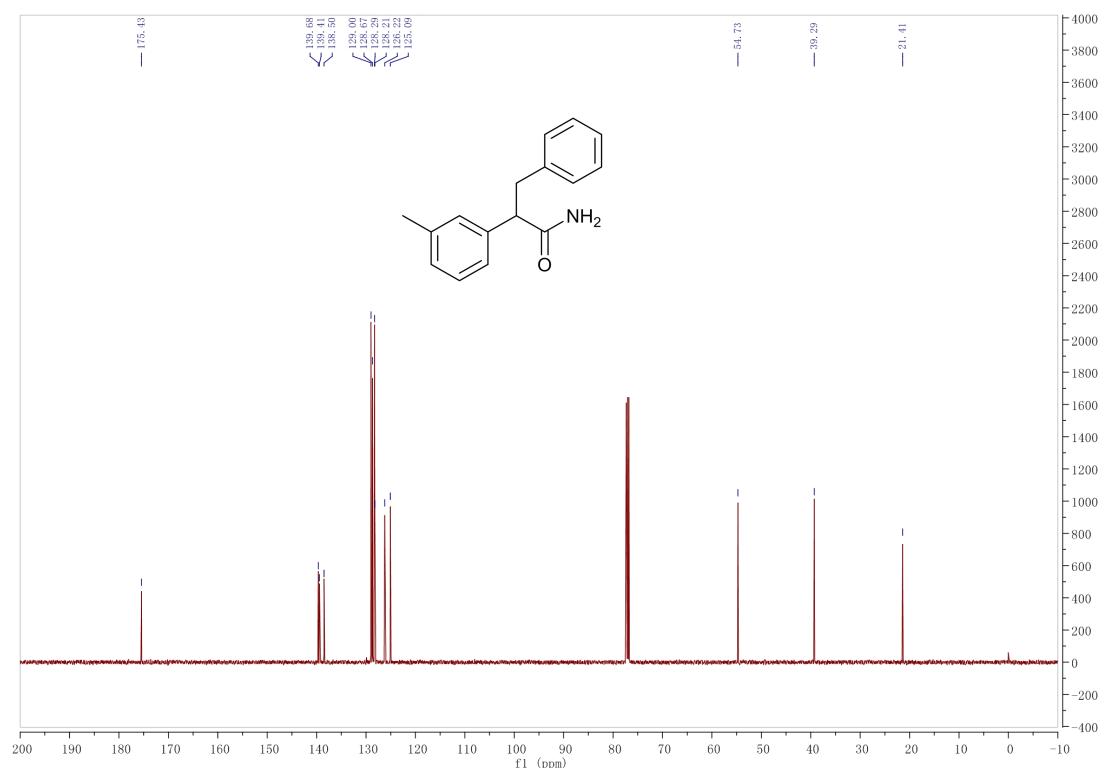
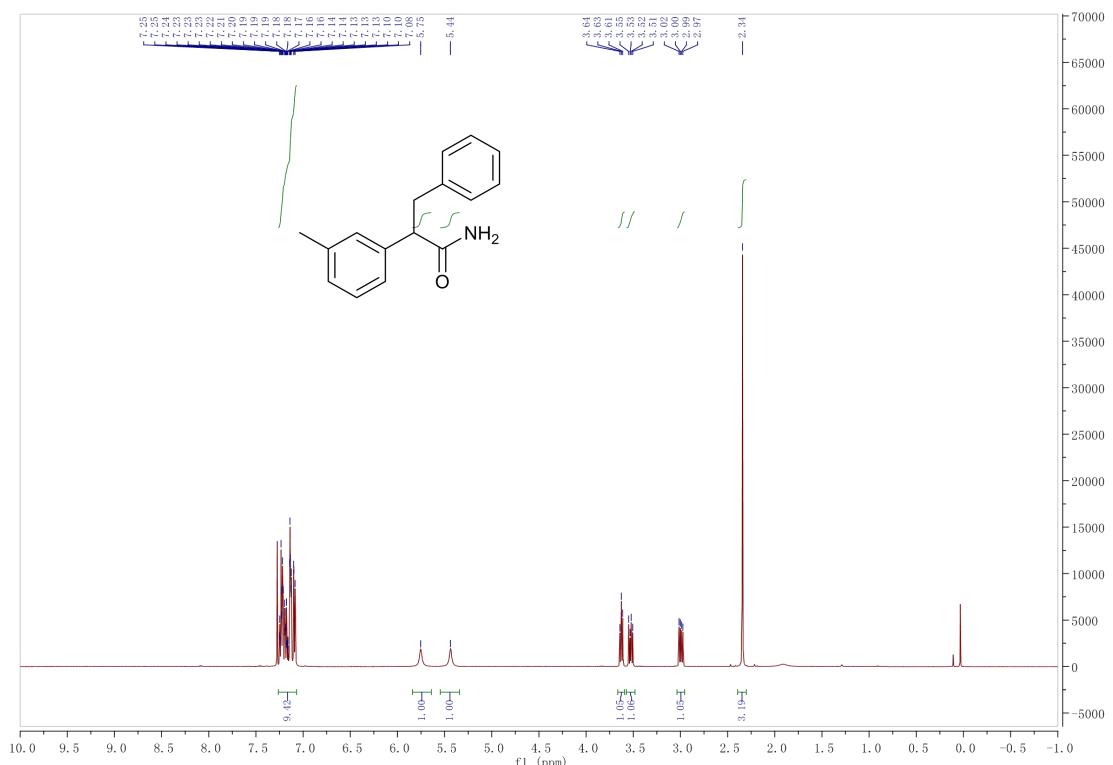
3-(adamantan-1-yl)-2-phenylpropanamide (4aw**)**



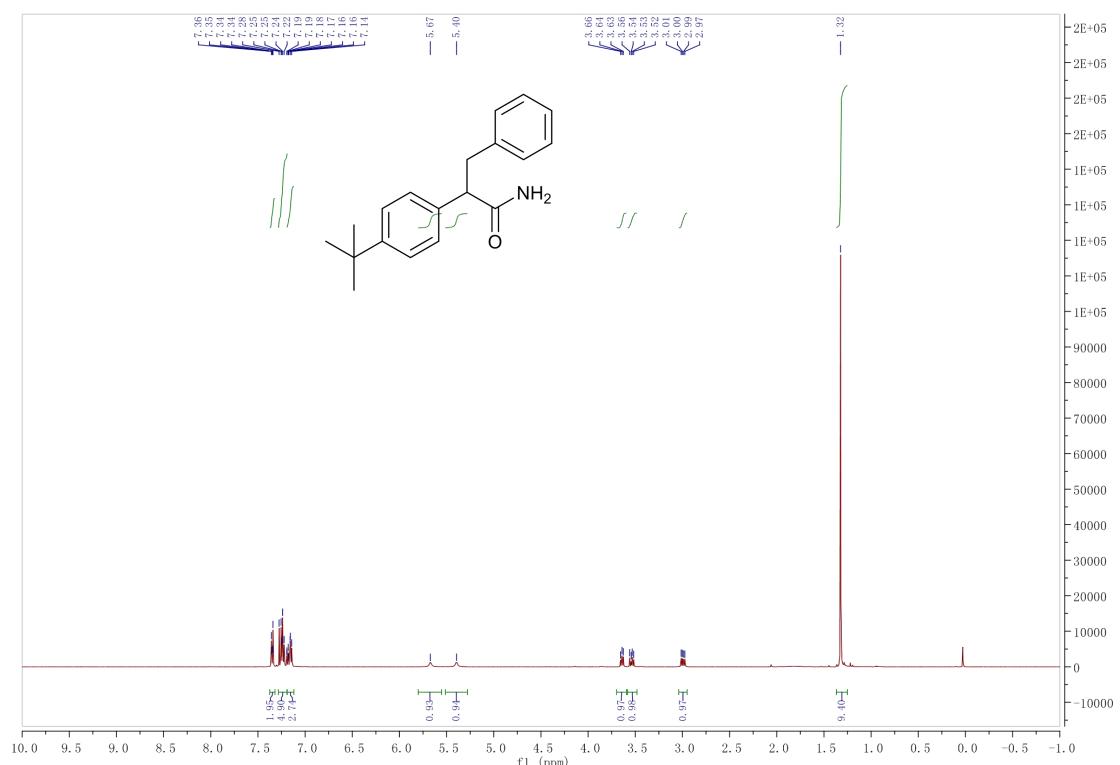
3-phenyl-2-(o-tolyl)propenamide (4ba**)**



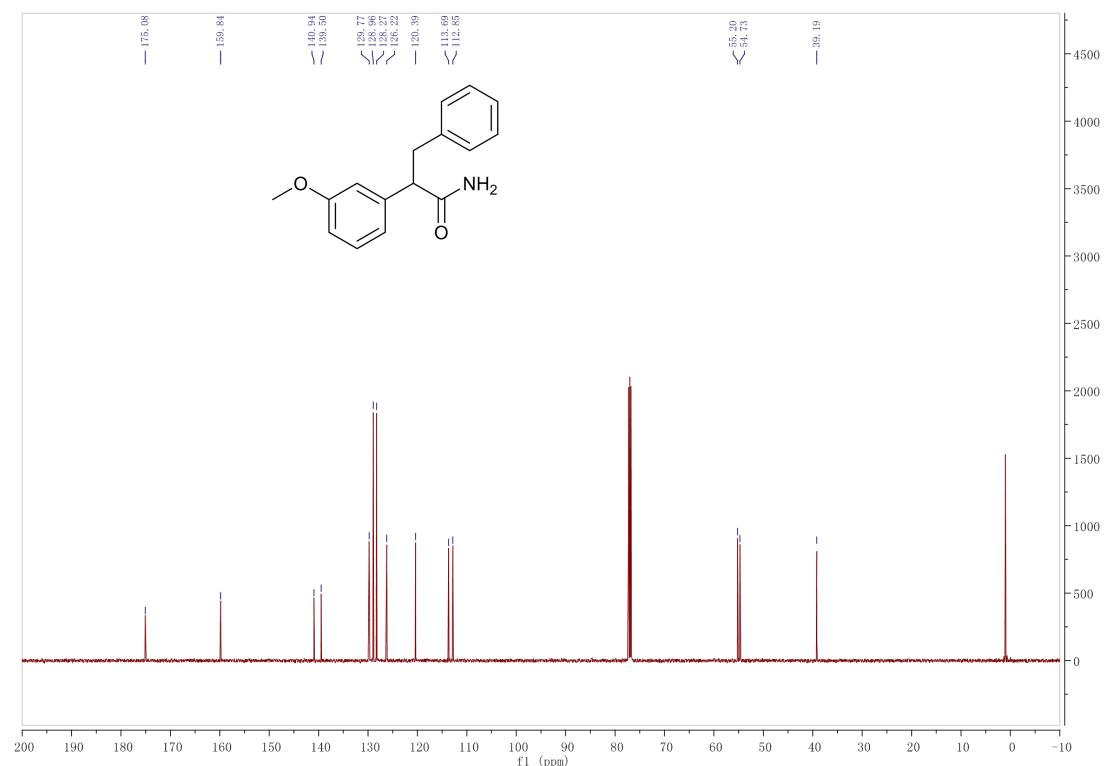
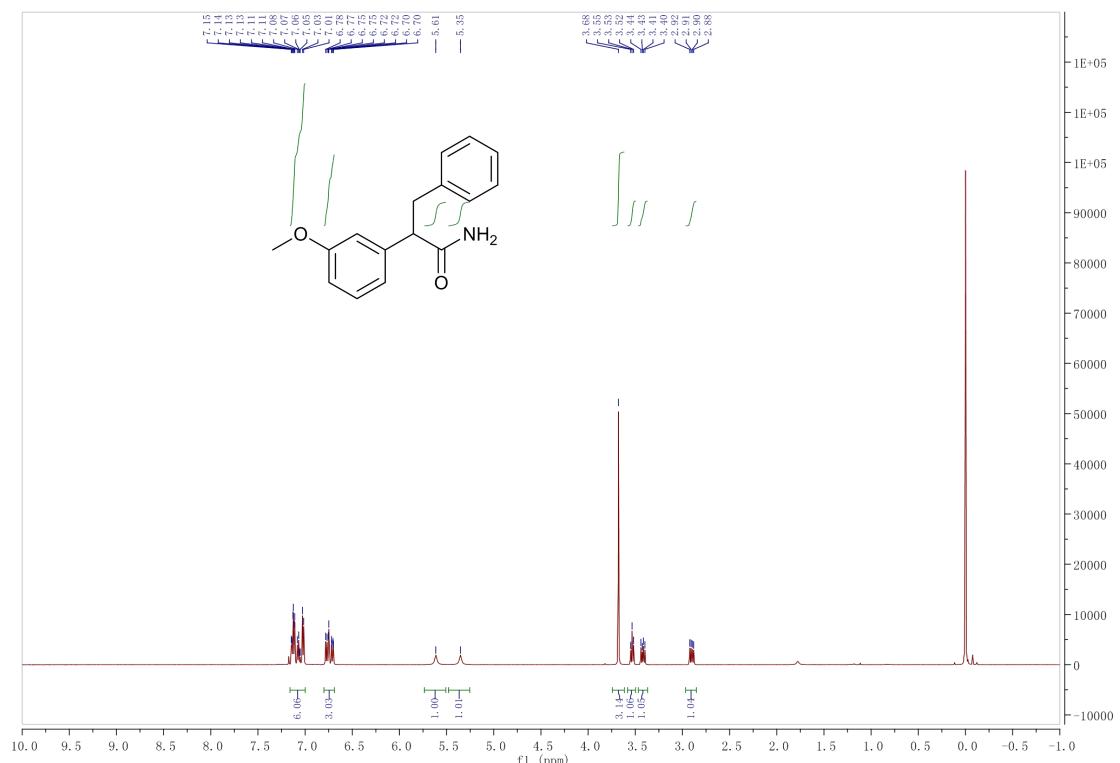
3-phenyl-2-(m-tolyl)propenamide (4bb**)**



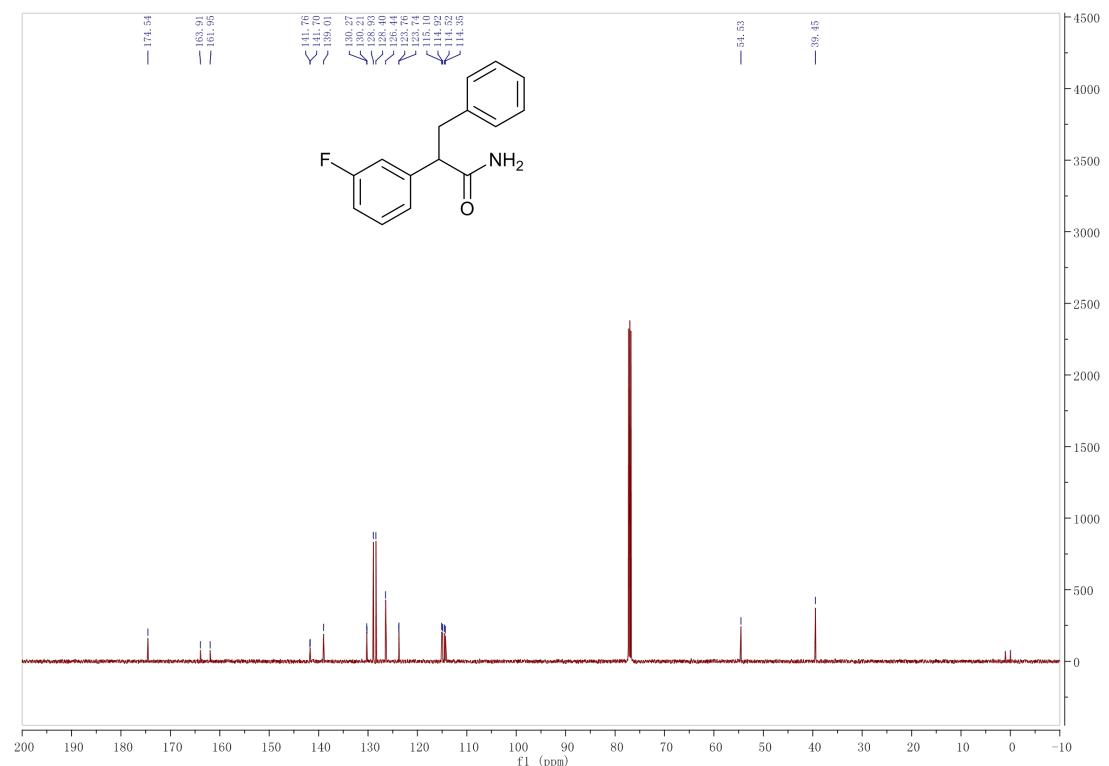
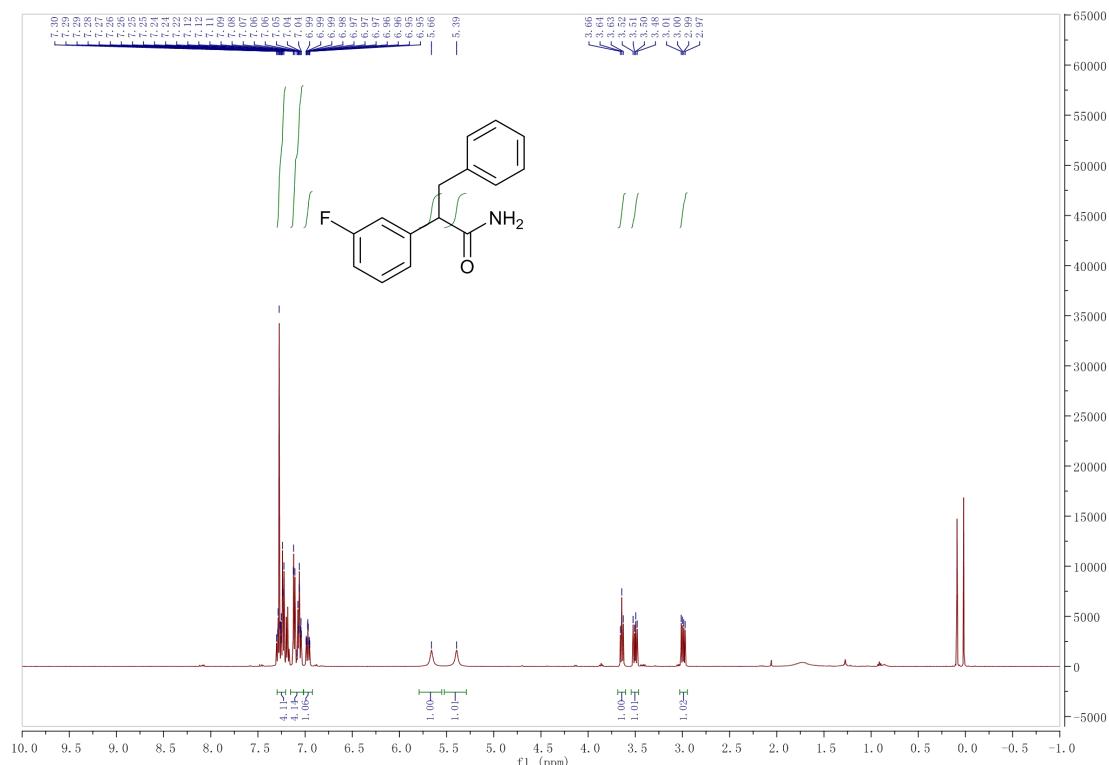
2-(4-(tert-butyl)phenyl)-3-phenylpropanamide (4bc**)**



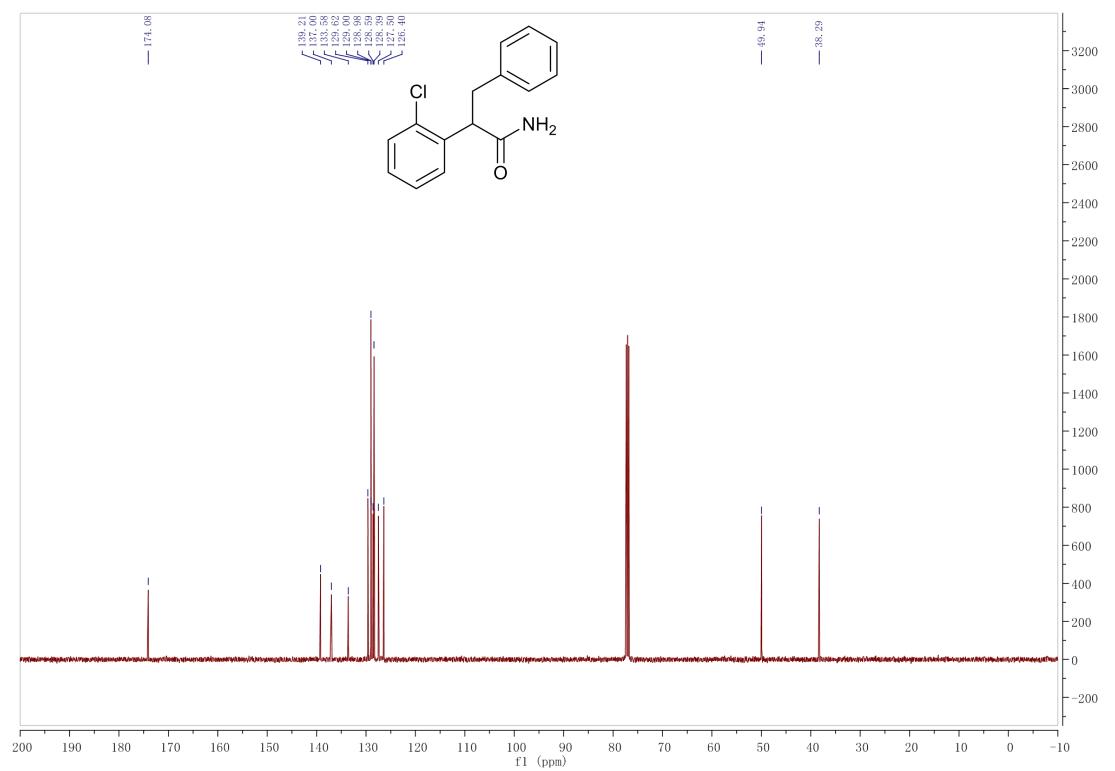
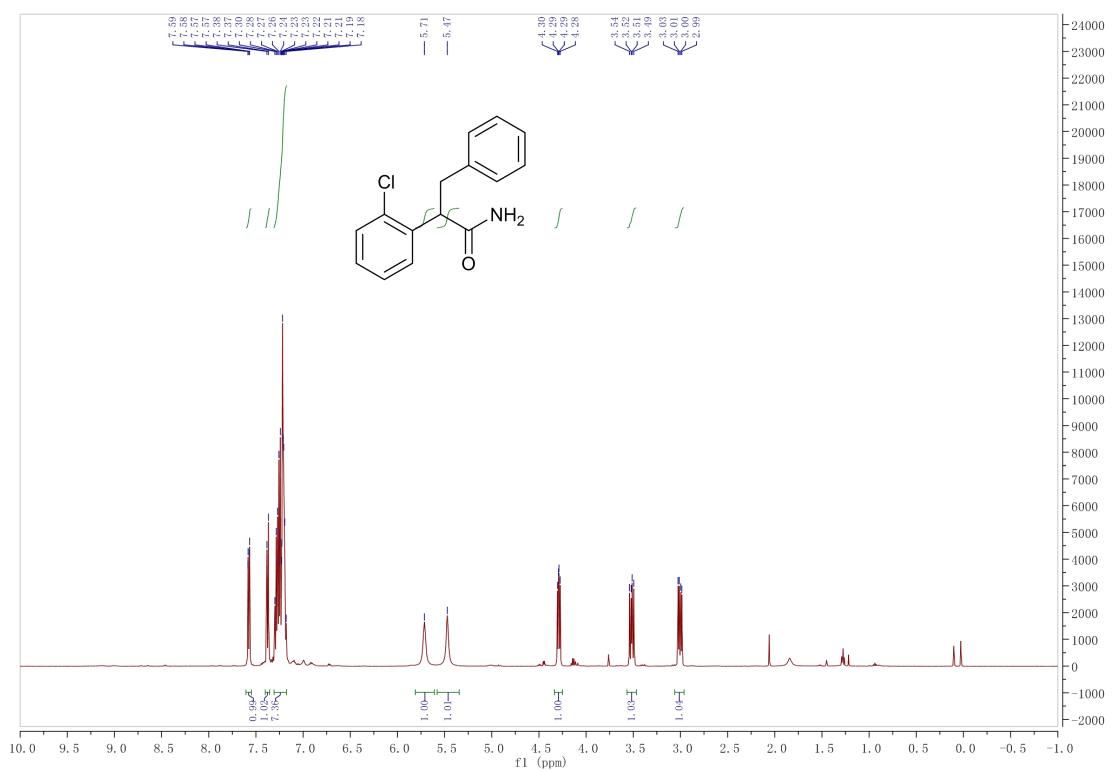
2-(3-methoxyphenyl)-3-phenylpropanamide (4bd**)**



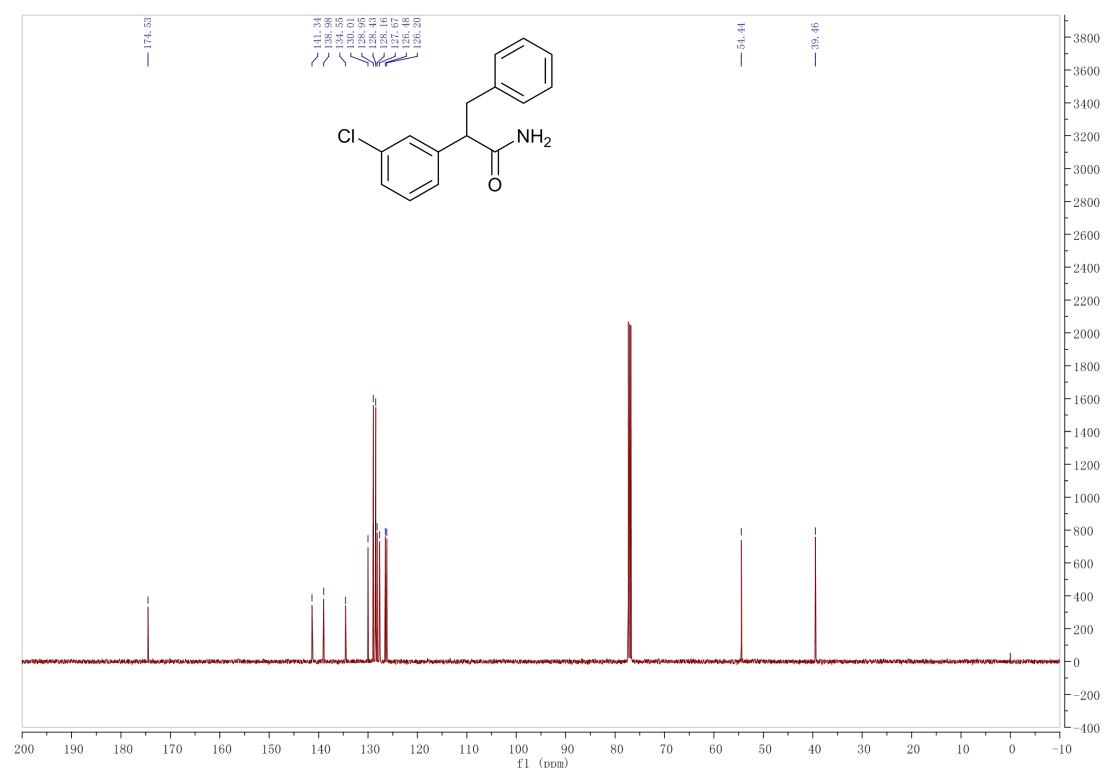
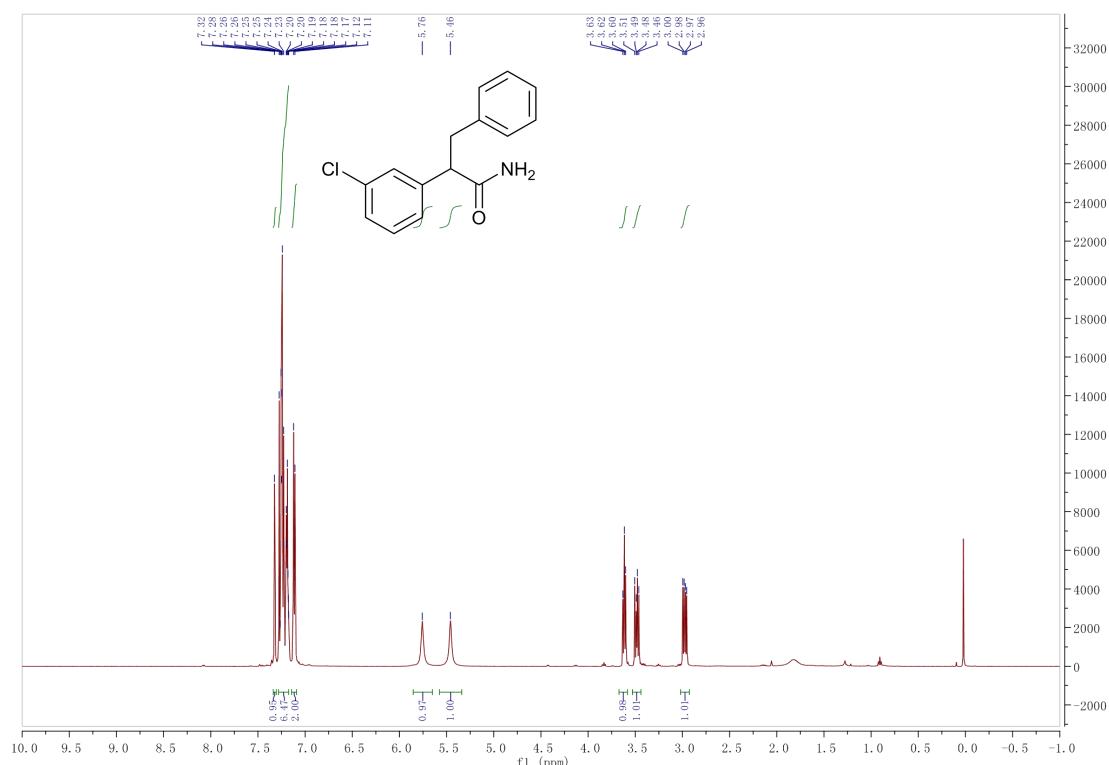
2-(3-fluorophenyl)-3-phenylpropanamide (4be**)**



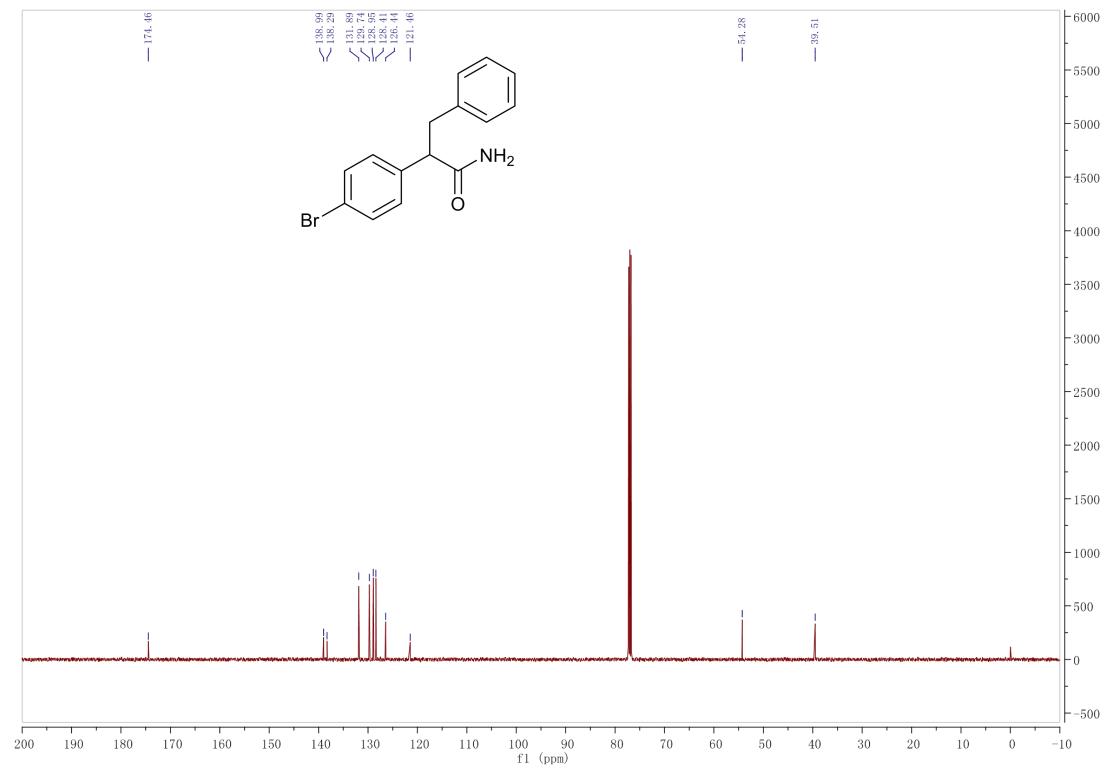
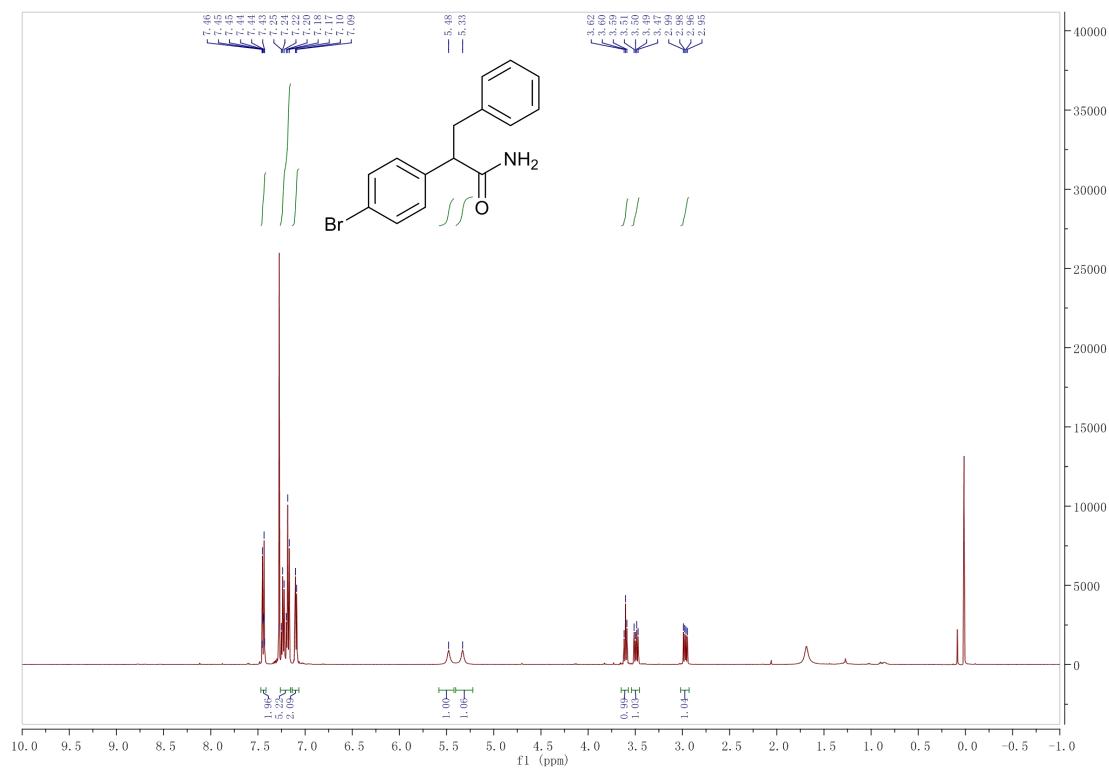
2-(2-chlorophenyl)-3-phenylpropanamide (**4bf**)



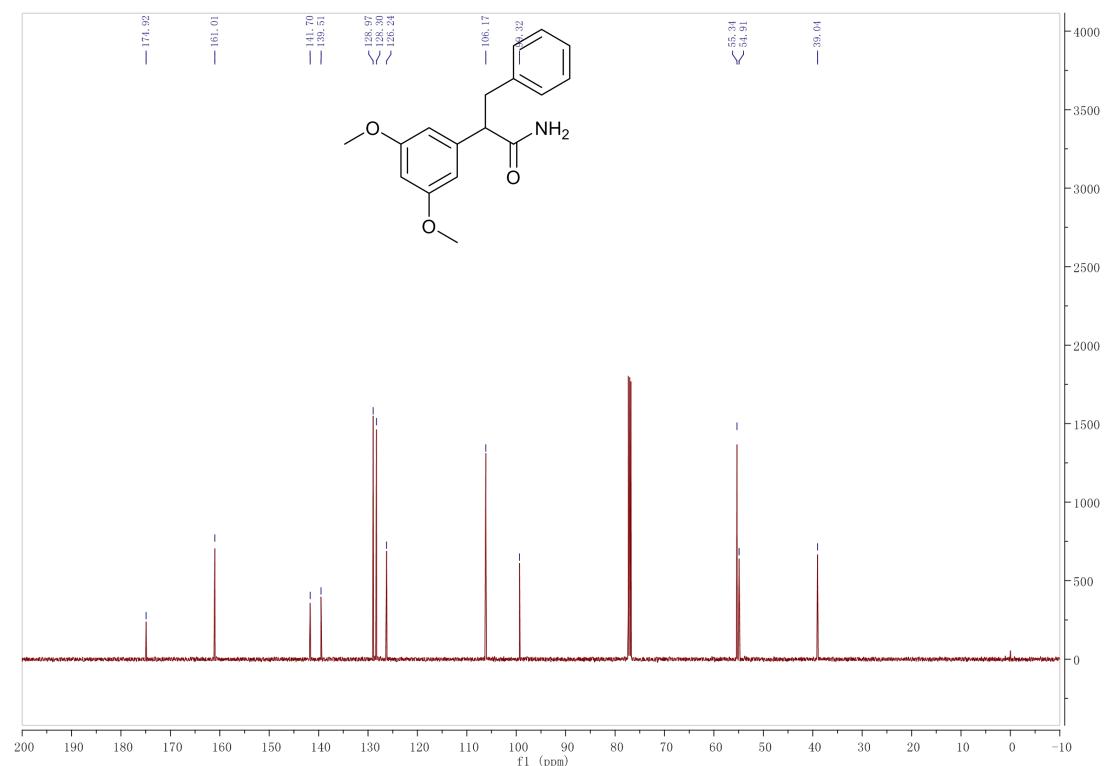
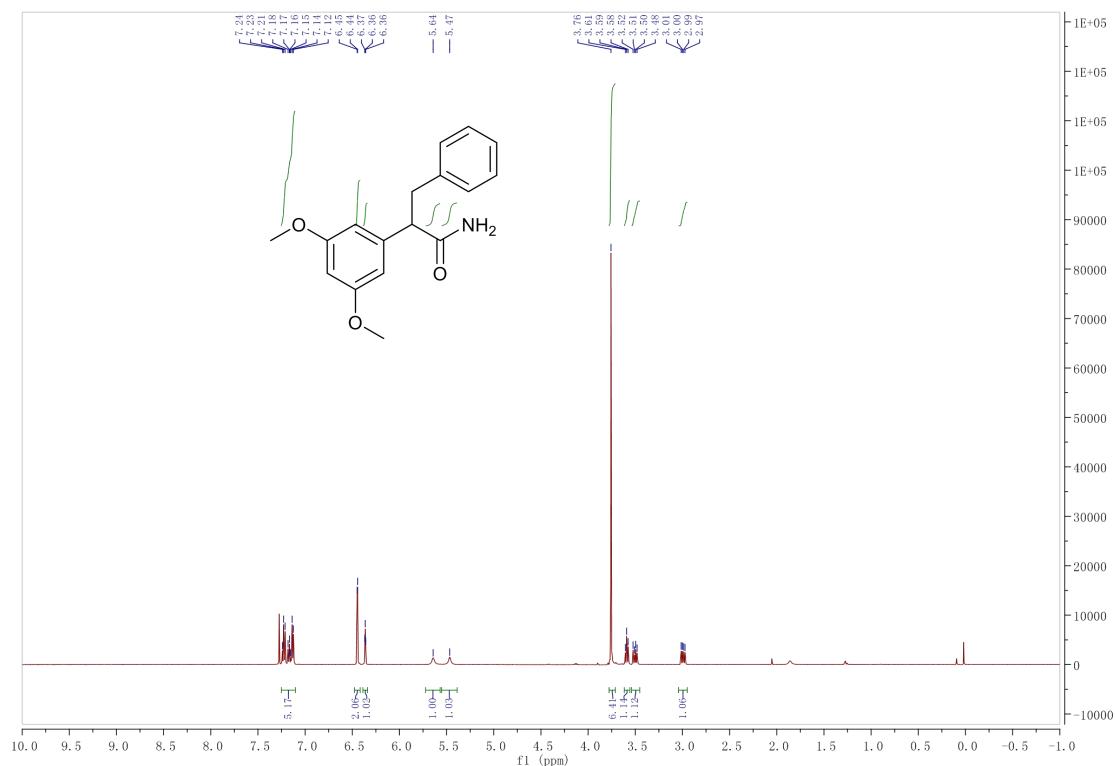
2-(3-chlorophenyl)-3-phenylpropanamide (4bg**)**



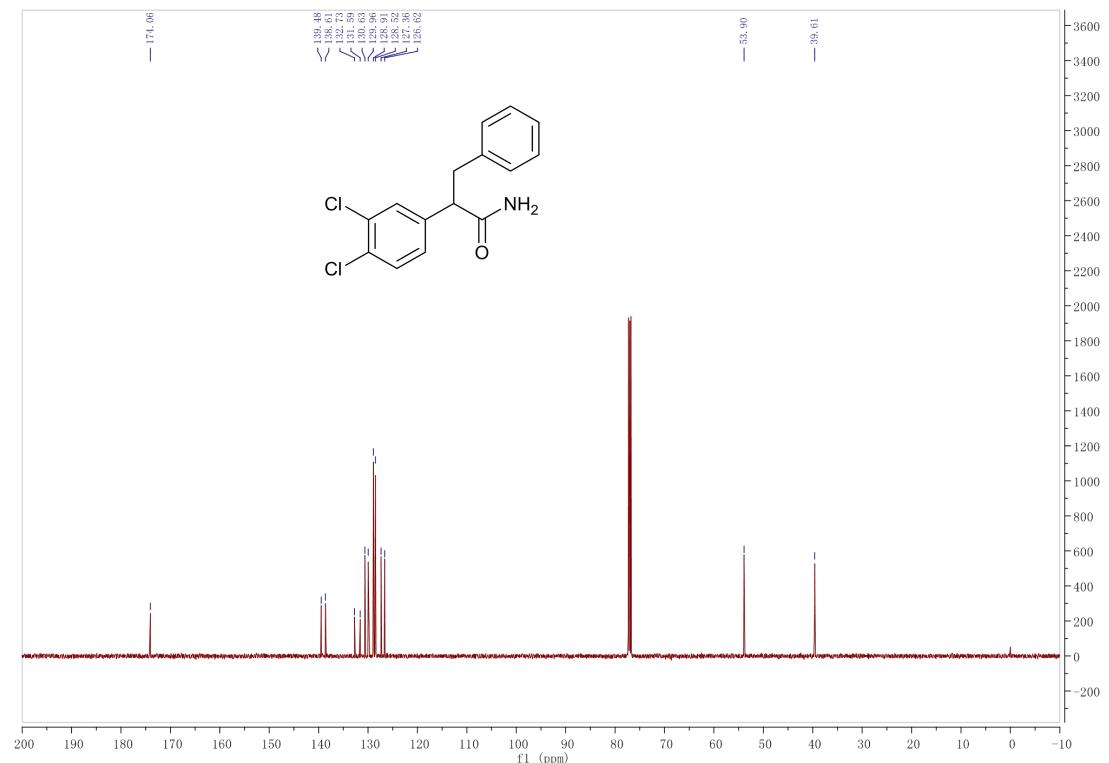
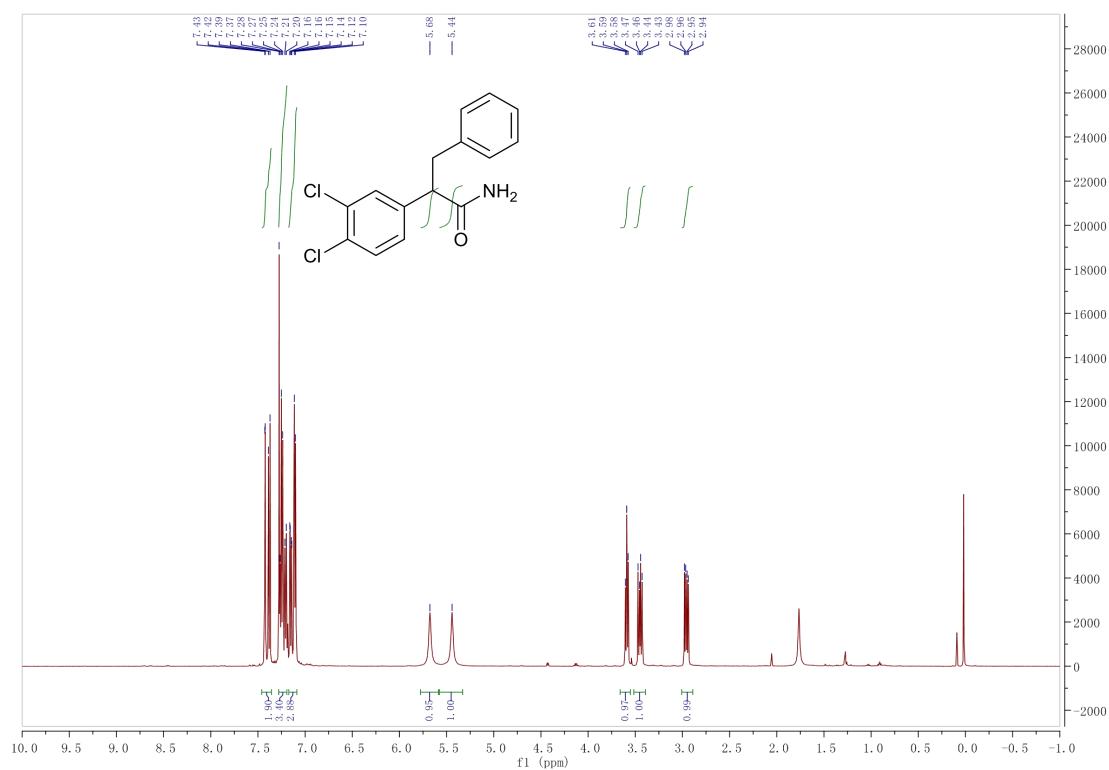
2-(4-bromophenyl)-3-phenylpropanamide (4bh**)**



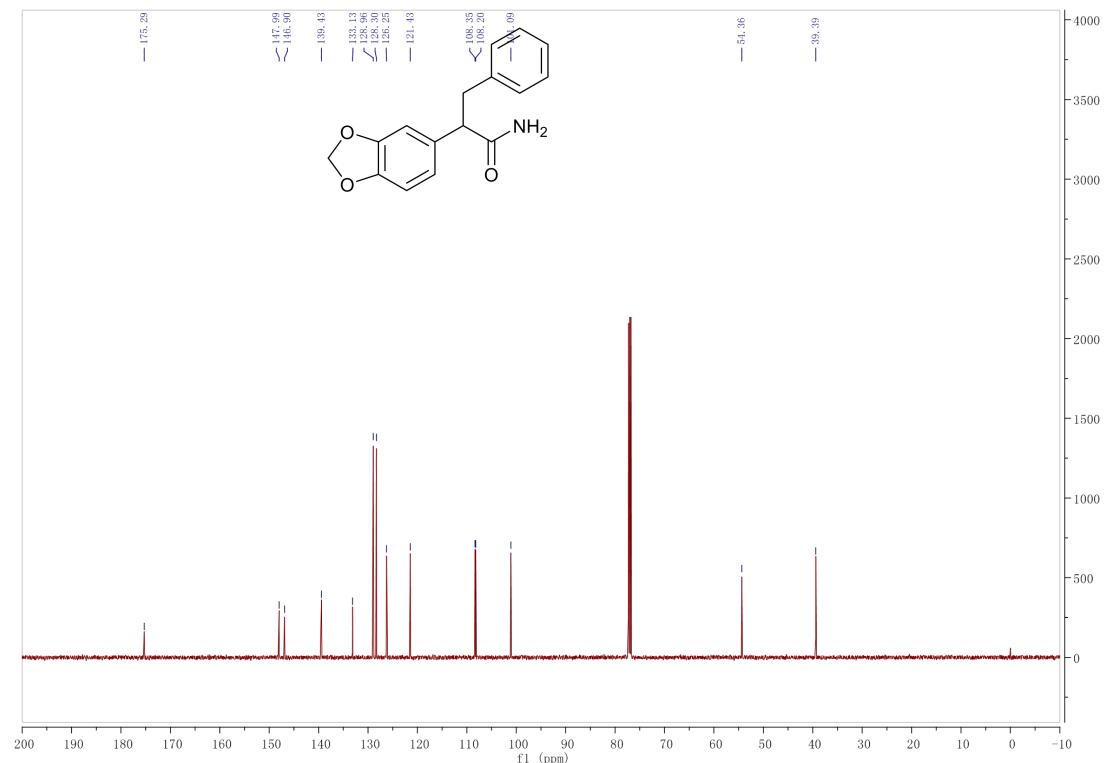
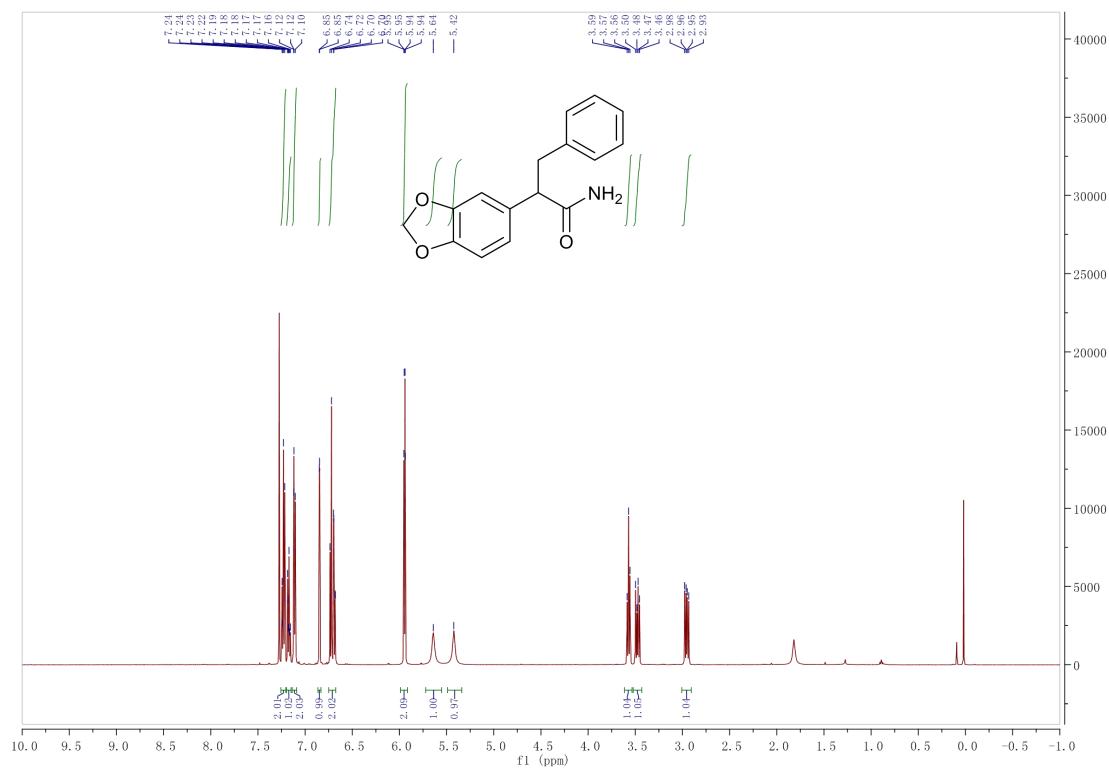
2-(3,5-dimethoxyphenyl)-3-phenylpropanamide (4bi**)**



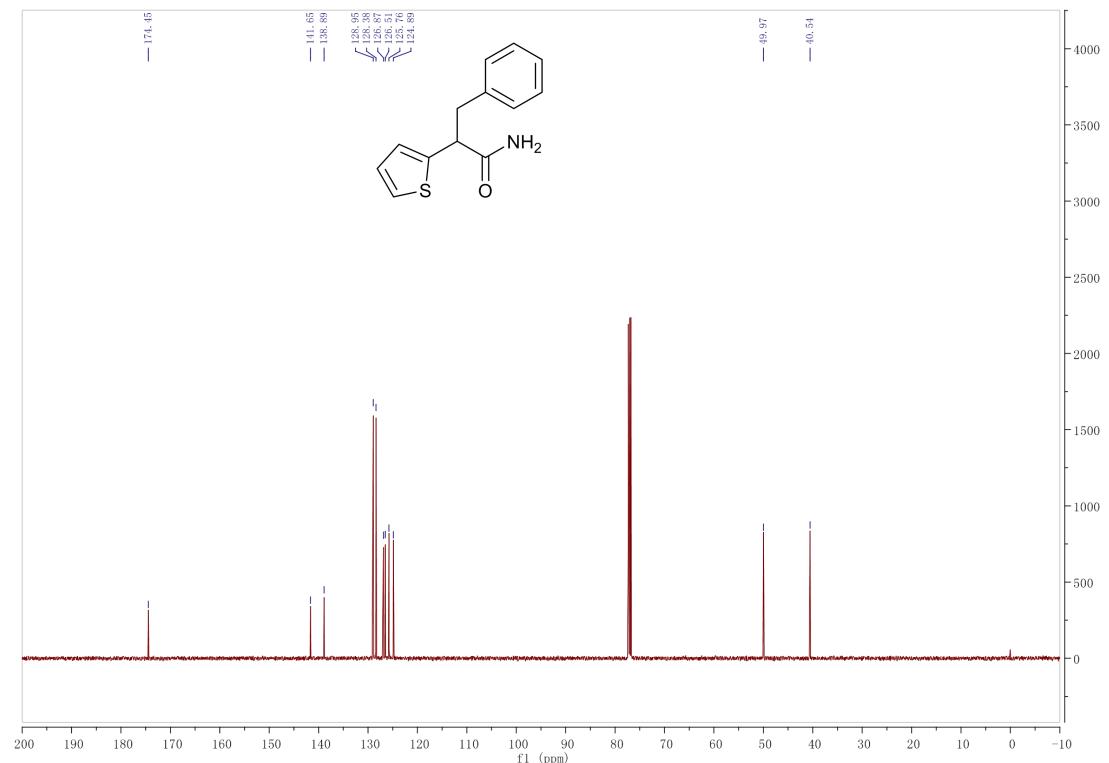
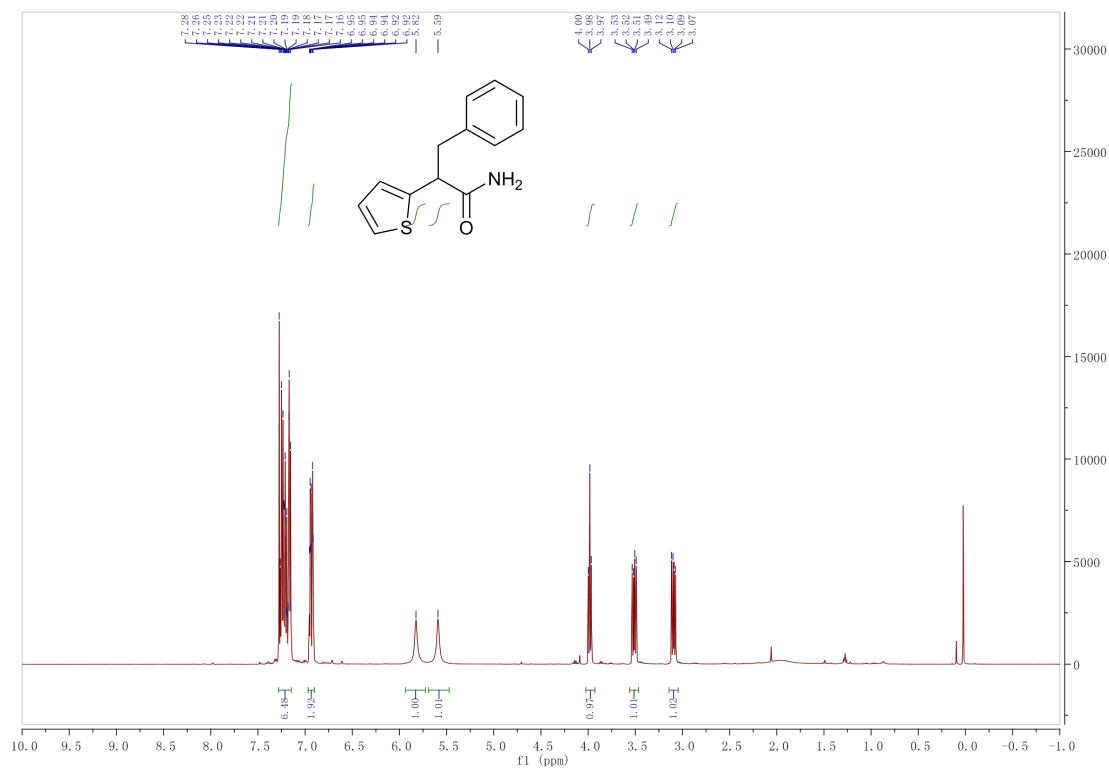
2-(3,4-dichlorophenyl)-3-phenylpropanamide (4bj**)**



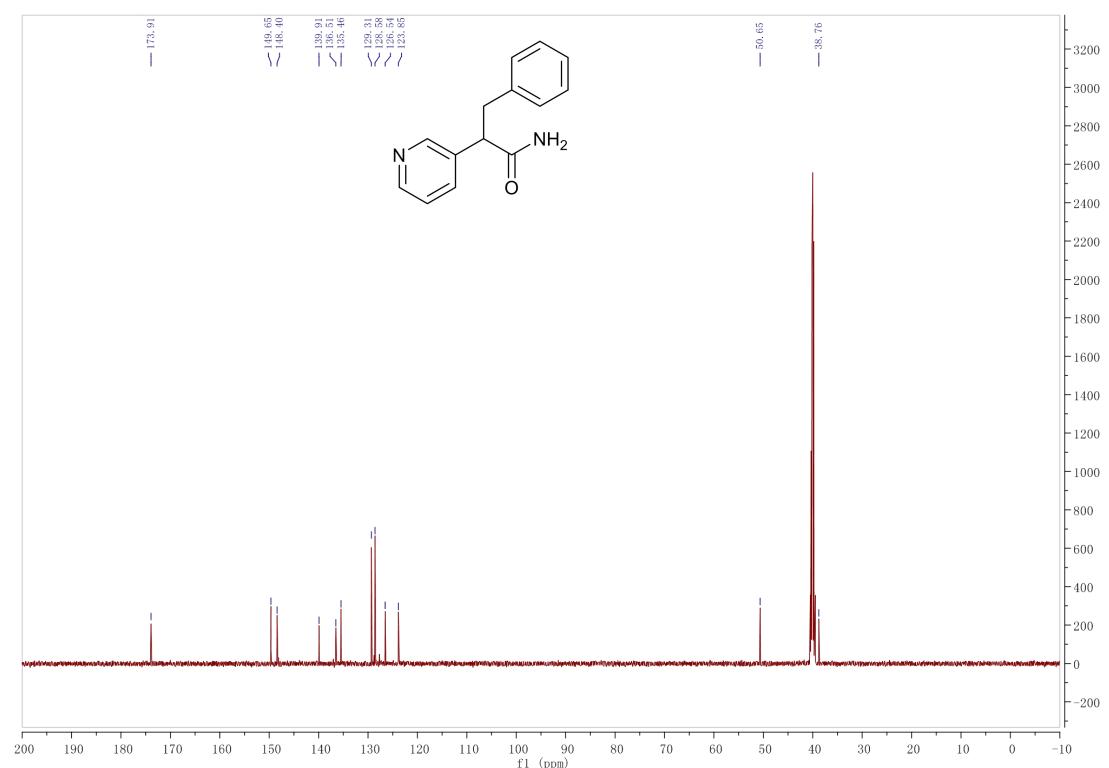
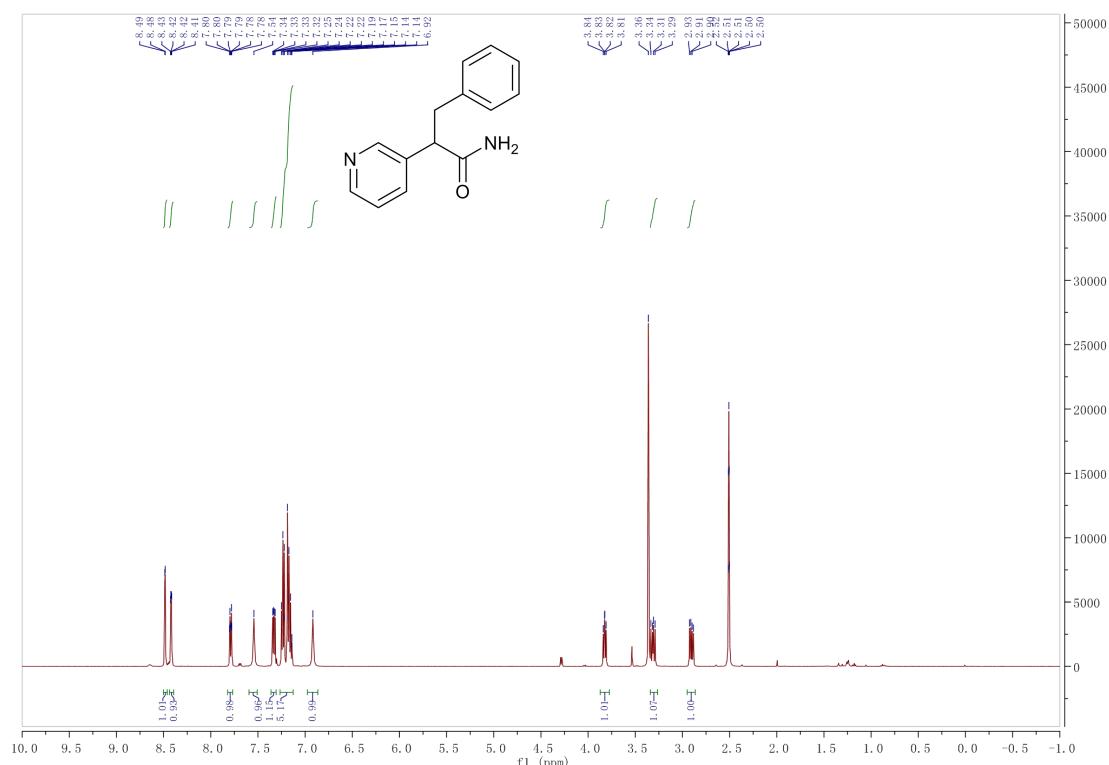
2-(benzo[d][1,3]dioxol-5-yl)-3-phenylpropanamide (4bk**)**



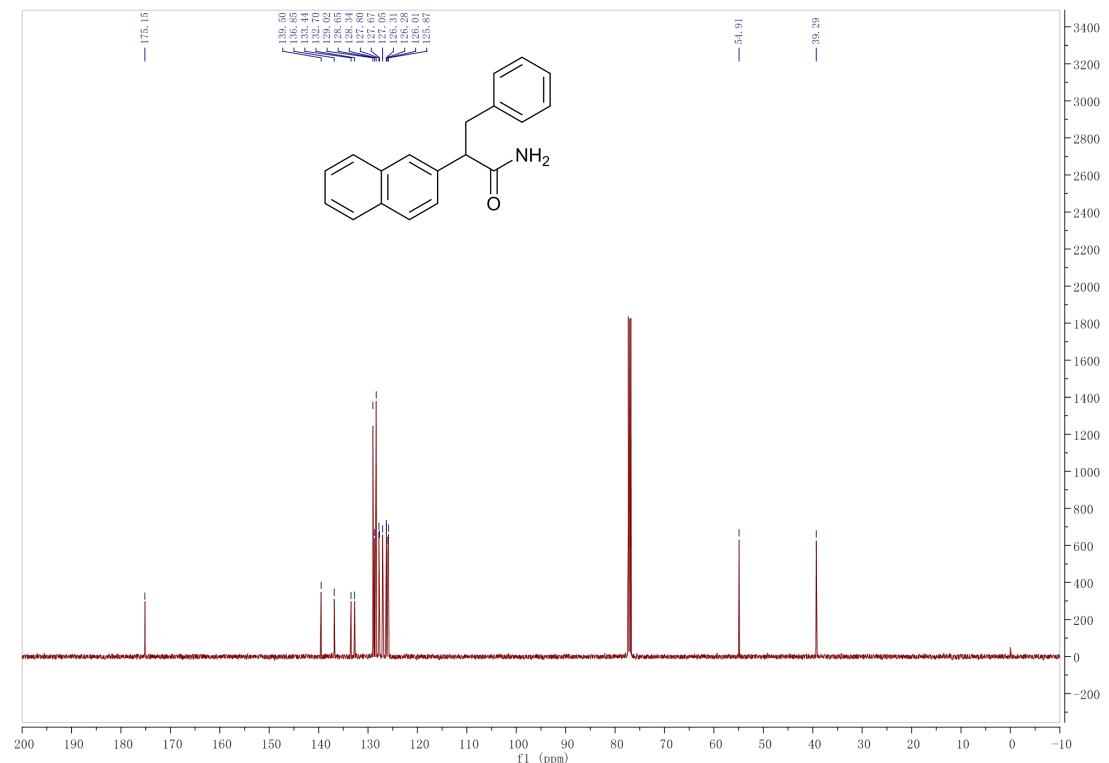
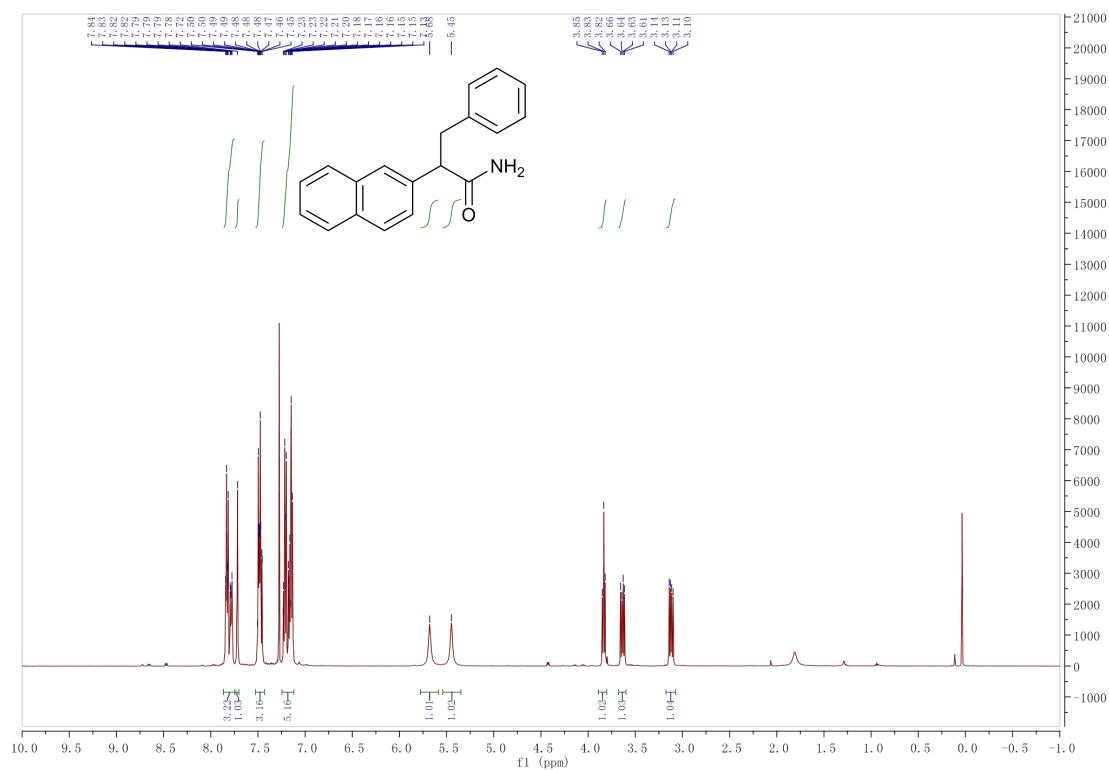
3-phenyl-2-(thiophen-2-yl)propenamide (4bl**)**



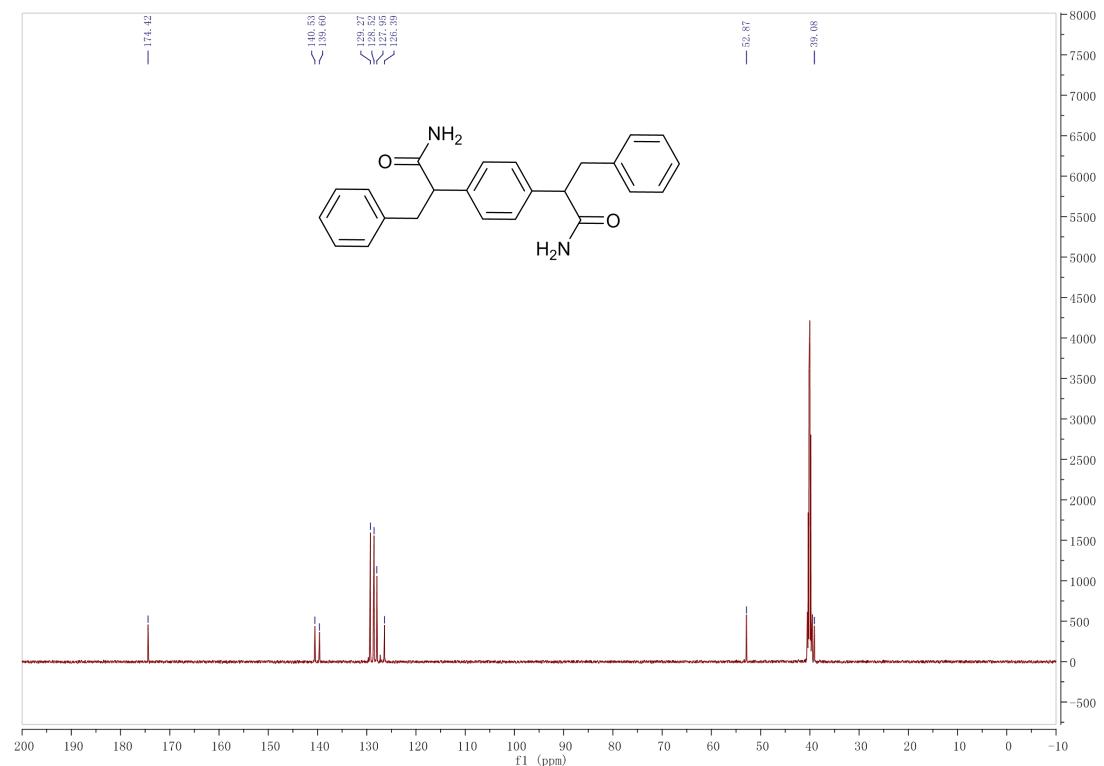
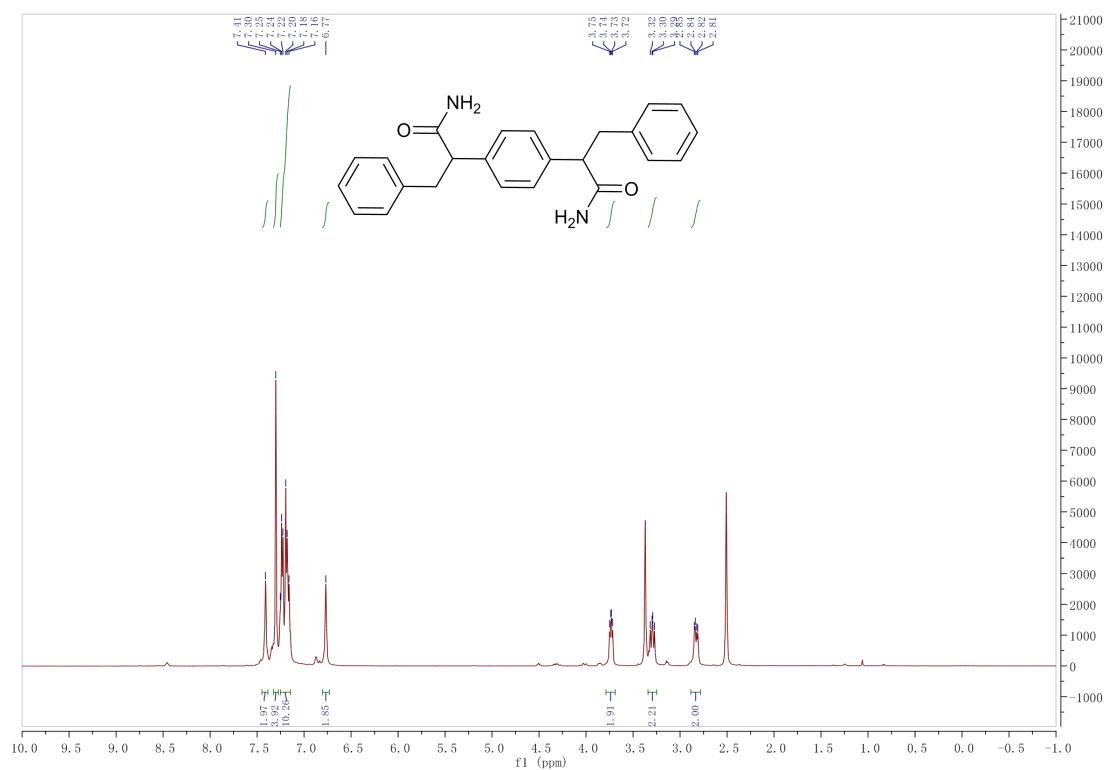
3-phenyl-2-(pyridin-3-yl)propenamide (4bm**)**



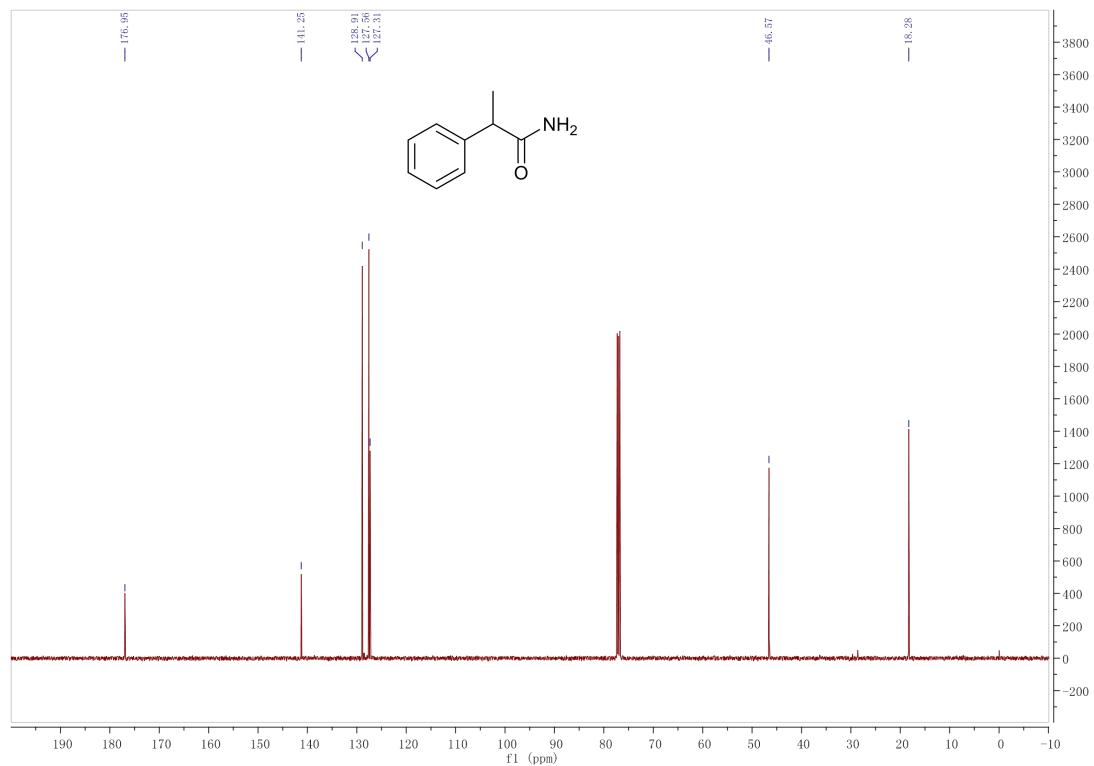
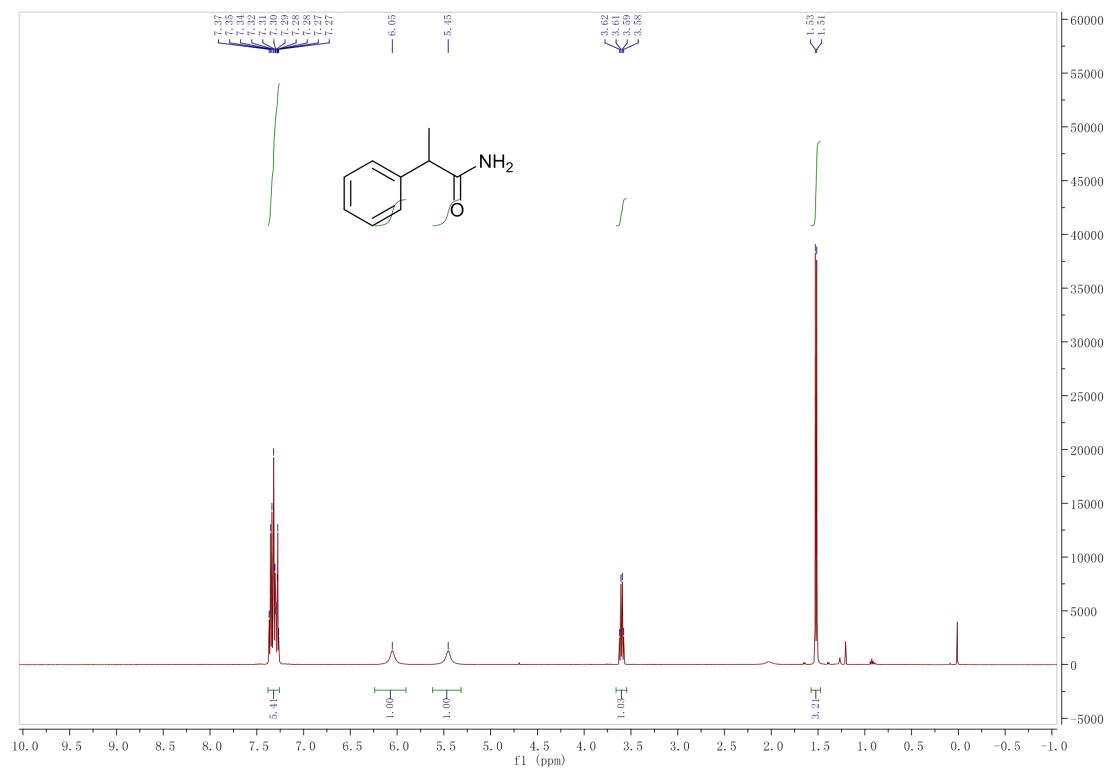
2-(naphthalen-2-yl)-3-phenylpropanamide (4bn**)**



2,2'-(1,4-phenylene)bis(3-phenylpropanamide) (4bo**)**

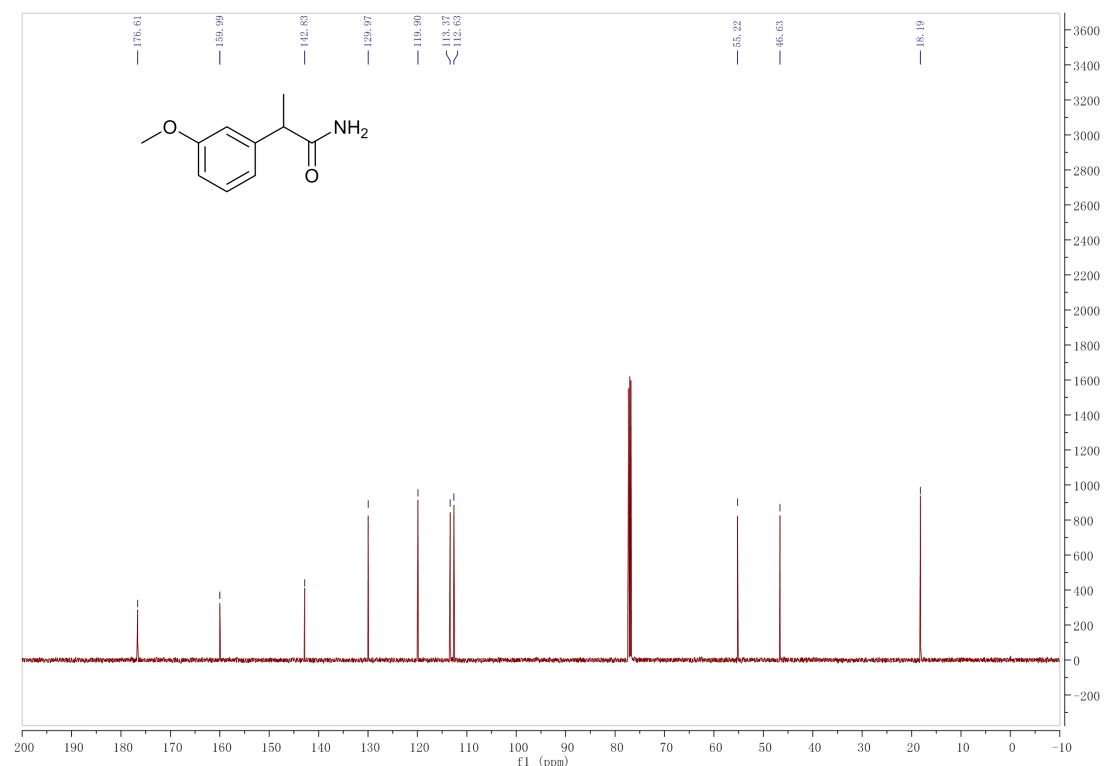
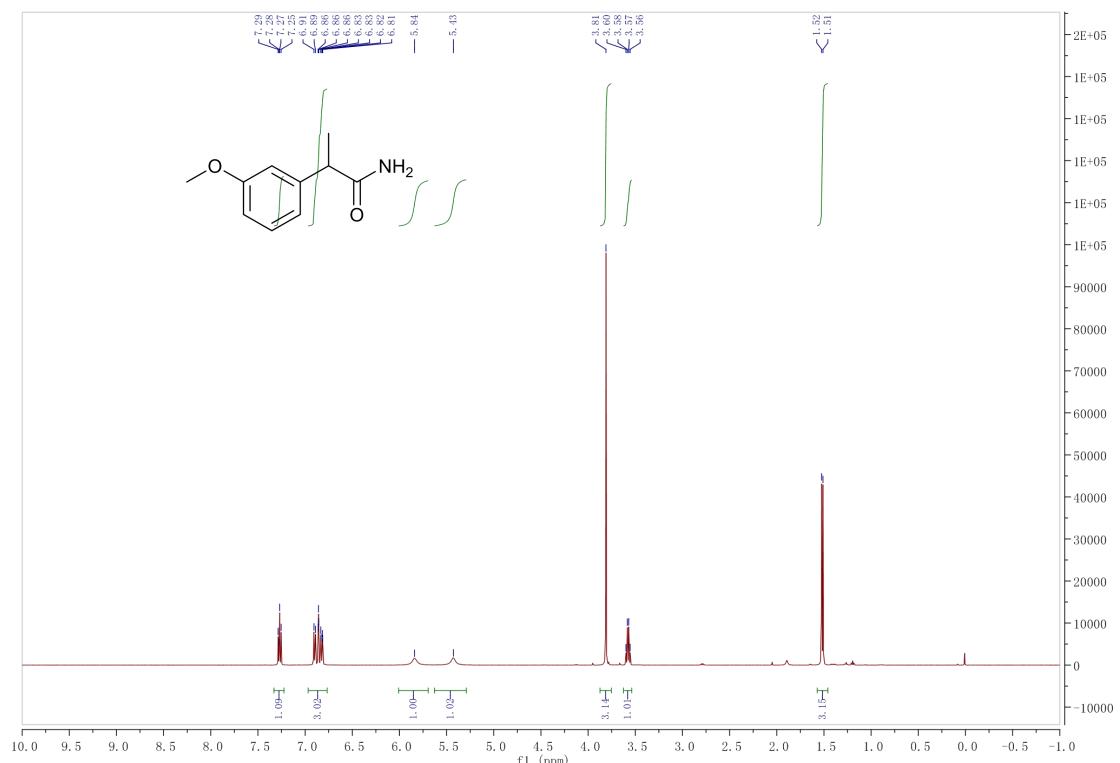


2-phenylpropanamide (4ca)

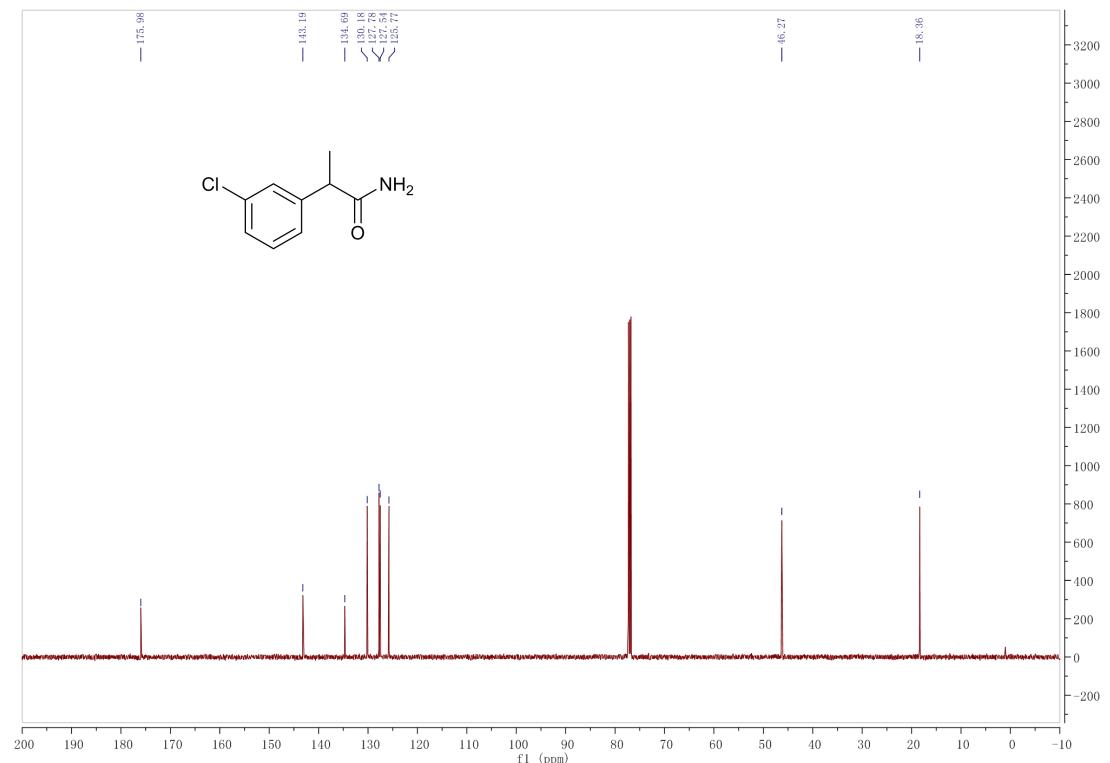
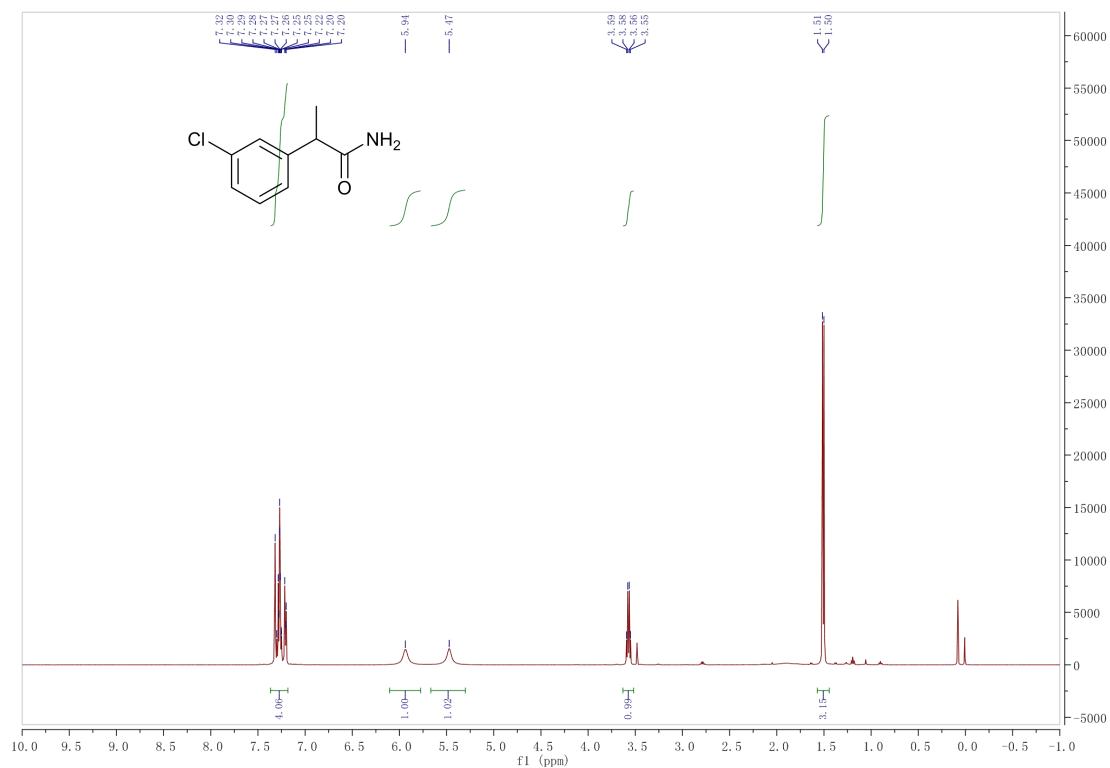


2-(3-methoxyphenyl)propenamide

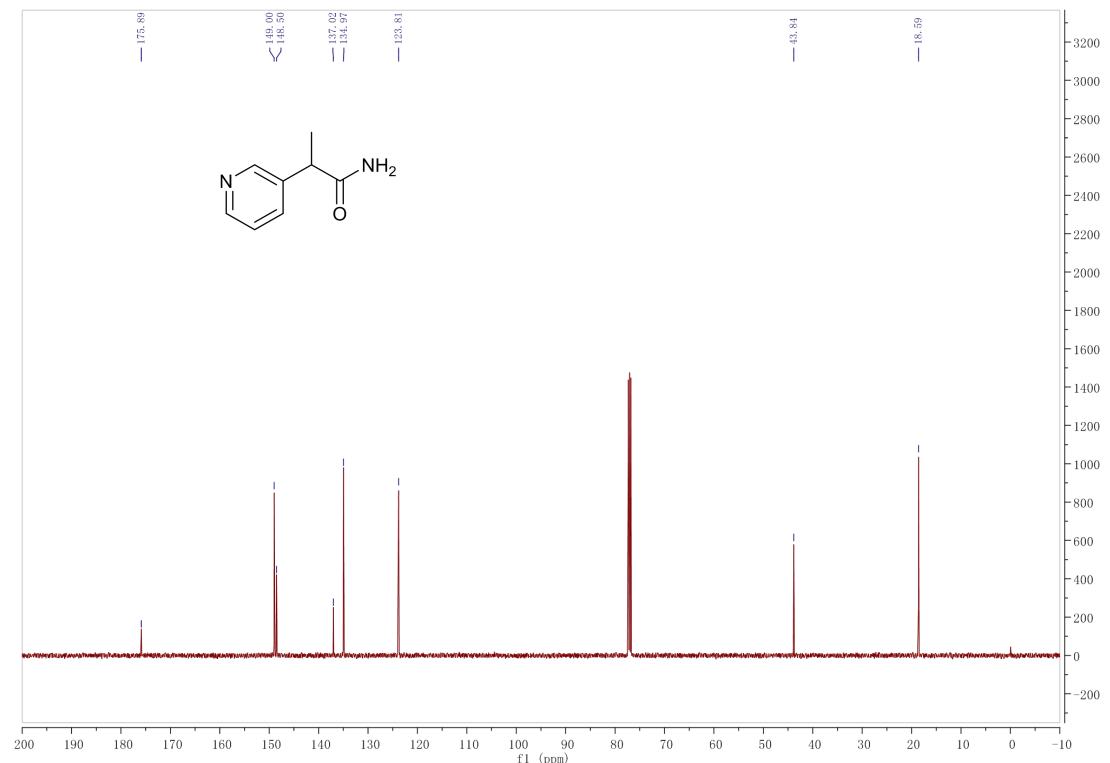
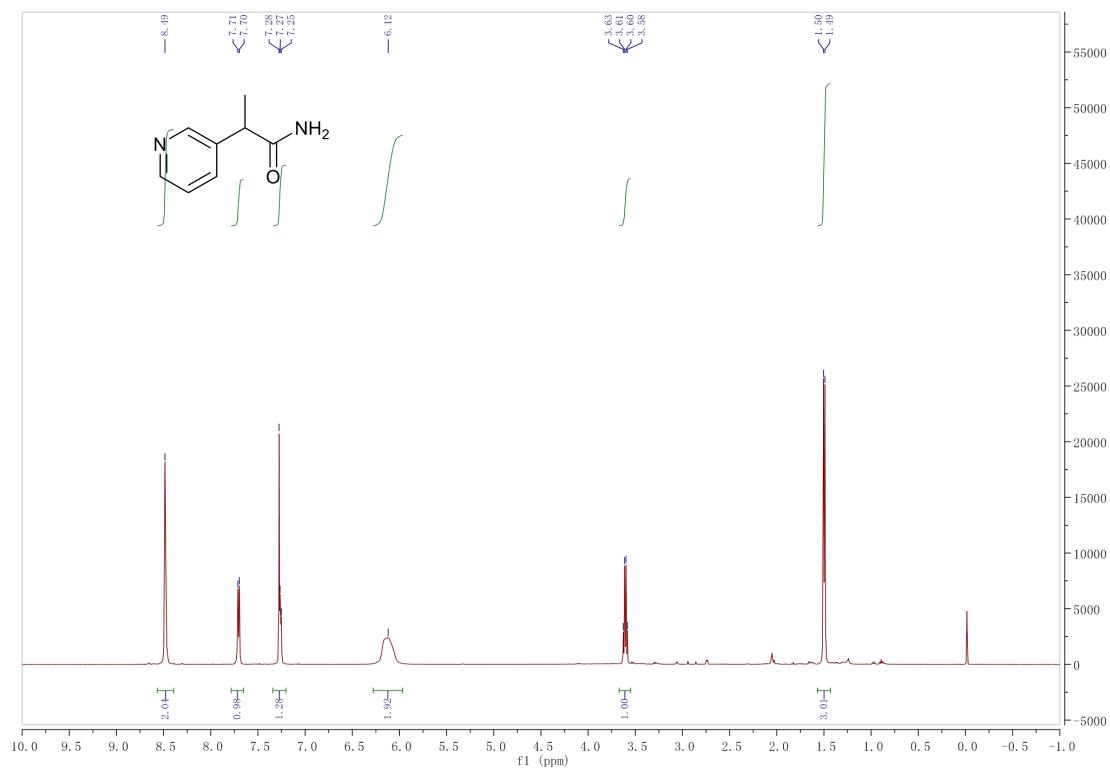
(4cb)



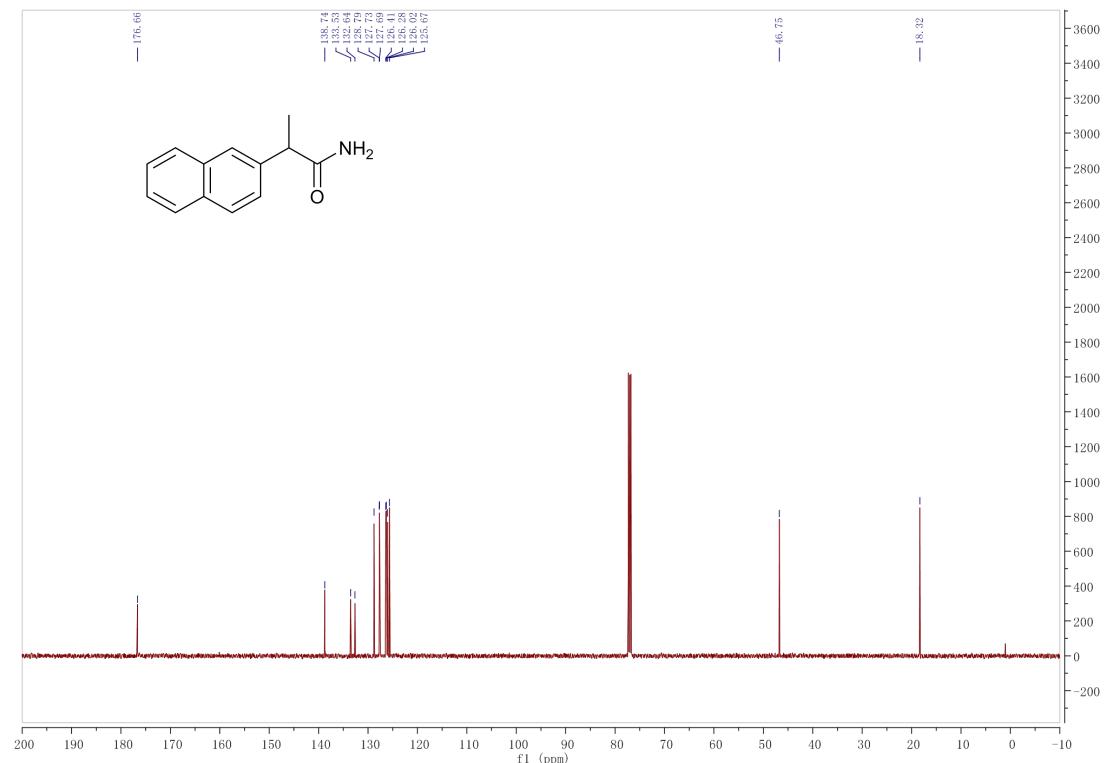
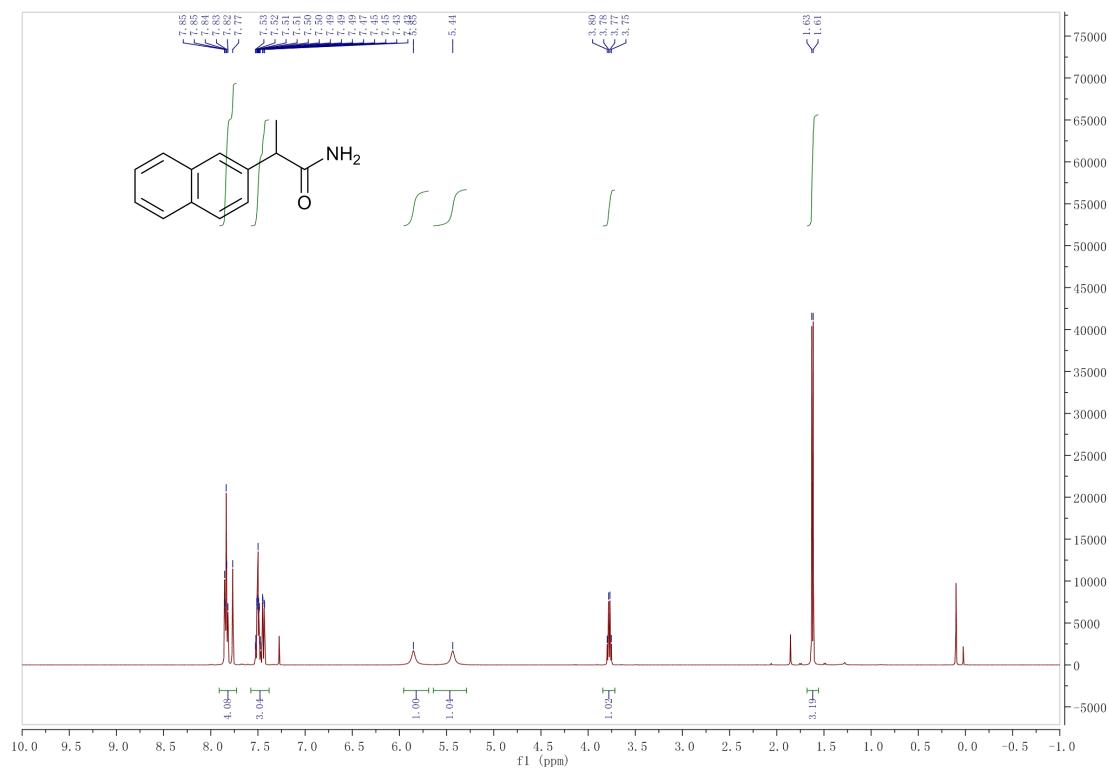
2-(3-chlorophenyl)propenamide (4cc**)**



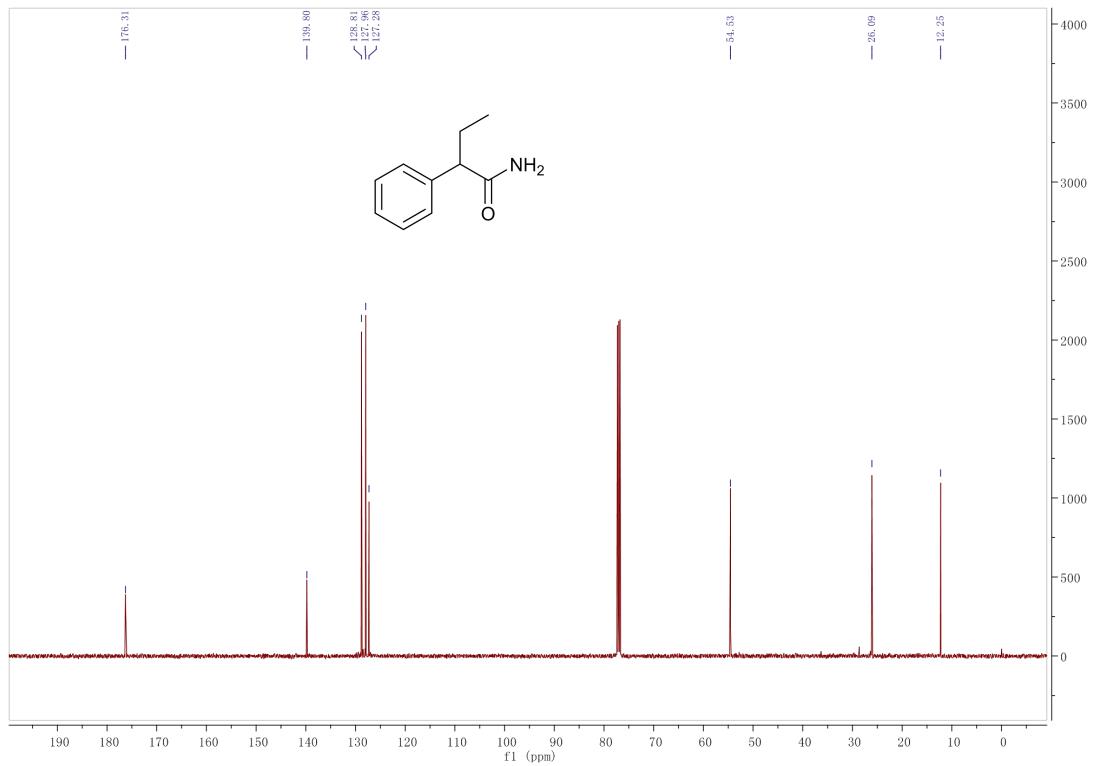
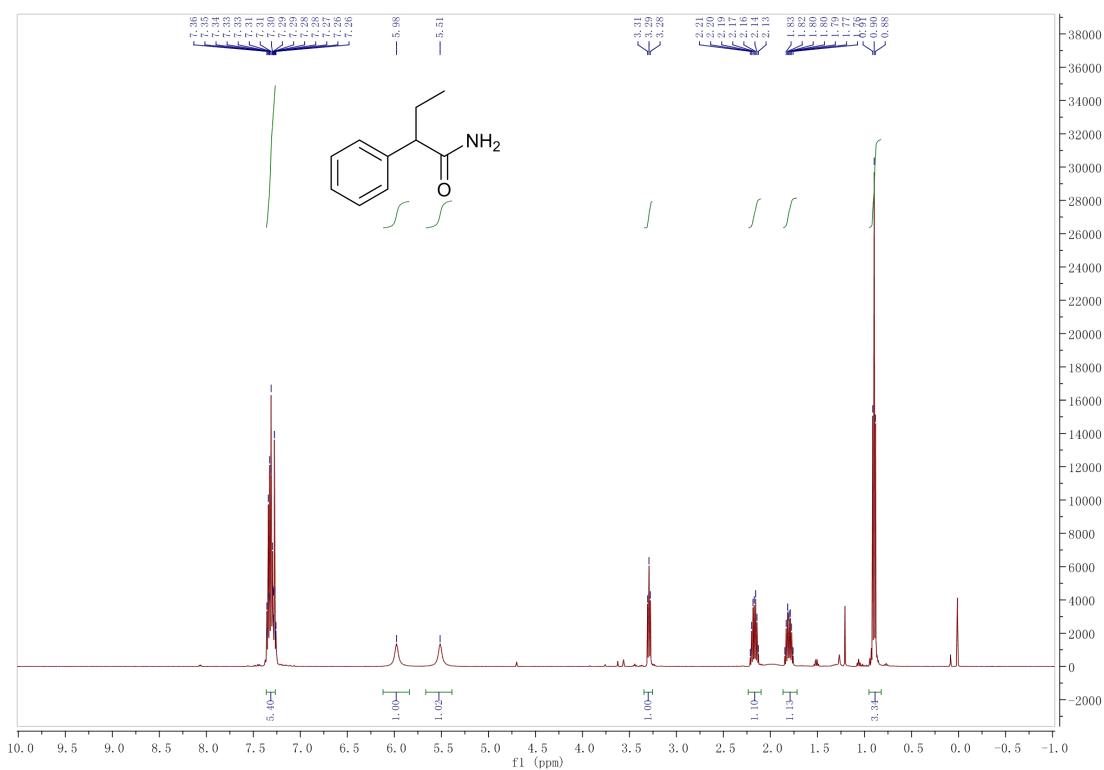
2-(pyridin-3-yl)propenamide (4cd**)**



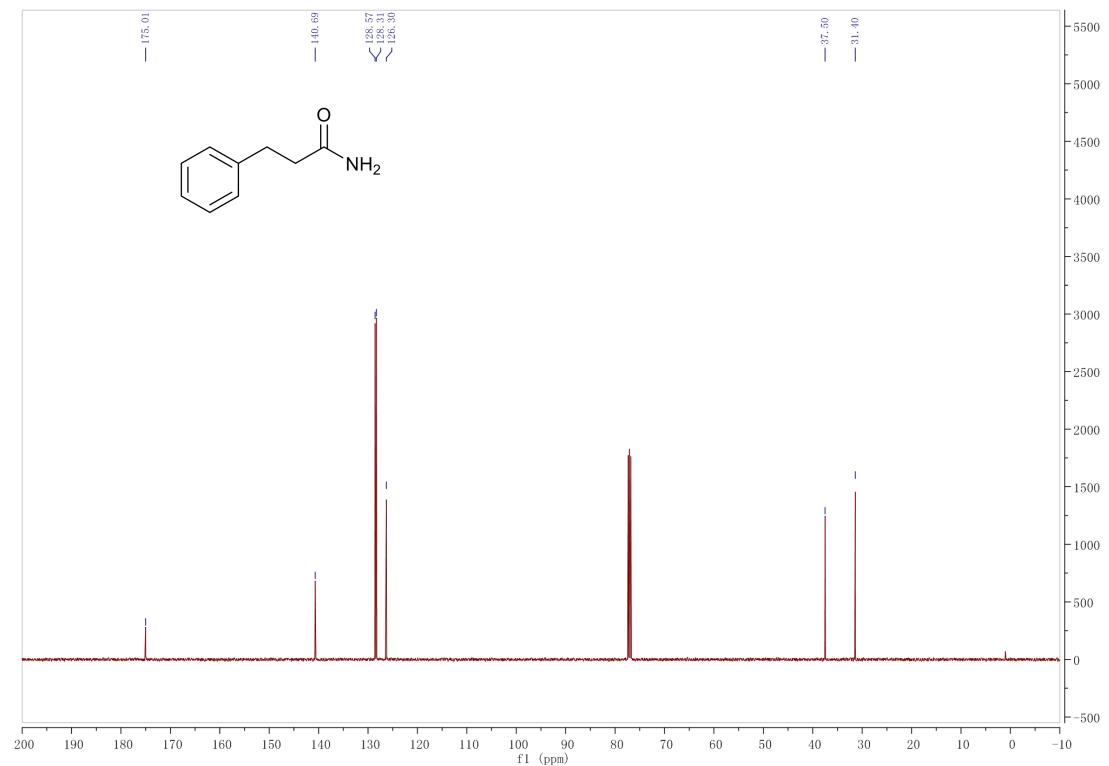
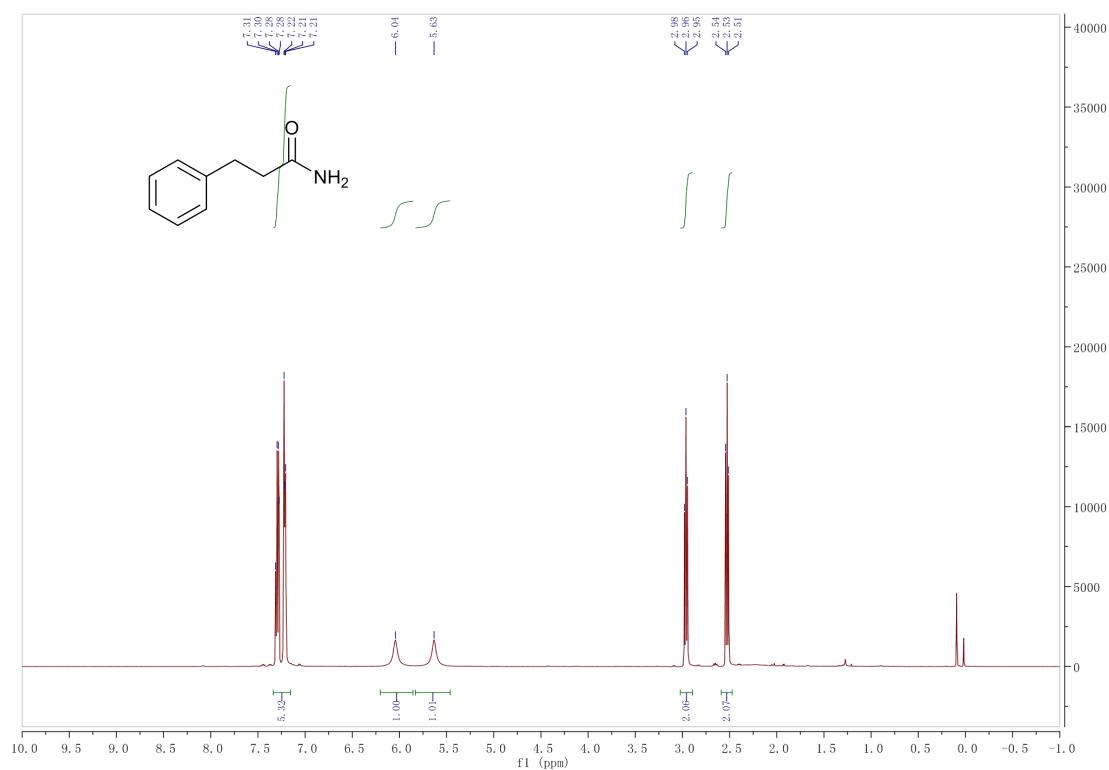
2-(naphthalen-2-yl)propenamide (4ce**)**



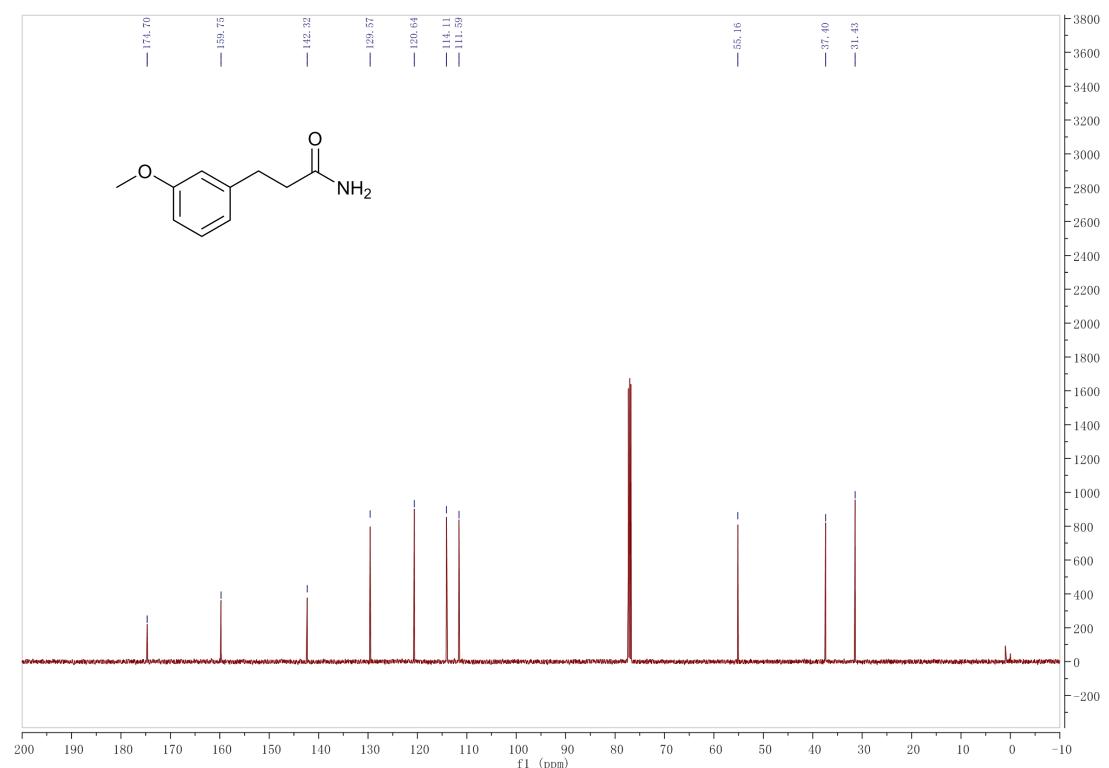
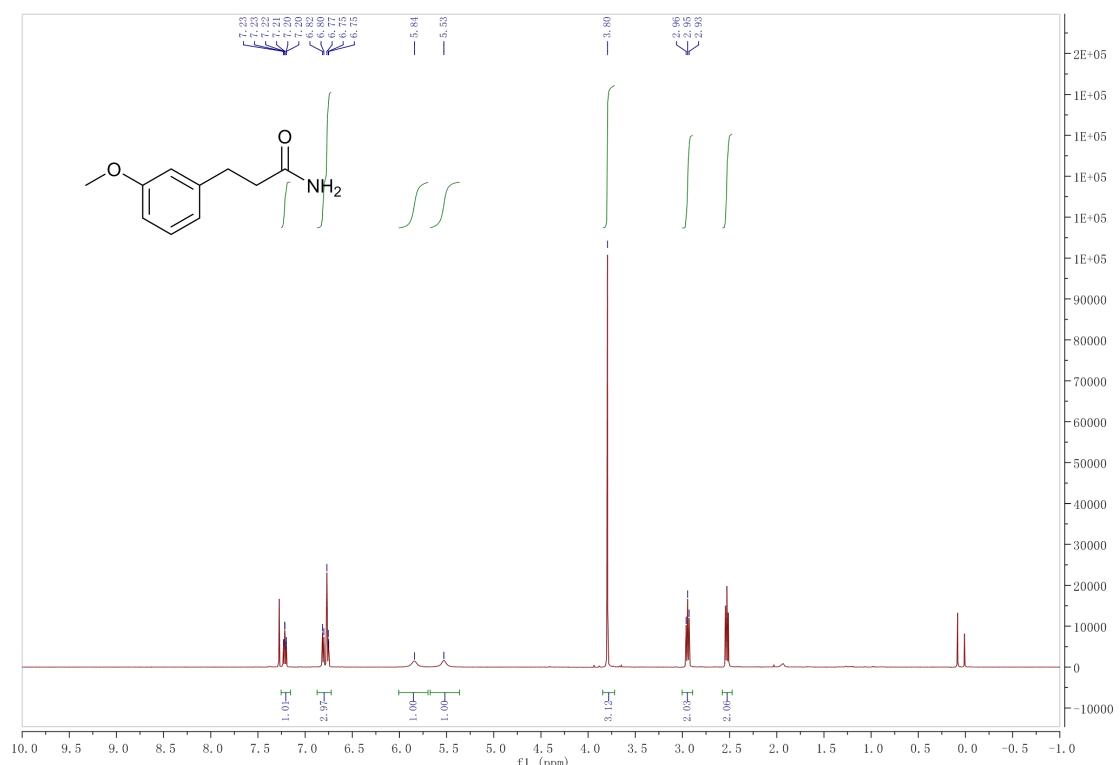
2-phenylbutanamide (4cf)



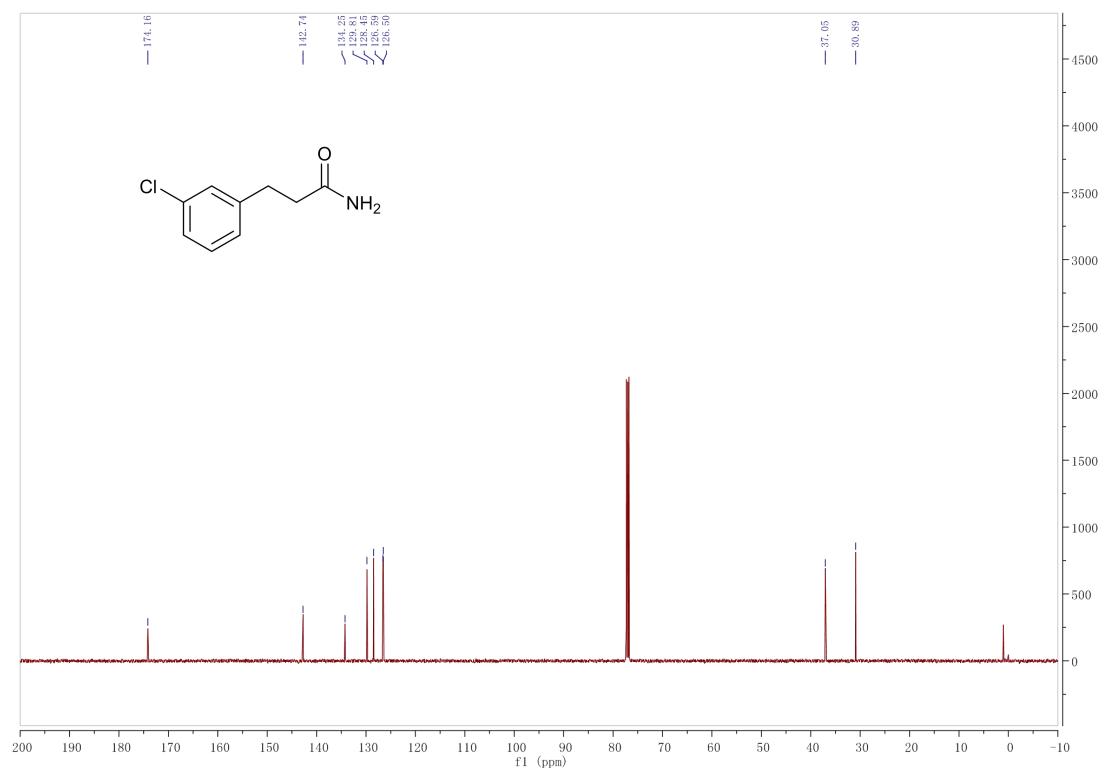
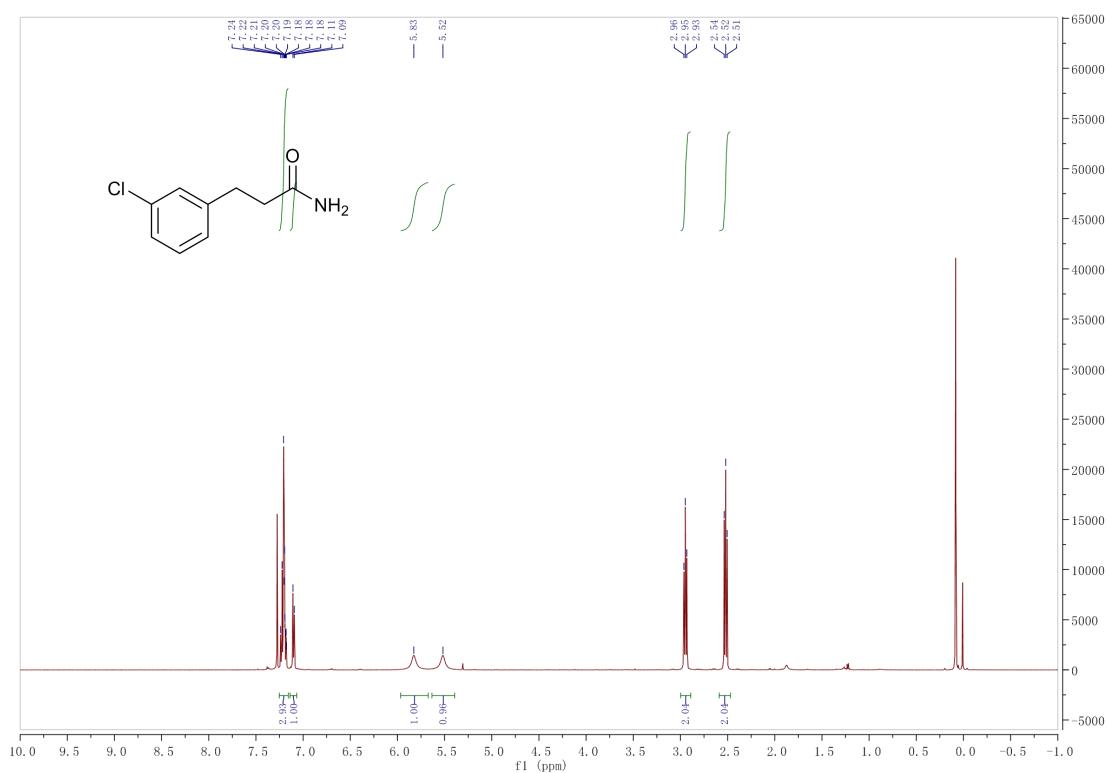
3-phenylpropanamide (4da)



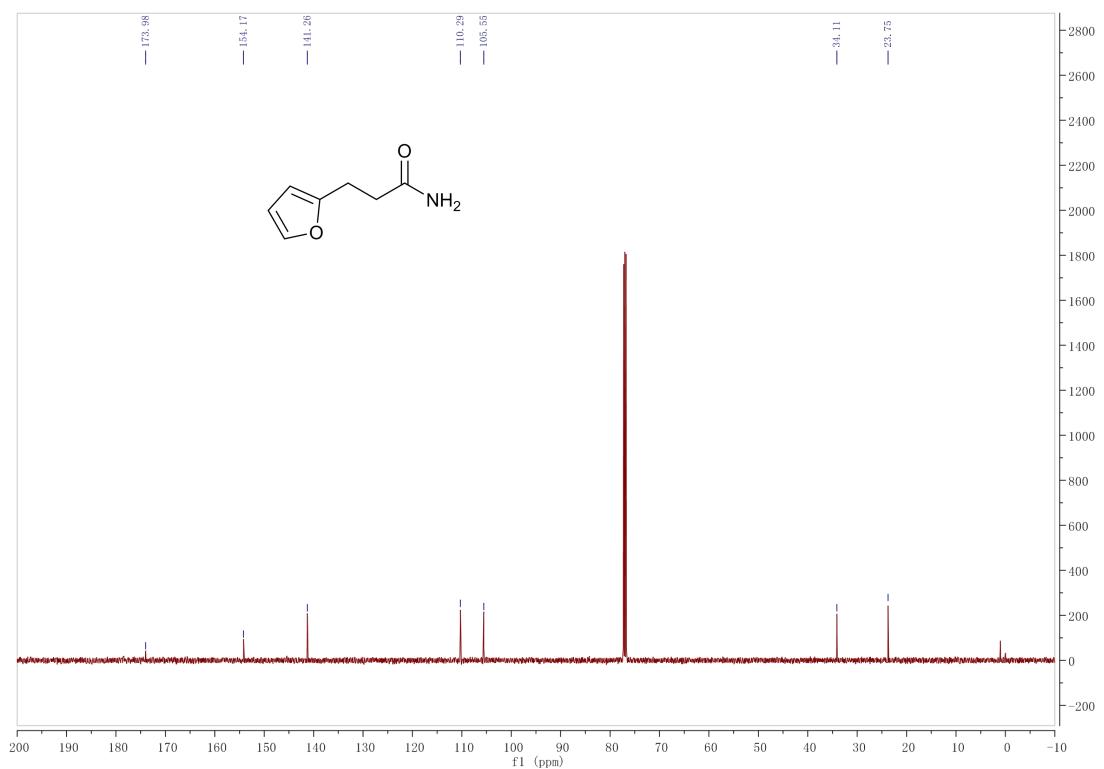
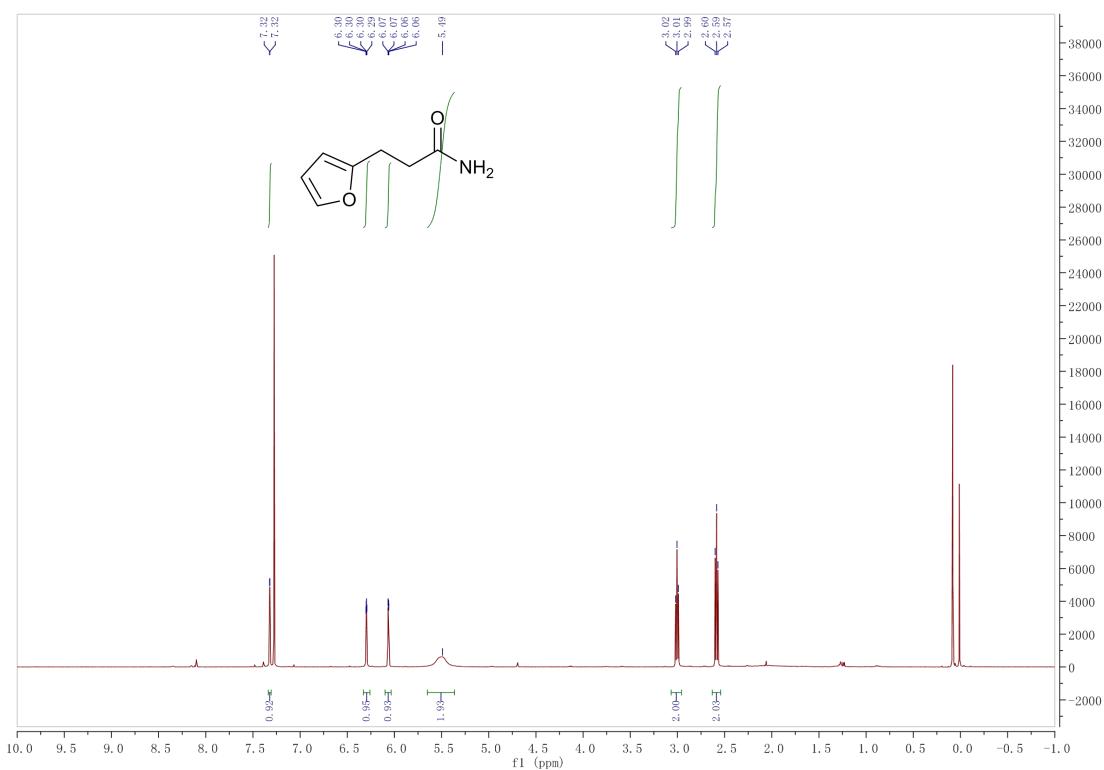
3-(3-methoxyphenyl)propenamide (4db**)**



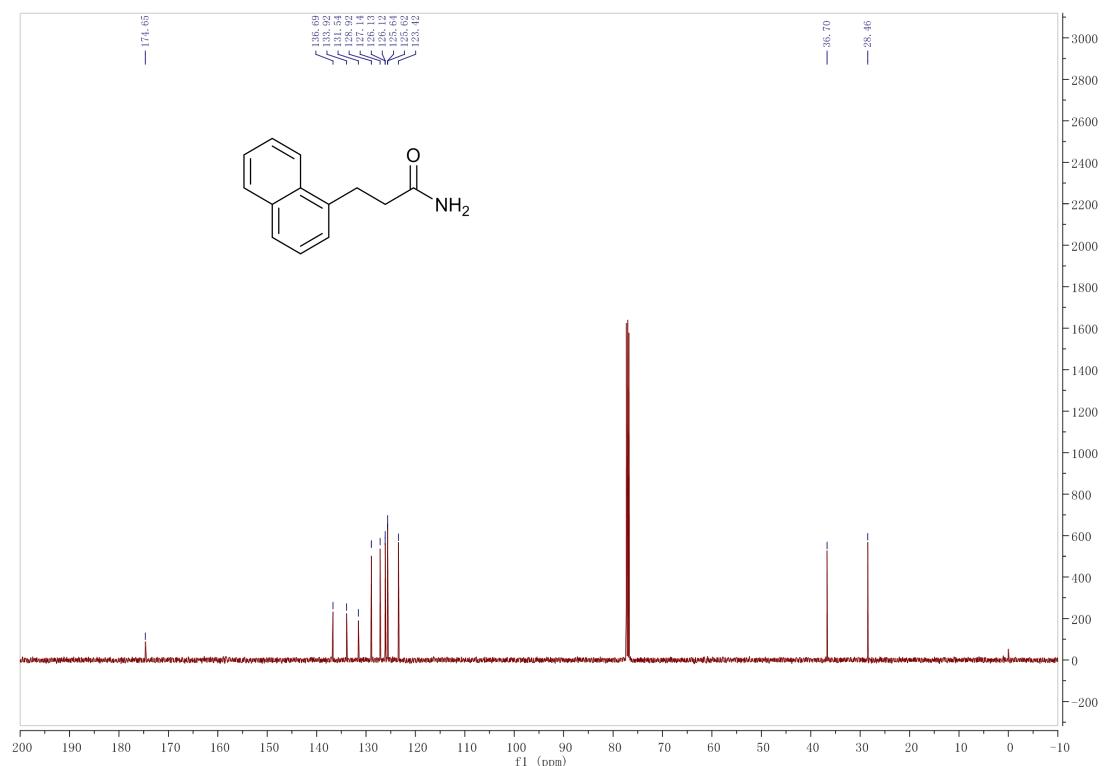
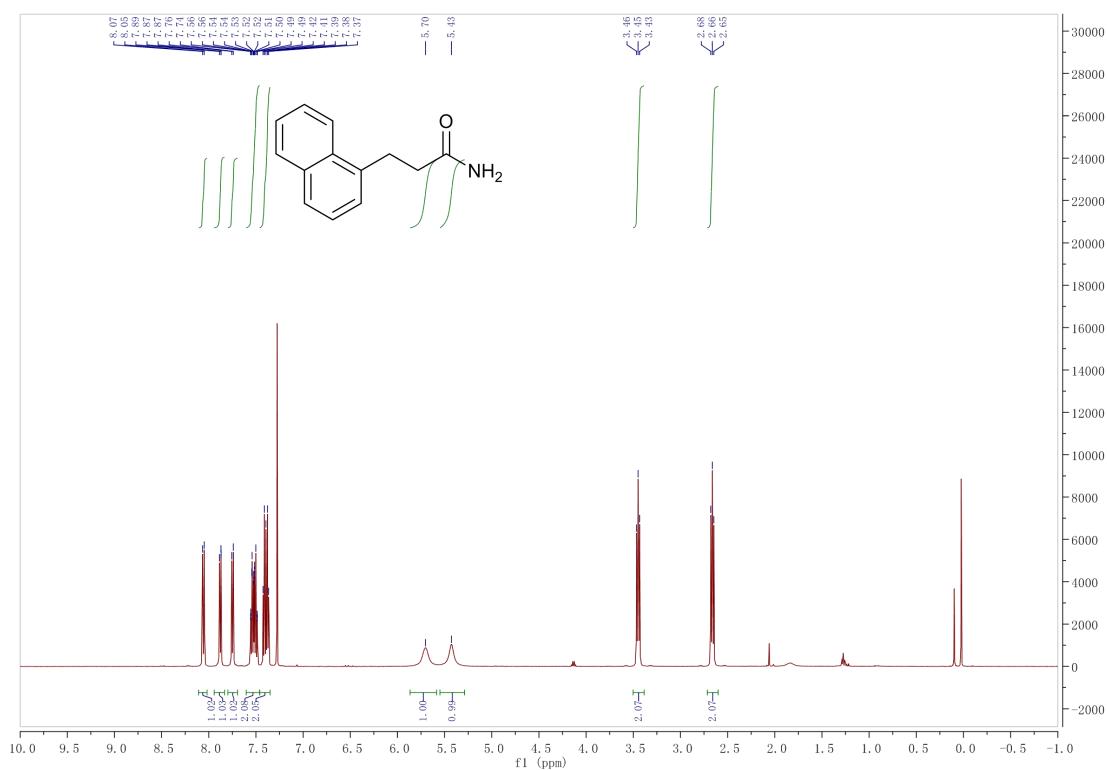
3-(3-chlorophenyl)propenamide (**4dc**)



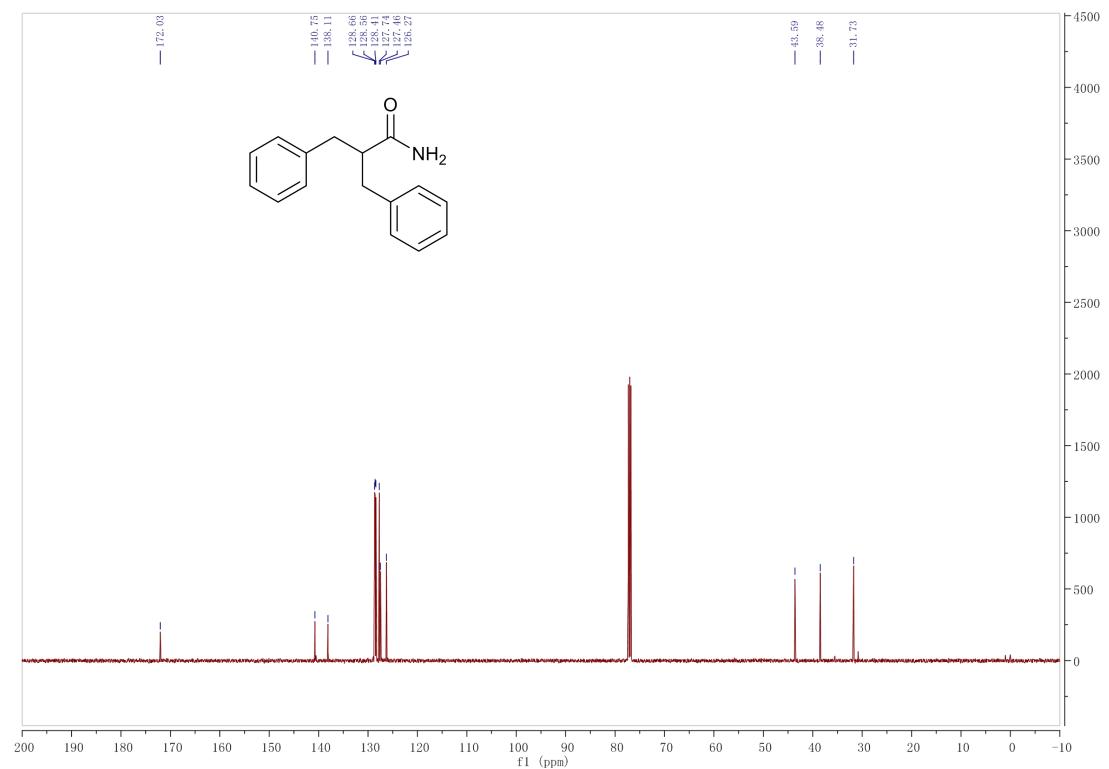
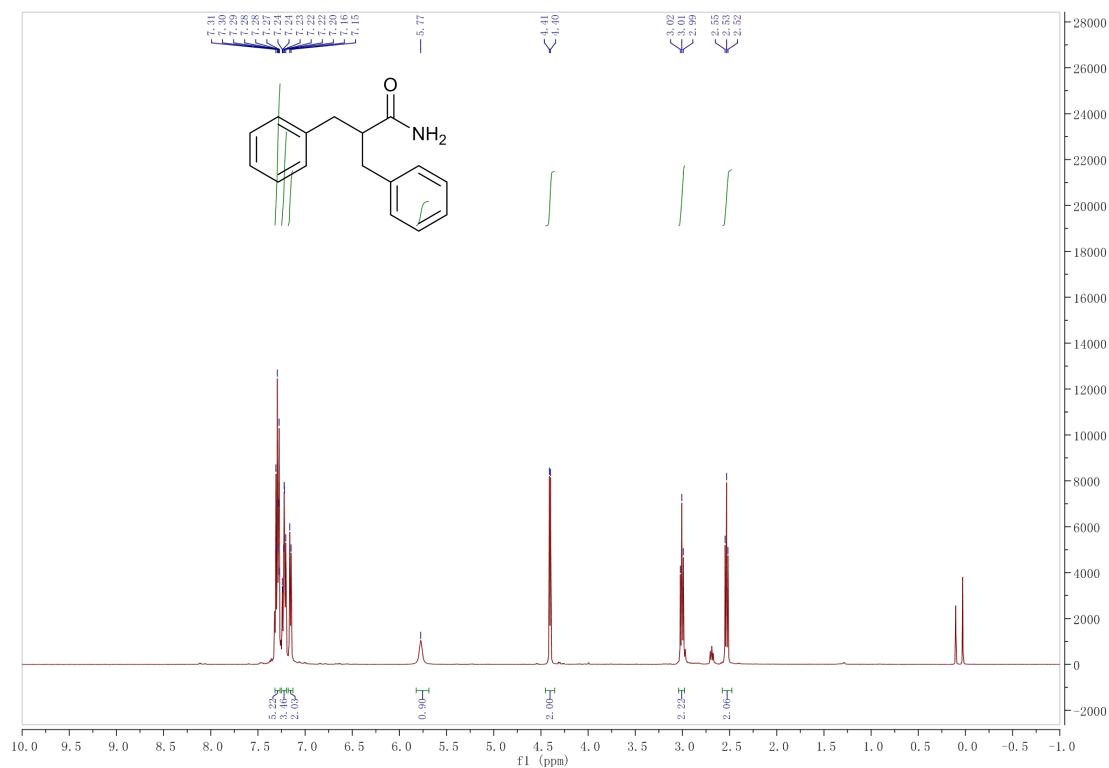
3-(furan-2-yl)propenamide (**4dd**)



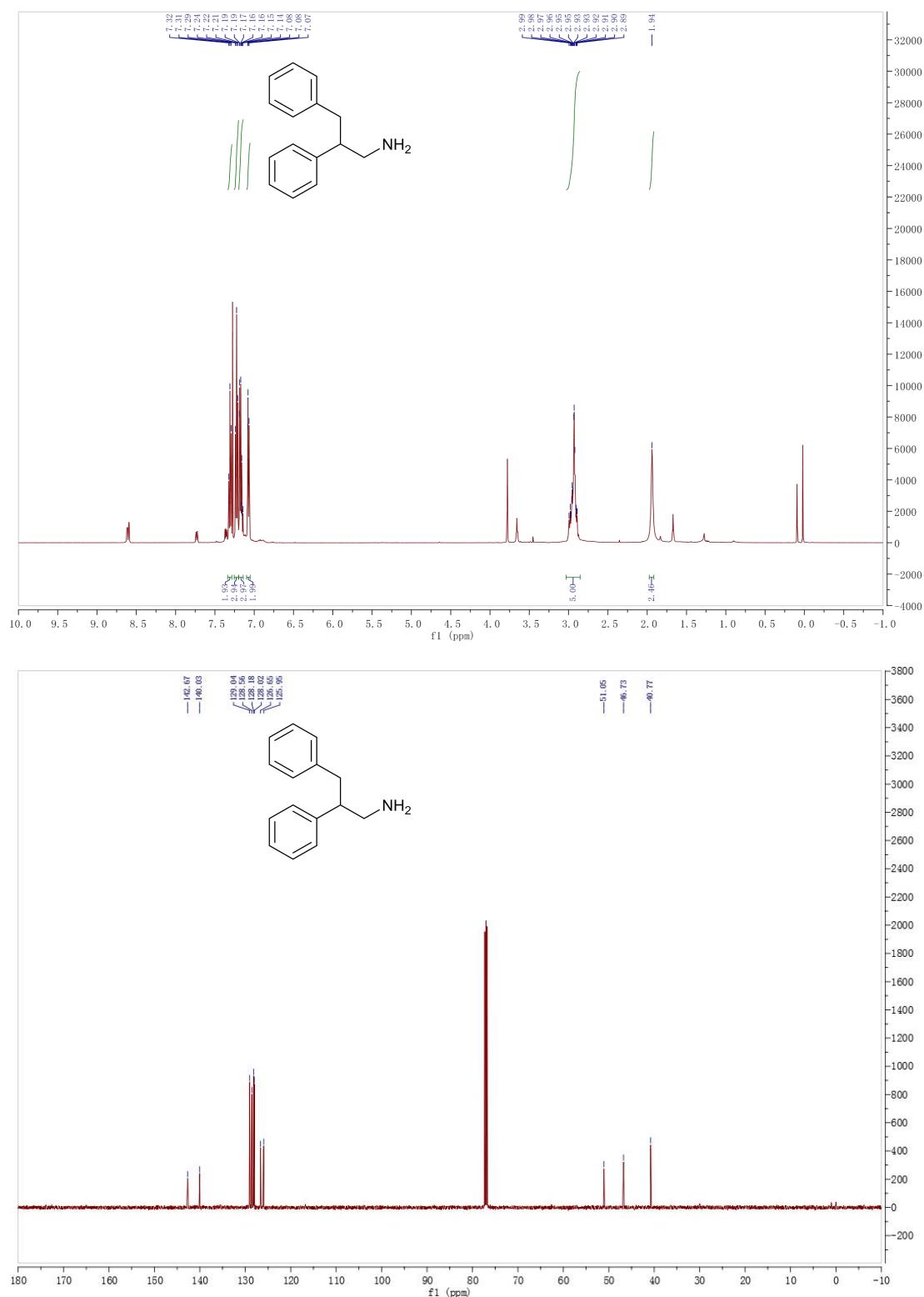
3-(naphthalen-1-yl)propenamide (4de**)**



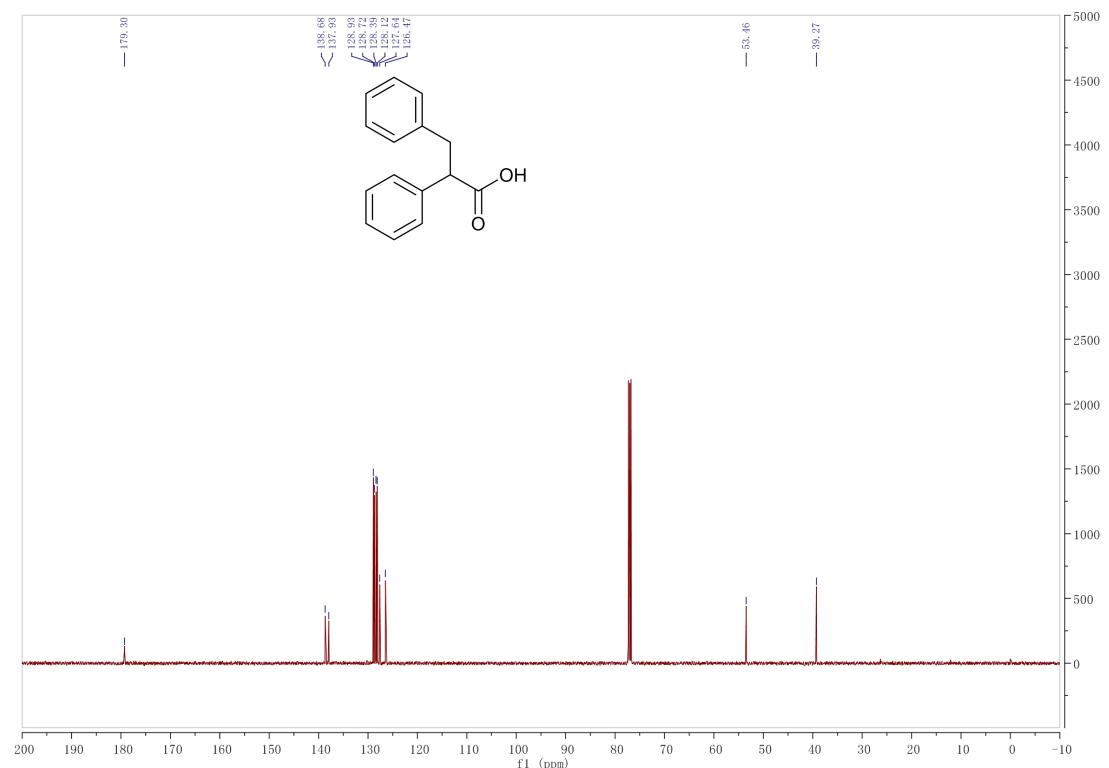
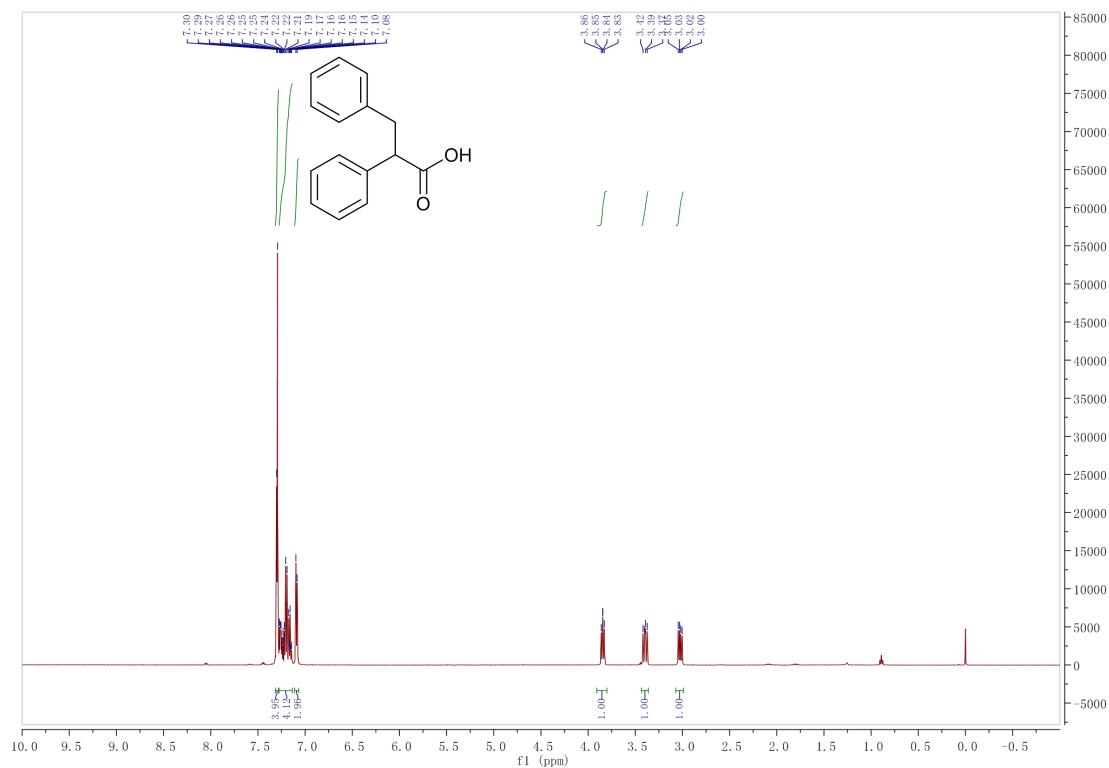
2-benzyl-3-phenylpropanamide (4df**)**



2,3-diphenylpropan-1-amine (5**)**



2,3-diphenylpropanoic acid (**6**)



2-phenyl-3,4-dihydronaphthalen-1(2H)-one (7)

